

**Typical Indoor Air Concentrations of Volatile Organic  
Compounds in Non-Smoking Montana Residences Not  
Impacted by Vapor Intrusion**

A Montana Indoor Air Quality Investigation

Prepared by the Montana Department of Environmental Quality

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## 1.0 INTRODUCTION

Vapor intrusion, the migration of volatile chemicals from a subsurface source to the indoor air of overlying or adjacent buildings, is a rapidly evolving science that has far-reaching implications to the investigation and remediation of sites impacted with volatile organic compounds (VOCs). A common confounding factor in the evaluation of data collected in vapor intrusion investigations is the separation of “background” (i.e. from typical activities and chemicals inside buildings) indoor air VOC concentrations from VOCs that may be present in indoor air due to subsurface vapor intrusion. This is especially difficult when background indoor air concentrations of VOCs exceed screening levels, such as the Environmental Protection Agency Regional Screening Levels (EPA RSLs) for contaminants of concern (COCs) during the course of vapor intrusion investigations. This scenario may result in additional sampling and data evaluation, and in difficulty explaining the investigation results to the building occupants, when ultimately the agency may determine that no vapor intrusion issue exists.

Numerous studies have documented the ubiquitous presence of some VOCs in indoor air due to human activities. Indoor sources of VOCs include consumer products (e.g., cleaners, solvents, strippers, polish, adhesives, water repellants, lubricants, air fresheners, aerosols, mothballs, scented candles, insect repellants, plastic products); combustion processes (e.g., smoking, cooking, home heating); fuels in attached garages; dry cleaned clothing or draperies; municipal tap water; or occupant activities (e.g., craft hobbies)<sup>1</sup>. In addition, sources of VOCs in outdoor air or attached garages may affect indoor air quality as the typical building experiences multiple exchanges of indoor air with attached garage or outdoor air each day.<sup>2</sup> The Montana Department of Environmental Quality (DEQ) does note that some sites investigated for vapor intrusion have demonstrated outdoor air concentrations of some VOCs greater than that observed in the indoor air.

In order to improve the understanding of typical concentrations of VOCs in the indoor air of residential homes in the state of Montana, the Department of Environmental Quality (DEQ) conducted indoor air sampling at 50 non-smoking residences in Montana that represent a cross section of “typical” residential building use (in the absence of smoking) in both urban and rural settings. Study participants were recruited based on the inclusion criteria presented in the Experimental Methods section. The goal of this study was to calculate tolerance limits and 25<sup>th</sup>, 50<sup>th</sup>, 75<sup>th</sup>, 90<sup>th</sup>, and 95<sup>th</sup> percentiles for the expected concentrations of VOCs in residential households, as measured by laboratory methods TO-15, TO-15 Selective Ion Monitoring (SIM), and hydrocarbon fractions measured by Massachusetts Air-Phase Petroleum Hydrocarbons (Mass APH).

The benefits of this study may include:

- Allowing for better evaluation of data obtained from vapor intrusion investigations by comparing measured indoor air concentrations of VOCs to the expected indoor air concentrations reported from this study;

- Providing possible alternate generic screening levels (DEQ currently uses the most current EPA RSLs) if it is found that some compounds are ubiquitously present above these EPA RSLs in indoor air;
- Providing possible alternate generic screening levels for EPA RSLs that are analytically unachievable;
- Eliminating the need to conduct more expensive TO-15 SIM analysis for some or all contaminants if it is found that these contaminants are ubiquitously present in indoor air at concentrations above standard TO-15 detection limits;
- Allowing DEQ and other environmental professionals to better explain measured indoor air concentrations of VOCs to the public by using the data obtained from this study as representative of typical concentrations of VOCs found in indoor air in Montana residences;
- Providing typical indoor air concentrations of VOCs that could be used as possible cleanup criteria for buildings requiring mitigation of subsurface vapor intrusion.

This study is not the first of its kind. Similar studies conducted to measure concentrations of VOCs in the indoor air of residential buildings include:

- Batterman et al. (2007)<sup>2</sup> – This study sampled the indoor air and garage air of 15 single family homes in southeast Michigan;
- Heroux et al. (2010)<sup>3</sup> – This study sampled the indoor air of 106 homes in the winter and 111 homes in summer of 2007 in Regina, Saskatchewan, Canada;
- Weisel et al. (2008)<sup>4</sup> – This study sampled the indoor air of 100 homes located in suburban and rural areas of New Jersey which were not expected to be impacted by vapor intrusion;
- Environmental Protection Agency (2011)<sup>1</sup> – The EPA reviewed a total of 18 studies of VOCs in the indoor air of residences of North America and developed summary ranges of VOCs present in background indoor air of North American residences between 1990 and 2005 from this compilation;
- Massachusetts Department of Environmental Protection (Mass DEP) (2008)<sup>5</sup> – Mass DEP developed criteria for reviewing available studies for inclusion in an updated 50<sup>th</sup>, 75<sup>th</sup>, and 90<sup>th</sup> percentile calculation of indoor air VOCs.
- New York State Department of Health (NYDOH) (2006)<sup>6</sup> – Appendix C of NYDOH's *Guidance for Evaluating Soil Vapor Intrusion in the State of New York* is a summary of five studies conducted between 1988 and 2005 with VOC concentration percentiles presented in this appendix.

While this study was conducted as a Montana-specific study, its findings may have application to other locations with similar urban development, building construction and building occupant habits. This study is limited only in the size of the sample (n=50). A larger sample size may provide a better range of expected indoor air concentrations of VOCs. Additionally, buildings that do not meet the inclusion criteria for the study (i.e. occupied by smokers, recent building renovations, commercial buildings, etc.) may have background concentrations of VOCs above the values calculated in this study.

## **2.0 DATA QUALITY OBJECTIVES**

Data Quality Objectives (DQOs) of this study were presented in Table 2.0 of the *Sampling, Analysis, and Quality Assurance Project Plan: Typical Indoor Air Concentrations of Volatile Organic Compounds in Non-Smoking Residences Not Impacted by Subsurface Vapor Intrusion*<sup>7</sup> Step 2, *Identifying the Decisions* of the DQOs is summarized as follows:

- Data from this study will be used to determine whether indoor air concentrations of VOCs in non-smoking residences not impacted by subsurface vapor intrusion typically exceed risk-based screening levels;
- The data will be used to determine whether alternate screening/cleanup levels for certain VOCs are appropriate because concentrations in background indoor air are higher;
- The data will be used to determine whether more costly analytical methods are necessary for site-specific vapor intrusion investigations;
- The data will be used to show members of the public who are concerned about vapor intrusion what concentrations of certain compounds are typically present in residences not impacted by vapor intrusion.

Section 5.1 provides further discussion regarding the data quality objectives.

## **3.0 EXPERIMENTAL METHODS**

### **3.1 Participant Selection Criteria**

The selection of residential buildings to include in this study was of critical importance as buildings with abnormally high levels of VOCs (i.e. cigarette smoking indoors or recent building renovations) have the potential to skew the data. This study sought to sample indoor air at a representative population of residential households according to the following criteria:

- The building must not be occupied by smokers currently, or in the past 12 months;
- The building must not have undergone any form of remodeling (including painting, wood treatment or finishing, new tile, new carpeting, etc.) within the past 12 months;
- The building must be a primary residence and occupied 12 months out of the year;
- The building must be in “sound condition” (no broken windows, missing doors, inadequate building heating, etc.);

- The building must not be used as a place of business;
- The building must not be located within ¼ mile of any known DEQ site, facility, or otherwise known VOC contamination existing in groundwater or subsurface soils.

Additional considerations in selecting the study participants were:

- A representation of buildings located in both urban and rural areas (as outdoor air VOC concentrations or building use patterns may differ between rural and urban areas);
- A representation of buildings located geographically across Montana;
- The sample population should contain a variety of building construction styles (slab on grade, subsurface basement, above ground crawlspace, etc.) and ages;
- The sample population should represent a variety of heating sources (indoor heating oil tank and furnace, natural gas fired stoves or furnaces, propane fired stove or furnace, wood stoves, etc.) as different heat sources may pose varying potential VOC impacts to indoor air;
- The sample population should include a significant population of buildings that have a garage attached to the living space. The presence of an attached garage has been identified as a major source of indoor air VOC contamination.<sup>2</sup> This study sought to include structures with attached garages due to the frequency of this type of building construction in Montana;
- Potential indoor air VOC sources were not removed from the buildings prior to the collection of indoor air samples as this study sought to quantify the impact to indoor air from these common household VOC sources.

This study employed a convenience sampling strategy; a non-probability based sampling that involves the selection of samples from a population readily available. It is believed that the population selected by convenience sampling is representative of the population as a whole.

### **3.2 Laboratory Methods and Constituents Analyzed**

Columbia Analytical Services (CAS) of Simi Valley, California was the sole laboratory for the study. CAS provided all of the 100% individually certified clean (for the entire constituent list) six-liter Summa canisters for sample collection. Table 3.2.1 lists the individual VOCs analyzed and provides their respective generic screening levels and laboratory detection limits.

**Table 3.2.1: Selected VOCs Analyzed, Generic Screening Levels (RSLs), and Laboratory Reporting Limits (RLs)**

Analyte	Chemical Abstract Services Number	Key - carcinogen (c) or non-carcinogen (n)	EPA BASED RSL* (residential air) ( $\mu\text{g}/\text{m}^3$ )	TO-15 RL ( $\mu\text{g}/\text{m}^3$ )	TO-15 SIM RL ( $\mu\text{g}/\text{m}^3$ )
1,1,1-Trichloroethane (TCA)	71-55-6	n	520	0.54	0.054
1,1,2,2-Tetrachloroethane	79-34-5	c	0.042	0.69	0.14
1,1,2-Trichloroethane	79-00-5	n	0.015	0.54	0.11
1,1-Dichloroethane (1,1-DCA)	75-34-3	c	1.5	0.40	0.081
1,1-Dichloroethene (1,1-DCE)	75-35-4	n	21	0.40	0.040
1,2,4-Trimethylbenzene	95-63-6	n	0.73	0.49	
1,2-Dibromoethane	106-93-4	c	0.0041	0.77	
1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	76-14-2		No value	0.70	
1,2-Dichlorobenzene	95-50-1	n	21	0.60	
1,2-Dichloroethane	107-06-2	c	0.094	0.40	0.081
1,2-Dichloropropane	78-87-5	c	0.24	0.46	
1,3,5-Trimethylbenzene	108-67-8		No value	0.49	
1,3-Butadiene	106-99-0	c	0.081	0.22	
1,3-Dichlorobenzene	541-73-1		No Value	0.60	
1,4-Dichlorobenzene	106-46-7	c	0.22	0.60	
2-Butanone (MEK)	78-93-3	n	520	1.5	
2-Hexanone	591-78-6	n	3.1	2.0	
4-Ethyltoluene	622-96-8		No value	0.49	
4-Methyl-2-pentanone	108-10-1	n	310	0.41	
Acetone	67-64-1	n	3200	1.2	
Benzene	71-43-2	c	0.31	0.32	0.16
Bromochloromethane	74-97-5	n	4.2		
Bromodichloromethane	75-27-4	c	0.066	0.67	
Bromoform	75-25-2	c	2.2	1.0	
Bromomethane	74-83-9	n	0.52	0.39	
Carbon Disulfide	75-15-0	n	73	1.6	
Carbon Tetrachloride	56-23-5	c	0.41	0.63	
Chlorobenzene	108-90-7	n	5.2	0.46	
Chloroethane (ethyl chloride)	75-00-3	n	1000	1.3	
Chloroform	67-66-3	c	0.11	0.49	
Chloromethane	74-87-3	n	9.4	0.21	
Cyclohexane	110-82-7	n	630	0.34	
Dibromochloromethane	124-48-1	c	0.09	0.81	
Dichlorodifluoromethane (CFC 12)	75-71-8	n	10	0.49	
Dichloromethane (methylene chloride)	75-09-2	c	96	0.69	



Analyte	Chemical Abstract Services Number	Key - carcinogen (c) or non-carcinogen (n)	EPA BASED RSL* (residential air) ( $\mu\text{g}/\text{m}^3$ )	TO-15 RL ( $\mu\text{g}/\text{m}^3$ )	TO-15 SIM RL ( $\mu\text{g}/\text{m}^3$ )
Ethylbenzene	100-41-4	c	0.97	0.43	0.087
Methyl tert-Butyl Ether	1634-04-4	c	9.4	0.36	0.36
Styrene	100-42-5	n	100	0.42	
Tetrachloroethene (PCE)	127-18-4	c	9.4	0.68	0.14
Tetrahydrofuran (THF)	109-99-9		210	1.5	
Toluene	108-88-3	n	520	0.38	
Trichloroethene (TCE)	79-01-6	n	0.43	0.54	0.11
Trichlorofluoromethane (CFC 11)	75-69-4	n	73	0.56	
Vinyl Chloride	75-01-4	c	0.16	0.26	0.026
cis-1,2-Dichloroethene	156-59-2		No value	0.40	0.079
cis-1,3-Dichloropropene (2)	10061-01-5	c	0.61	0.45	
m,p-Xylenes (1)	179601-23-1	n	10	0.43	0.17
n-Heptane	142-82-5		No value	0.41	
n-Hexane	110-54-3	n	73	0.35	
o-Xylene	95-47-6	n	10	0.43	0.087
trans-1,2-Dichloroethene	156-60-5	n	6.3	0.40	0.40
trans-1,3-Dichloropropene (2)	10061-02-6	c	0.61	0.45	
2-Propanol	67-63-0	n	730	1.2	
Ethanol	64-17-5		No value	0.94	
Hexachlorobutadiene	87-68-3	c	0.11	0.81	
Naphthalene	91-20-3	c	0.072		0.16
Cumene (isopropylbenzene)	98-82-8	n	42	0.49	
3-Chloropropene	107-05-1		No value	1.6	
2,2,4-trimethylpentane	540-84-1		No value	2.3	
1,4-Dioxane	123-91-1	c	0.32	0.36	
Propylbenzene	103-65-1		No value	0.49	
1,2,4-Trichlorobenzene	120-82-1	n	2.1	3.7	

Notes:

\* = DEQ divides non-carcinogenic RSLs by 10 to ensure a total hazard index of 1 or less. Carcinogenic RSLs are not modified

1 – m,p xylenes are listed separately in the 2012 EPA RSL table

2 – DEQ adjusted EPA RSL is for 1,3-dichloropropene

$\mu\text{g}/\text{m}^3$  – micrograms per cubic meter

Shaded compounds have RSLs that are less than the lowest practicable reporting limit

Table 3.2.2 lists the hydrocarbon fractions analyzed and provides their respective generic screening level and laboratory detection limits:

**Table 3.2.2: Hydrocarbon Fractions Analyzed by Massachusetts Air-Phase Petroleum Hydrocarbons (Mass APH), DEQ Screening Levels, and Laboratory Reporting Limits**

Analyte	CAS Number	DEQ Screening Level (residential air) ( $\mu\text{g}/\text{m}^3$ )	Reporting Limit ( $\mu\text{g}/\text{m}^3$ )
C <sub>5</sub> -C <sub>8</sub> Aliphatic Hydrocarbons	Not Available	62.6	27
C <sub>9</sub> -C <sub>12</sub> Aliphatic Hydrocarbons	Not Available	10.4	13
C <sub>9</sub> -C <sub>10</sub> Aromatic Hydrocarbons	Not Available	10.4	6.7

Notes:

**SHADED** compounds have RSLs that are less than the lowest practicable reporting limit  $\mu\text{g}/\text{m}^3$  – micrograms per cubic meter

### 3.3 Field Sampling Methodology

All samples were collected in laboratory supplied, individually certified clean six-liter Summa canisters with a dedicated vacuum gauge and flow regulator set to collect the sample over a 24-hour period. All Summa canisters were tested with a digital vacuum gauge upon receipt from the laboratory to verify that the canisters had adequate initial negative pressure (between -24” Hg and -29” Hg, adjusted for feet above sea level).

All samples were collected between March 5 and March 31, 2012. Samples were collected during a cold weather time period in order to collect samples during what has been identified as a potential worst case time period for indoor air quality.<sup>8</sup> Additionally, sampling during a cold weather time period ensured that the results from this study seasonally correlated with most results from vapor intrusion investigations in Montana.

One indoor air sample was collected per building. The placement of the Summa canister for collection of the indoor air sample adhered to the following selection criteria:

- Summa canisters were placed in a central location on what the building occupant or DEQ determined to be the main living floor of the building (central gathering or food preparation areas used most frequently by occupants of the building);
- Summa canisters were not placed in kitchens when possible so as to not preferentially sample VOCs generated by cooking activities;
- Summa canisters were placed so that the intake to the Summa canister was located between three and five feet above the floor to collect the indoor air sample from the typical breathing zone area;
- Summa canisters were not placed in closets or similar storage areas to avoid preferentially sampling common locations of VOC-containing household products;

- Summa canisters were not placed in a breezeway, doorway, or close to any other opening to the outdoors or other ventilation; and
- Summa canisters were not placed in a workshop, garage, or hobby room.

At the time of the sampling, participants were asked to complete DEQ's Occupied Building Questionnaire (Reference 8, Attachment A) which describes building construction, heat sources, occupant habits, and consumer product storage. Sampling teams assisted participants with the completion of the questionnaire when necessary and also completed indoor air sampling forms (<http://deq.mt.gov/StateSuperfund/viguide.mcp>) documenting sampling equipment serial numbers, initial and final vacuums, weather conditions, building layout, building GPS coordinates, sample location within the building, and any other relevant field observations for each location.

The sample teams were instructed to avoid using perfumes, colognes, hair spray, fueling vehicles before the collection or deployment of samples, smoking at any time during the sampling event, or engaging in any other activity that may introduce VOCs to indoor air.

Following collection, all samples were shipped back to the laboratory in their original packages under chain of custody. No more than ten days elapsed between sample collection and shipment (sample hold time is 30 days). Samples remained under DEQ custody at all times prior to shipment to the laboratory.

## 4.0 RESULTS

Data from the 50 samples, plus five field duplicates, were validated using DEQ protocols.<sup>9</sup> In brief, all data, some with qualification, are usable for the purpose of this study. In addition, all the data are assumed to be representative for the purpose of the study given that each building sampled met the previously identified selection criteria. Therefore, no outliers are assumed to be present in the data.

Data from all 50 residential buildings for each of the 77 VOCs were input into EPA ProUCL software (version 4.1.01). The 25<sup>th</sup>, 50<sup>th</sup>, 75<sup>th</sup>, 90<sup>th</sup>, and 95<sup>th</sup> percentiles were calculated using raw data for each VOC. This means that non-detect values were included at the reporting limit; no conversion or substitution of data occurred to determine percentiles. For cases where non-detect values constituted an entire percentile, the statistic is presented as less than the reporting limit at the specified percentile. For example, the 50<sup>th</sup> percentile for C9-C10 aromatics is "<9.2 µg/m<sup>3</sup>."

In addition to percentiles, upper tolerance limits (UTLs) on the median (i.e., 50<sup>th</sup> percentile) with 95% confidence (95% UTL-50<sup>th</sup>) were calculated for VOCs with detected values for at least 50 percent of each data set. In other words, if non-detect values constituted at least 50 percent of the data set for a specific VOC, then a 95% UTL-50<sup>th</sup> statistic was not calculated. A UTL is a statistic of proportion for a data set; therefore, it is a useful metric for estimating an upper limit of a percentile of a population. In particular, the 95% UTL-50<sup>th</sup> parametric is pertinent for a study of "typical" indoor air concentrations because the metric represents a statistically-based upper limit on the median of expected indoor air concentrations.

50 percent was utilized as the limit of allowable non-detect values in a data set for calculating 95% UTL-50<sup>th</sup> statistics because when calculating a statistically-based upper limit for a proportion, i.e., median, the data set should desirably contain quantified results within the respective proportion to validate the statistic. In other words, a 95% UTL-50<sup>th</sup> statistic for a data set that has 50 percent or more qualified values, i.e., non-detect values, has no quantified measurements to confirm the statistic.

In other words, DEQ calculated the average indoor air concentration of each VOC (the 50<sup>th</sup> percentile value). For those VOCs that had a non-detect value as the average (or 50<sup>th</sup> percentile) concentration, no additional statistics could be performed since the average concentration was an unquantified value (i.e. <0.84 µg/m<sup>3</sup>). For VOCs with a quantified value as the 50<sup>th</sup> percentile (i.e. benzene 50<sup>th</sup> percentile = 0.90 µg/m<sup>3</sup>) DEQ calculated the upper tolerance limit (UTL) of this value. The UTL is an estimate of what the highest value that a calculated average could be in reality, if one were to sample the entire population. The UTL typically is calculated with a level of confidence. For this study a 95% level of confidence was used when calculating the UTL. This 95%-UTL value is an estimation for what the highest real value that the calculated averages could be, with a 95% certainty. For example, the 95%-UTL-50<sup>th</sup> for benzene (1.3 µg/m<sup>3</sup>) is the estimate of the highest value that the average concentration of benzene found in the 50 houses sampled in this study (0.90 µg/m<sup>3</sup>) could be if one were to sample a much larger population.

Table 3 contains percentiles for all 77 VOCs and 95%UTL-50<sup>th</sup> statistics for the 21 VOCs that had less than 50 percent non-detect values.

**Table 4.0.1: Concentration Percentiles**

Compounds	Frequency of Non-Detects	25th Percentile (µg/m <sup>3</sup> )	50th Percentile (µg/m <sup>3</sup> )	75th Percentile (µg/m <sup>3</sup> )	90th Percentile (µg/m <sup>3</sup> )	95th Percentile (µg/m <sup>3</sup> )	95% UTL on 50th Percentile (µg/m <sup>3</sup> )
1,1,1-Trichloroethane	96%	<0.77	<0.84	<0.93	<1.0	<1.7	NA
1,1,2-Trichloroethane	100%	<0.15	<0.17	<0.19	<0.20	<0.25	NA
1,1,2-Trichlorotrifluoroethane	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
1,1-Dichloroethane	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
1,1-Dichloroethene	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
1,2,4-Trichlorobenzene	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
1,2,4-Trimethylbenzene	52%	<0.86	<1.1	1.9	4.9	8.7	NA
1,2-Dibromo-3-chloropropane	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
1,2-Dibromoethane	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
1,2-Dichlorobenzene	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
1,2-Dichloroethane	0%	0.11	0.17	0.48	0.82	1.2	0.23
1,2-Dichloropropane	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
1,3,5-Trimethylbenzene	82%	<0.78	<0.87	<0.97	2.0	2.9	NA
1,3-Butadiene	98%	<0.77	<0.84	<0.93	<0.99	<1.3	NA
1,3-Dichlorobenzene	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA

Compounds	Frequency of Non-Detects	25th Percentile ( $\mu\text{g}/\text{m}^3$ )	50th Percentile ( $\mu\text{g}/\text{m}^3$ )	75th Percentile ( $\mu\text{g}/\text{m}^3$ )	90th Percentile ( $\mu\text{g}/\text{m}^3$ )	95th Percentile ( $\mu\text{g}/\text{m}^3$ )	95% UTL on 50th Percentile ( $\mu\text{g}/\text{m}^3$ )
1,4-Dichlorobenzene	96%	<0.77	<0.84	<0.93	<1.1	<2.6	NA
1,4-Dioxane	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
2-Butanone	86%	<7.7	<8.5	<9.6	13	18	NA
2-Hexanone	90%	<0.77	<0.84	<0.93	<0.98	1.5	NA
2-Propanol	0%	11	27	46	77	95	29
Allyl Chloride	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
4-Ethyltoluene	86%	<0.78	<0.87	<0.97	2.0	2.9	NA
4-Methyl-2-pentanone	74%	<0.78	<0.88	1.0	2.1	2.3	NA
Acetone	0%	31	46	65	83	103	51
Acetonitrile	94%	<0.77	<0.84	<0.93	<1.0	1.8	NA
Acrolein	84%	<3.1	<3.5	<3.9	4.8	6.8	NA
Acrylonitrile	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
alpha-Pinene	4%	1.7	3.8	5.5	13	25	4.7
Benzene	2%	0.52	0.90	2.5	8.4	12	1.3
Benzyl Chloride	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
Bromodichloromethane	94%	<0.77	<0.84	<0.93	<1.0	1.7	NA
Bromoform	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
Bromomethane	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
C5 - C8 Aliphatics	10%	46	81	160	224	414	94
C9 - C12 Aliphatics	16%	22	33	73	100	196	44
C9 - C10 Aromatics	60%	<8.4	<9.2	12	24	38	NA
Carbon Disulfide	100%	<7.7	<8.3	<9.3	<9.8	<13	NA
Carbon Tetrachloride	92%	<0.77	<0.85	<0.93	<1.0	1.5	NA
Chlorobenzene	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
Chloroethane	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
Chloroform	64%	<0.82	<0.93	1.7	2.7	3.6	NA
Chloromethane	94%	<0.77	<0.84	<0.93	<1.0	1.3	NA
cis-1,2-Dichloroethene	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
cis-1,3-Dichloropropene	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
Cyclohexane	76%	<1.6	<1.8	<2.2	3.8	5.1	NA
Dibromochloromethane	98%	<0.77	<0.84	<0.93	<1.0	<1.4	NA
Dichlorodifluoromethane (CFC 12)	2%	1.9	2.0	2.1	2.2	2.2	2.1
Methylene Chloride	62%	<0.78	<0.93	1.5	9.0	29	NA
d-Limonene	0%	11	16	28	46	95	22
Ethanol	0%	288	625	1,375	2,020	2,930	769
Ethyl Acetate	10%	3.8	6.3	15	25	31	8.6
Ethylbenzene	6%	0.41	0.78	2.1	4.0	6.0	1.1
Hexachlorobutadiene	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
Cumene	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
m,p-Xylenes	34%	<1.7	2.7	6.6	8.9	24	3.6
Methyl Methacrylate	96%	<1.5	<1.7	<1.9	<2.0	<2.5	NA
Methyl tert-Butyl Ether	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
Naphthalene	20%	0.21	0.36	0.51	0.82	1.3	0.39
n-Butyl Acetate	38%	<0.85	1.3	2.5	4.4	5.7	1.5
n-Heptane	38%	<0.87	1.1	3.2	5.0	6.0	1.6
n-Hexane	38%	<0.91	1.5	3.2	11	16	1.9
n-Nonane	62%	<0.80	<0.93	1.8	3.0	4.0	NA
n-Octane	60%	<0.82	<0.93	1.2	2.0	2.3	NA

Compounds	Frequency of Non-Detects	25th Percentile ( $\mu\text{g}/\text{m}^3$ )	50th Percentile ( $\mu\text{g}/\text{m}^3$ )	75th Percentile ( $\mu\text{g}/\text{m}^3$ )	90th Percentile ( $\mu\text{g}/\text{m}^3$ )	95th Percentile ( $\mu\text{g}/\text{m}^3$ )	95% UTL on 50th Percentile ( $\mu\text{g}/\text{m}^3$ )
n-Propylbenzene	88%	<0.77	<0.85	<0.93	1.5	1.8	NA
o-Xylene	58%	<0.82	<0.94	2.4	5.2	7.6	NA
Propene	4%	2.2	6.0	8.7	18	35	6.5
Styrene	66%	<0.82	<0.93	1.2	2.2	2.4	NA
Tetrachloroethene	16%	0.061	0.099	0.24	2.3	2.8	0.14
Tetrahydrofuran	70%	<0.80	<0.90	1.6	3.2	4.2	NA
Toluene	2%	5.0	8.3	25	49	60	13
trans-1,2-Dichloroethene	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
trans-1,3-Dichloropropene	100%	<0.77	<0.83	<0.93	<0.98	<1.3	NA
Trichloroethene	56%	<0.042	<0.048	0.096	0.58	1.3	NA
Trichlorofluoromethane (CFC 11)	10%	1.1	1.2	1.8	2.7	3.6	1.3
Vinyl Acetate	98%	<7.7	<8.4	<9.3	<10	<15	NA
Vinyl Chloride	98%	<0.038	<0.042	<0.046	<0.049	<0.064	NA

## 5.0 DISCUSSION

### 5.1 Data Quality Objectives Discussion

The results of this study provide useful additional lines of evidence in evaluating the vapor intrusion pathway, particularly in terms of comparison of measured indoor air results from vapor intrusion investigations to the percentiles calculated from this study. The results of this study are discussed in terms of the study's data quality objectives (DQOs) as follows:

*Data from this study will be used to determine whether indoor air concentrations of VOCs in non-smoking residences not impacted by subsurface vapor intrusion typically exceed risk-based screening levels.*

Twenty-seven of 77 VOCs had non-detect values for all 50 samples (100% non-detect). For these 27 VOCs, it may be assumed that these should not normally be detected in residential indoor air. These VOCs include many compounds that are not typically contaminants of concern (COCs) in vapor intrusion investigations (i.e. bromoform, hexachlorobutadiene); however, some compounds that may be COCs for a vapor intrusion investigation (methyl tert-butyl ether (MTBE) or 1,2-dichlorobenzene) are also included in this group.

Twenty-nine of the 77 VOCs had non-detect values for greater than 50% of the samples collected, but were detected in a minimum of one sample. Together with the 27 VOCs that were not detected in any sample (discussed above), these 56 VOCs were not able to have 95%UTL-50<sup>th</sup> statistics calculated due to the fact that the 50<sup>th</sup> percentile for these compounds represents a non-detect result and not a quantified indoor air concentration.

Six of the 56 VOCs are compounds for which no screening level exists.

Twenty-nine of the 56 compounds have estimated 50<sup>th</sup> percentiles below a detection limit that are also below the respective screening level for each VOC; for these VOCs, typical indoor air concentrations in Montana residences are not expected to exceed screening levels. Of particular note in this set of 29 VOCs are C9-C10 aromatics, trichloroethene (TCE), and vinyl chloride, which are frequently compounds of potential concern for vapor intrusion investigations.

The remaining 21 VOCs from the set of 56 have estimated 50<sup>th</sup> percentiles below a detection limit that is above the respective screening levels for each VOC. This study provides the least utility for these 21 VOCs because typical indoor air concentrations in Montana residences cannot be quantified for these compounds and the qualified values of the 50<sup>th</sup> percentiles do not allow for a meaningful comparison to generic risk screening levels. The most useful finding for these 21 VOCs is that they should not typically be measured in indoor air unless they are related to subsurface contamination or they have an indoor or ambient source not usually present in the 50 residences sampled during this study.

Table 5.1.1 below summarizes the 56 VOCs that were not able to have 95%UTL-50<sup>th</sup> statistics calculated due to the fact that the 50<sup>th</sup> Percentile for these compounds represents a non-detect result and not a quantified indoor air concentration.

**Table 5.1.1 VOCs with Non-Detect Value as 50<sup>th</sup> Percentile**

Compounds	DEQ-ADJUSTED EPA RSL* (residential air) ( $\mu\text{g}/\text{m}^3$ )	50 <sup>th</sup> Percentile ( $\mu\text{g}/\text{m}^3$ )
1,1,1-Trichloroethane	520	<0.84
1,1,2-Trichloroethane	0.15	<0.17
1,1,2-Trichlorotrifluoroethane	No value	<0.83
1,1-Dichloroethane	1.5	<0.83
1,1-Dichloroethene	21	<0.83
1,2,4-Trichlorobenzene	0.21	<0.83
1,2,4-Trimethylbenzene	0.73	<1.1
1,2-Dibromo-3-chloropropane	0.00016	<0.83
1,2-Dibromoethane	0.0041	<0.83
1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	No value	<0.83
1,2-Dichlorobenzene	21	<0.83
1,2-Dichloropropane	0.24	<0.83
1,3,5-Trimethylbenzene	No value	<0.87
1,3-Butadiene	0.081	<0.84
1,3-Dichlorobenzene	No Value	<0.83
1,4-Dichlorobenzene	0.22	<0.84
1,4-Dioxane	0.32	<0.83
2-Butanone	520	<8.5
2-Hexanone	3.1	<0.84
Allyl Chloride	0.41	<0.83
4-Ethyltoluene	No value	<0.87
4-Methyl-2-pentanone	310	<0.88
Acetonitrile	6.3	<0.84
Acrolein	0.0021	<3.5
Acrylonitrile	0.036	<0.83
Benzyl Chloride	0.05	<0.83

Compounds	DEQ-ADJUSTED EPA RSL* (residential air) ( $\mu\text{g}/\text{m}^3$ )	50 <sup>th</sup> Percentile ( $\mu\text{g}/\text{m}^3$ )
Bromodichloromethane	0.066	<0.84
Bromoform	2.2	<0.83
Bromomethane	0.52	<0.83
C9 - C10 Aromatics	10.4**	<9.2
Carbon Disulfide	73	<8.3
Carbon Tetrachloride	0.41	<0.85
Chlorobenzene	5.2	<0.83
Chloroethane	1000	<0.83
Chloroform	0.11	<0.93
Chloromethane	9.4	<0.84
cis-1,2-Dichloroethene	No value	<0.83
cis-1,3-Dichloropropene	0.61	<0.83
Cyclohexane	630	<1.8
Dibromochloromethane	0.09	<0.84
Methylene Chloride	96	<0.93
Hexachlorobutadiene	0.11	<0.83
Cumene	42	<0.83
Methyl Methacrylate	73	<1.7
Methyl tert-Butyl Ether	9.4	<0.83
n-Nonane	21	<0.93
n-Octane	No value	<0.93
n-Propylbenzene	No value	<0.85
o-Xylene	10	<0.94
Styrene	100	<0.93
Tetrahydrofuran	210	<0.90
trans-1,2-Dichloroethene	No value	<0.83
trans-1,3-Dichloropropene	0.61	<0.83
Trichloroethene	0.43	<0.048
Vinyl Acetate	21	<8.4
Vinyl Chloride	0.16	<0.042

Notes:

\* = DEQ divides non-carcinogenic RSLs by 10 to ensure a total hazard index of 1 or less. Carcinogenic RSLs are not modified

\*\*= The screening level for C9-C10 aromatic hydrocarbons is the DEQ residential air screening level.

$\mu\text{g}/\text{m}^3$  – micrograms per cubic meter

Shaded compounds have a 50<sup>th</sup> percentile non-detect value above their respective screening level.

*The data will be used to determine whether alternate screening/cleanup levels for certain VOCs are appropriate because concentrations in background indoor air are higher*

Twenty-one of the 77 VOCs measured had non-detect results constitute less than 50 percent of all results for each VOC, and therefore 95%UTL-50<sup>th</sup> statistics, as previously detailed, were calculated for these VOCs. The 95%UTL-50<sup>th</sup> statistics for these 21 VOCs may be used as a line of evidence for vapor intrusion investigations when considering typical indoor air concentrations in Montana residences. These 21 VOCs are summarized in Table 5.1.2:



**Table 5.1.2 VOCs with Calculated 95% UTL-50<sup>th</sup> Percentile Statistics**

Compounds	DEQ-ADJUSTED EPA RSL* (residential air) (µg/m <sup>3</sup> )	Calculated 95% UTL-50 <sup>th</sup> Percentile (µg/m <sup>3</sup> )
1,2-Dichloroethane	0.094	0.23
2-Propanol	730	29
Acetone	3200	51
alpha-Pinene	No value	4.7
Benzene	0.31	1.3
C5 - C8 Aliphatics	62.6**	94
C9 - C12 Aliphatics	10.4**	44
Dichlorodifluoromethane (CFC 12)	10	2.1
d-Limonene	No value	22
Ethanol	No value	769
Ethyl Acetate	No value	8.6
Ethylbenzene	0.97	1.1
m,p-Xylenes	10***	3.6
Naphthalene	0.072	0.39
n-Butyl Acetate	No value	1.5
n-Heptane	No value	1.6
n-Hexane	73	1.9
Propene	310	6.5
Tetrachloroethene	9.4	0.14
Toluene	520	13
Trichlorofluoromethane (CFC 11)	73	1.3

Notes:

\* = DEQ divides non-carcinogenic RSLs by 10 to ensure a total hazard index of 1 or less. Carcinogenic RSLs are not modified

\*\*= The screening level for C5-C8 and C9-C12 aliphatics are the DEQ residential air screening levels.

\*\*\*= m,p-Xylenes are listed separately in the 2012 EPA RSL table

µg/m<sup>3</sup> – micrograms per cubic meter

Shaded compounds have a calculated 95% UTL-50<sup>th</sup> percentile value above their respective screening level.

Six VOCs (alpha-pinene, d-limonene, ethanol, ethyl acetate, n-butyl acetate, and n-heptane) in the group of 21 are compounds for which no EPA RSLs exist. These compounds are not usually COCs for vapor intrusion investigations. Regardless, the 95%UTL-50<sup>th</sup> statistics provide a reference for typical levels of these compounds in residential indoor air.

Nine VOCs (2-propanol, acetone, CFC-11, CFC-12, m&p-xylenes, n-hexane, propene, tetrachloroethene [PCE], and toluene) of these 21 VOCs have 95%UTL-50<sup>th</sup> statistics that are less than their respective RSLs; therefore, these VOCs are assumed to have common indoor and ambient sources, but typical indoor air concentrations in Montana residences related to these indoor and ambient sources are less than generic risk screening levels. Of particular note in this group of compounds are PCE, toluene, and m&p-xylenes, which are frequent COCs in vapor intrusion investigations.

For the remaining six VOCs of this group of 21 VOCs, the 95%UTL-50<sup>th</sup> statistics are greater than their respective screening levels. The six VOCs are 1,2-dichloroethane (1,2-DCA), benzene, C5-C8 aliphatics, C9-C12 aliphatics, ethylbenzene, and naphthalene. All of these compounds are common COCs in vapor intrusion investigations. These VOCs may be expected to normally be

found above their generic screening levels in the indoor air of residential homes. These compounds are compared to their respective generic screening levels (RSLs) in Table 5.1.3 for clarity.

**Table 5.1.3 VOCs with Calculated 95% UTL-50<sup>th</sup> Percentile Statistics Greater than Generic Screening Levels**

Compounds	DEQ-ADJUSTED EPA RSL* (residential air) (µg/m <sup>3</sup> )	Calculated 95% UTL-50 <sup>th</sup> Percentile (µg/m <sup>3</sup> )
1,2-Dichloroethane	0.094	0.23
Benzene	0.31	1.3
C5 - C8 Aliphatics	62.6**	94
C9 - C12 Aliphatics	10.4**	44
Ethylbenzene	0.97	1.1
Naphthalene	0.072	0.39

Notes:

\* = DEQ divides non-carcinogenic RSLs by 10 to ensure a total hazard index of 1 or less. Carcinogenic RSLs are not modified

\*\*= The screening level for C5-C8 and C9-C12 aliphatics are the DEQ residential air screening levels.

µg/m<sup>3</sup> – micrograms per cubic meter

*The data will be used to determine whether more costly analytical methods are necessary for site-specific vapor intrusion investigations*

If DEQ determines that as an outcome of this study, alternative screening levels may be appropriate for certain compounds (such as the six compounds discussed in the paragraph above and listed in table 5.1.3) then the use of the more expensive TO-15 SIM analysis could be eliminated for those compounds as long as standard TO-15 detection limits are below these alternative screening levels.

*The data will be used to show members of the public who are concerned about vapor intrusion what concentrations of certain compounds are typically present in residences not impacted by vapor intrusion*

By publishing the results of this study, DEQ provides the public with percentiles of measured indoor air concentrations of VOCs that may be expected in non-smoking, residential households that have not undergone recent remodeling. These percentiles may also be provided to members of the public who are involved in vapor intrusion investigations so that their measured indoor air results may be easily compared to results from residences known to not be impacted by subsurface vapor intrusion.

## 5.2 Additional Discussion

It is worth re-stating that evaluation of the vapor intrusion pathway requires multiple lines of evidence. The analytical results of this study demonstrate that typical indoor air concentrations in residences not impacted by subsurface vapor intrusion may vary by an order of magnitude, and measured concentrations sometimes exceed generic screening levels. As necessary, DEQ may consider the evaluation of each building as a unique situation, and allow for consideration of possibly higher typical indoor air concentrations as demonstrated by concentrations in the 75th,

90th and 95th percentiles of this study. Additionally, since the study design excluded residences that have undergone remodeling within 12 months, commercial structures, and residences occupied by smokers, these types of buildings may exhibit indoor air concentrations of VOCs equal to or greater than the upper percentiles of this study. However, variations in evaluating the vapor intrusion pathway should only occur after sufficient data have been collected for evaluation purposes and there are multiple lines of evidence indicating an indoor source.

Another line of evidence for which this study could be used is to determine whether certain building characteristics may significantly affect typical indoor air concentrations. A full analysis of this line of evidence has not been performed; however, DEQ did consider whether the presence of an attached garage affects the indoor air concentrations for the six VOCs (1,2-DCA, benzene, C5-C8 aliphatics, C9-C12 aliphatics, ethylbenzene, and naphthalene) with 95%UTL-50<sup>th</sup> statistics greater than their respective screening levels. For each VOC, the full data set may be divided into data subsets for residences with attached garages (27 data points) and residences without attached garages (23 data points). The reasoning behind the data subdivision is that people generally store many products in garages (automobiles, power equipment, building supplies, household effects, etc) that are likely sources of the six VOCs identified above, and for residences with attached garages there is assumed air flow that occurs from the garage to the occupied living space that may impact indoor air quality.<sup>2</sup>

For each of the six VOCs, a hypothesis test was performed in ProUCL to determine whether a significant difference exists in the median concentration for the attached garage subset and no attached garage subset. A non-parametric Wilcoxon-Mann-Whitney test was utilized so that potential differences or uncertainty in the underlying distributions of the data subsets was not a concern. The results indicate that for 1,2-DCA, benzene, and ethylbenzene the 95%UTL-50<sup>th</sup> statistic for the attached garage subset is significantly greater with 95% confidence than the no attached garage subset. For C5-C8 aliphatics, C9-C12 aliphatics, and naphthalene, there was no observed statistical difference at 95% confidence between the 95%UTL-50<sup>th</sup> statistics of the attached garage and no attached garage subsets.

The 95%UTL-50<sup>th</sup> statistics for 1,2-DCA, benzene, C5-C8 aliphatics, C9-C12 aliphatics, ethylbenzene, and naphthalene are identified above; however, in advanced situations involving multiple lines of evidence, DEQ may utilize the following 95%UTL-50<sup>th</sup> statistics for 1,2-DCA, benzene, and ethylbenzene to accommodate the presence of an attached garage. Table 5.2.1 highlights the comparisons of percentiles and 95%UTL-50<sup>th</sup> statistics between the full data set and attached garage and no attached garages data subsets for 1,2-DCA, benzene, and ethylbenzene. Lastly, it is noted that the 95%UTL-50<sup>th</sup> statistic for ethylbenzene for residences without an attached garage is less than the EPA RSL.

**Table 5.2.1: Comparison of Data Subsets for Residences with Attached Garages and No Attached Garages**

Compounds	Data Set	Frequency of Non-Detects	25th Percentile ( $\mu\text{g}/\text{m}^3$ )	50th Percentile ( $\mu\text{g}/\text{m}^3$ )	75th Percentile ( $\mu\text{g}/\text{m}^3$ )	90th Percentile ( $\mu\text{g}/\text{m}^3$ )	95th Percentile ( $\mu\text{g}/\text{m}^3$ )	95% UTL on 50th Percentile ( $\mu\text{g}/\text{m}^3$ )
1,2-DCA	Full	0%	0.11	0.17	0.48	0.82	1.2	0.23
	Attached Garage	0%	0.13	0.35	0.60	1.0	1.2	0.40
	No Attached Garage	0%	0.089	0.13	0.28	0.48	0.77	0.17
Benzene	Full	2%	0.52	0.90	2.5	8.4	12	1.3
	Attached Garage	0%	1.1	1.9	6.0	12	14	3.1
	No Attached Garage	4%	0.48	0.52	0.75	1.1	1.4	0.64
Ethylbenzene	Full	6%	0.41	0.78	2.1	4.0	6.0	1.1
	Attached Garage	0%	0.79	1.4	2.9	5.9	6.1	2.3
	No Attached Garage	13%	0.22	0.48	0.67	1.5	1.6	0.53

## 6.0 COMPARISON TO OTHER STUDIES

The Montana Vapor Intrusion Guide states that the range of indoor air concentrations for VOCs typically found in Montana is more narrow and falls at or below the low end of nationwide background ranges.<sup>8</sup> DEQ has previously supported this position by citing Montana's low population, rural conditions, small manufacturing base, and fewer contaminant sources of VOCs than locations predominantly included in other background studies. This indoor air study has provided data to empirically evaluate the issue.

Table 5 contains 50th percentiles from similar indoor air studies conducted in Massachusetts (2008), New York (2005), New Jersey (Weisel 2008), and an EPA (2011) compilation that combines indoor air data from 15 different studies (including the studies from Massachusetts, New York, and New Jersey) for which 50<sup>th</sup> percentiles were available from at least one other study.

**Table 6.0.1: Comparison of Typical Montana Indoor Air Concentrations to Typical Indoor Air Concentrations from Other Studies**

Compounds	Montana 50 <sup>th</sup> Percentile (µg/m <sup>3</sup> )	MA DEP 50 <sup>th</sup> Percentile <sup>1</sup> (µg/m <sup>3</sup> )	NYSDOH 50 <sup>th</sup> Percentile <sup>2</sup> (µg/m <sup>3</sup> )	New Jersey 50 <sup>th</sup> Percentile <sup>3</sup> (µg/m <sup>3</sup> )	EPA Background 50 <sup>th</sup> Percentile <sup>4</sup> (µg/m <sup>3</sup> )
1,1,1-Trichloroethane	<0.84	0.5	0.33	<2.7	<RL - 5.9
1,1,2-Trichloroethane	<0.17	ND	<0.25	No Value	No Value
1,1,2-Trichlorotrifluoroethane	<0.83	No Value	0.54	No Value	<RL - 0.5
1,1-Dichloroethane	<0.83	ND	<0.25	No Value	<RL
1,1-Dichloroethene	<0.83	ND	<0.25	No Value	No Value
1,2,4-Trichlorobenzene	<0.83	ND	<0.25	No Value	No Value
1,2,4-Trimethylbenzene	<1.1	No Value	1.9	2.5	No Value
1,2-Dibromoethane	<0.83	No Value	<0.25	No Value	No Value
1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	<0.83	No Value	<0.25	No Value	No Value
1,2-Dichlorobenzene	<0.83	ND	<0.25	No Value	No Value
1,2-Dichloroethane	0.17	ND	<0.25	No Value	<RL
1,2-Dichloropropane	<0.83	ND	<0.25	No Value	No Value
1,3,5-Trimethylbenzene	<0.87	No Value	0.64	<2.5	No Value
1,3-Butadiene	<0.84	No Value	No Value	<1.1	No Value
1,3-Dichlorobenzene	<0.83	ND	<0.25	No Value	No Value
1,4-Dichlorobenzene	<0.84	0.5	<0.25	3.0	No Value
1,4-Dioxane	<0.83	ND	No Value	No Value	No Value
2-Butanone	<8.5	3.4	3.4	3.5	No Value
4-Ethyltoluene	<0.87	No Value	No Value	<2.5	No Value
4-Methyl-2-pentanone	<0.88	No Value	No Value	<2.0	No Value
Acetone	46	26	21	35	No Value
alpha-Pinene	3.8	No Value	1.5	No Value	No Value
Benzene	0.90	2.3	2.1	1.8	<RL - 4.7
Bromodichloromethane	<0.84	ND	No Value	No Value	No Value
Bromoform	<0.83	ND	No Value	No Value	No Value
Bromomethane	<0.83	ND	<0.25	No Value	No Value
C5 - C8 Aliphatics	81	58	No Value	No Value	No Value
C9 - C12 Aliphatics	33	68	No Value	No Value	No Value
C9 - C10 Aromatics	<9.2	ND	No Value	No Value	No Value
Carbon Disulfide	<8.3	No Value	No Value	<1.6	No Value
Carbon Tetrachloride	<0.85	ND	<0.25	No Value	<RL - 0.68
Chlorobenzene	<0.83	ND	<0.25	No Value	No Value
Chloroethane	<0.83	No Value	<0.25	No Value	No Value
Chloroform	<0.93	1.9	<0.25	2.4	<RL - 2.4
Chloromethane	<0.84	No Value	0.50	1.4	No Value
cis-1,2-Dichloroethene	<0.83	ND	<0.25	No Value	<RL
cis-1,3-Dichloropropene	<0.83	ND	<0.25	No Value	No Value
Cyclohexane	<1.8	No Value	0.81	1.7	No Value
Dibromochloromethane	<0.84	ND	No Value	No Value	No Value
Dichlorodifluoromethane (CFC 12)	2.0	No Value	<0.25	3.3	No Value
Methylene Chloride	<0.93	1.4	1.4	<1.7	0.68 - 61

Compounds	Montana 50 <sup>th</sup> Percentile (µg/m <sup>3</sup> )	MA DEP 50 <sup>th</sup> Percentile <sup>1</sup> (µg/m <sup>3</sup> )	NYSDOH 50 <sup>th</sup> Percentile <sup>2</sup> (µg/m <sup>3</sup> )	New Jersey 50 <sup>th</sup> Percentile <sup>3</sup> (µg/m <sup>3</sup> )	EPA Background 50 <sup>th</sup> Percentile <sup>4</sup> (µg/m <sup>3</sup> )
d-Limonene	16	No Value	2.8	No Value	No Value
Ethanol	625	No Value	160	No Value	No Value
Ethylbenzene	0.78	1.5	1.0	2.2	1 - 3.7
Hexachlorobutadiene	<0.83	ND	<0.25	No Value	No Value
Cumene	<0.83	No Value	<0.25	No Value	No Value
m,p-Xylenes	2.7	5.9	1.5	3.8	1.5 - 14
Methyl Methacrylate	<1.7	No Value	<0.25	No Value	No Value
Methyl tert-Butyl Ether	<0.83	3.5	0.79	3.45	0.025 - 3.5
Naphthalene	0.36	ND	No Value	No Value	No Value
n-Heptane	1.1	No Value	2.8	2.0	No Value
n-Hexane	1.5	No Value	1.6	2.8	No Value
n-Nonane	<0.93	No Value	1.3	No Value	No Value
n-Octane	<0.93	No Value	0.89	No Value	No Value
n-Propylbenzene	<0.85	No Value	<0.25	No Value	No Value
o-Xylene	<0.94	5.9	1.1	2.2	1.1 - 3.6
Styrene	<0.93	0.63	0.30	<2.1	No Value
Tetrachloroethene	0.099	1.4	0.34	<3.4	<RL - 2.2
Tetrahydrofuran	<0.90	No Value	<0.25	No Value	No Value
Toluene	8.3	11	9.6	13	4.8 - 24
trans-1,2-Dichloroethene	<0.83	ND	No Value	No Value	No Value
trans-1,3-Dichloropropene	<0.83	ND	<0.25	No Value	No Value
Trichloroethene	<0.048	0.29	<0.25	<2.7	<RL - 1.1
Trichlorofluoromethane (CFC 11)	1.2	No Value	2.9	2.8	No Value
Vinyl Chloride	<0.042	ND	<0.25	No Value	<RL

Notes:

ND = not detected (reporting limit unknown)

RL = reporting limit

µg/m<sup>3</sup> – micrograms per cubic meter

The following list highlights the comparisons of significant VOCs for vapor intrusion investigations:

- 1,2,4-Trimethylbenzene – The Montana median concentration is non-detect at 1.1 µg/m<sup>3</sup>, while New York's and New Jersey's are 1.9 and 2.5 µg/m<sup>3</sup>, respectively.
- Benzene – The Montana median concentration is 0.90 µg/m<sup>3</sup>, while Massachusetts', New York's, and New Jersey's are 2.3, 2.1, and 1.8 µg/m<sup>3</sup>, respectively. The range of 50th percentiles from the EPA compilation is less than the reporting limit (not identified) to 4.7 µg/m<sup>3</sup>.
- Chloroform – The Montana median concentration is non-detect at 0.93 µg/m<sup>3</sup>, while Massachusetts', New York's, and New Jersey's are 1.9, non-detect at 0.25, and 2.4 µg/m<sup>3</sup>, respectively. The range of 50th percentiles from the EPA compilation is less than the reporting limit (not identified) to 2.4 µg/m<sup>3</sup>.
- Ethylbenzene – The Montana median concentration is 0.78 µg/m<sup>3</sup>, while Massachusetts', New York's, and New Jersey's are 1.5, 1.0, and 2.2 µg/m<sup>3</sup>, respectively. The range of 50th percentiles from the EPA compilation is 1 to 3.7 µg/m<sup>3</sup>.

- m,p-Xylenes – The Montana median concentration is 2.7  $\mu\text{g}/\text{m}^3$ , while Massachusetts', New York's, and New Jersey's are 5.9, 1.5, and 3.8  $\mu\text{g}/\text{m}^3$ , respectively. The range of 50th percentiles from the EPA compilation is 1.5 to 14  $\mu\text{g}/\text{m}^3$ .
- n-Hexane – The Montana median concentration is 1.5  $\mu\text{g}/\text{m}^3$ , while New York's and New Jersey's are 1.6 and 2.8  $\mu\text{g}/\text{m}^3$ , respectively.
- o-Xylene – The Montana median concentration is non-detect at 0.94  $\mu\text{g}/\text{m}^3$ , while Massachusetts', New York's, and New Jersey's are 5.9, 1.1, and 2.2  $\mu\text{g}/\text{m}^3$ , respectively. The range of 50th percentiles from the EPA compilation is 1.1 to 3.6  $\mu\text{g}/\text{m}^3$ .
- PCE – The Montana median concentration is 0.099  $\mu\text{g}/\text{m}^3$ , while Massachusetts', New York's, and New Jersey's are 1.4, 0.34, and non-detect at 3.4  $\mu\text{g}/\text{m}^3$ , respectively. The range of 50th percentiles from the EPA compilation is less than the reporting limit (not identified) to 2.2  $\mu\text{g}/\text{m}^3$ .

In general, this comparison demonstrates that indoor air concentrations for VOCs typically found in Montana residences fall at or below the low end of nationwide background ranges; however, the results of this study do not allow a blanket statement that all typical indoor air quality in Montana residences is better than other more urbanized and industrial States. For instance, C5-C8 aliphatics, acetone, and ethanol are VOCs for which typical indoor air concentrations in Montana exceed levels from the other studies (though C5-C8 aliphatics and ethanol were only analyzed in one of the four studies referenced). Additionally, other background indoor air studies may not have utilized the same building selection criteria, and as such may include background indoor air VOC concentrations that represent a building use not captured by this background study (e.g., cigarette smoke in indoor air).

## 7.0 SUMMARY

DEQ undertook a study of typical concentrations of VOCs in the indoor air of non-smoking residential homes in Montana not impacted by subsurface vapor intrusion. The study design included building selection criteria to control potential variation in analytical results. The following key points are outcomes of the study:

- Typical indoor air concentrations in Montana residences for 56 out of the 77 VOCs evaluated in this study are expected to be less than normally achievable laboratory reporting limits;
- The study identified 21 VOCs for which 95%UTL-50<sup>th</sup> statistics provide a useful line of evidence for vapor intrusion investigations when considering typical indoor air concentrations in Montana residences;
- Six VOCs (1,2-DCA, benzene, C5-C8 aliphatics, C9-C12 aliphatics, ethylbenzene, and naphthalene) exhibited 95%UTL-50<sup>th</sup> statistics of typical indoor air concentrations in Montana residences above their respective EPA RSLs;
- For advanced, building-specific evaluations of the vapor intrusion pathway, Table 5.2.1 provides useful typical background indoor air concentrations of 1,2-DCA, benzene, and ethylbenzene based on whether a residence has an attached garage or not;
- The upper end of the percentiles (75<sup>th</sup>, 90<sup>th</sup> or 95<sup>th</sup> percentiles) reported in Table 4.0.1 still represent concentrations of VOCs from sources other than vapor intrusion;
- The use of TO-15 SIM, depending on individual laboratory reporting limits, may not be necessary for 1,2-DCA, benzene, ethylbenzene, and naphthalene;

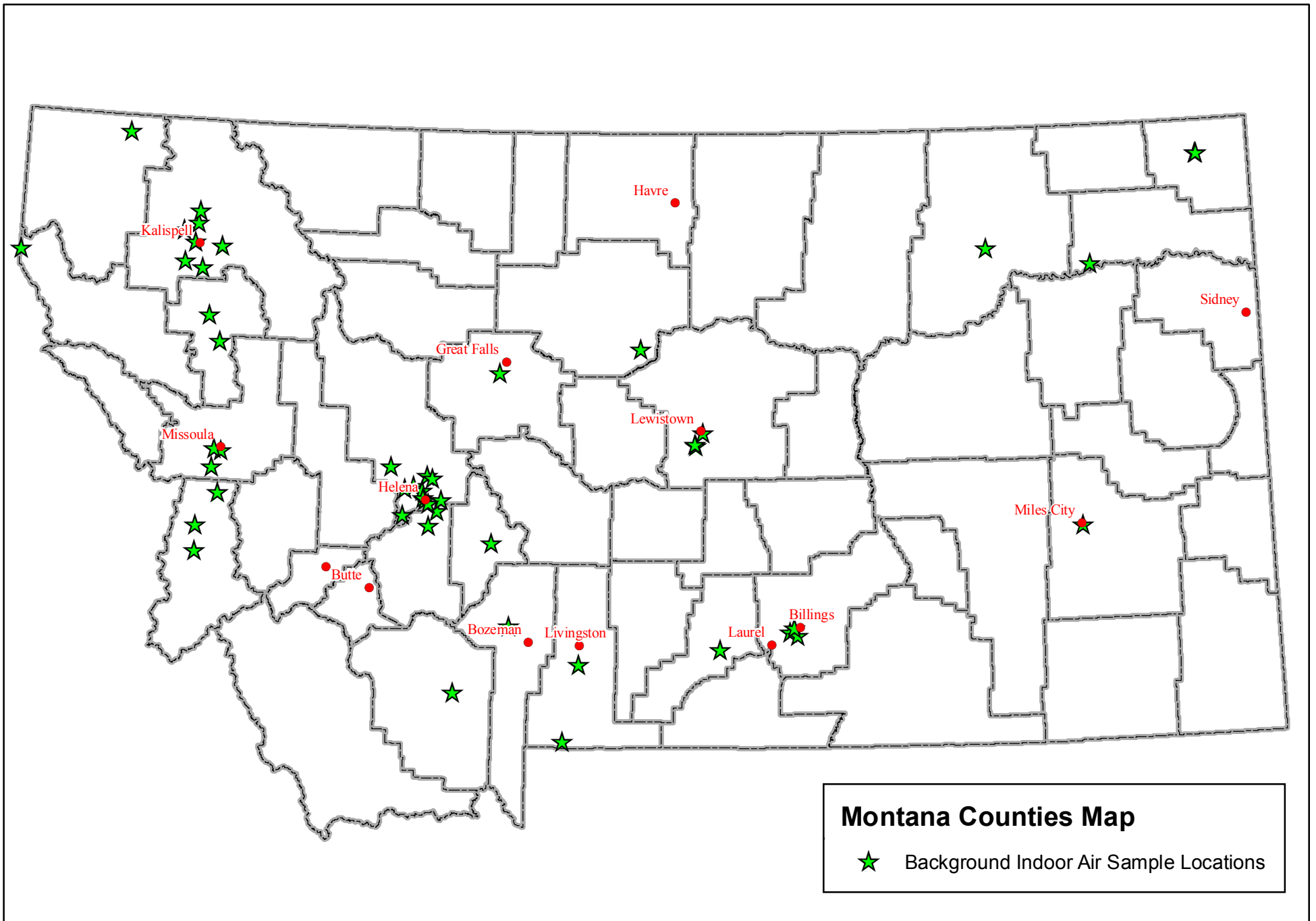
- DEQ will utilize the results of this study to inform the public involved in vapor intrusion assessments about typical indoor air concentrations in Montana residences not affected by vapor intrusion;
- DEQ will allow others (states, industries, consultants) access to the database for further evaluation of typical indoor air concentrations in Montana (e.g., evaluation of potential differences in typical indoor air concentrations based on heating fuel source);
- While additional data gathering has the potential to expand the evaluation for some VOCs or refine the evaluation of data subsets, the 50 samples collected during this study have provided a significant additional line of evidence for assessing the vapor intrusion pathway. No additional sampling is required to utilize the findings of this study.

## 8.0 REFERENCES

1. *Background Indoor Air Concentrations of Volatile Organic Compounds in North American Residences (1990-2005): A Compilation of Statistics for Assessing Vapor Intrusion*; U.S. Environmental Protection Agency; U.S. Government Printing Office: Washington, DC, 2010; EPA/530/R-10/001.
2. Batterman, S.; Jia, C.; Hatzivasilis, G. *Environ Res.* **2007**, *104*, 224-240.
3. Heroux, M.E.; Clark, N.; Ryswyk, K.V.; Ranjeeta, M.; Gilber, N.L.; Harrison, I.; Rispler, K.; Wang, D.; Anastassopoulos, A.; Guay, M.; MacNeill, M. Wheeler, A. *Int. J. Environ. Res. Public Health.* **2010**, *7*, 3080-3099.
4. Weisel, C.P.; Alimokhtari, S.; Sanders, P.F. *Environ. Sci. Technol.* **2008**, *43*, 290-307.
5. *Residential Typical Indoor Air Concentrations*; Massachusetts Department of Environmental Protection. **2008**. See <http://www.mass.gov/dep/cleanup/laws/policies.htm> (accessed July 2012).
6. *Guidance for Evaluation Soil Vapor Intrusion in the State of New York: Appendix C Volatile Organic Chemicals in Air – Summary of Background Databases.* **2006**. See [http://www.health.ny.gov/environmental/investigations/soil\\_gas/svi\\_guidance](http://www.health.ny.gov/environmental/investigations/soil_gas/svi_guidance) (accessed July 2012).
7. *Sampling, Analysis, and Quality Assurance Project Plan: Typical Indoor Air Concentrations of Volatile Organic Compounds in Non-Smoking Residences Not Impacted by Subsurface Vapor Intrusion*; Montana Department of Environmental Quality. **2011**.
8. *Montana Vapor Intrusion Guide*; Montana Department of Environmental Quality. **2011**. See <http://deq.mt.gov/StateSuperfund/viguide.mcp> (accessed July 2012).
9. *Data Validation Guidelines for Evaluating Analytical Data*; Montana Department of Environmental Quality. **2010**. See <http://deq.mt.gov/StateSuperfund/default.mcp> (accessed July 2012).



Figure



Montana Indoor Air Background Sample Locations  
in the State of Montana

# Appendix A

## LABORATORY REPORT

March 26, 2012

Christoher Cote  
Montana DEQ  
1100 N. Last Chance Gulch  
Helena, MT 59601

### **RE: Background Indoor Air Study**

Dear Christoher:

Enclosed are the results of the samples submitted to our laboratory on March 12, 2012. For your reference, these analyses have been assigned our service request number P1200947.

All analyses were performed according to our laboratory's NELAP and DoD-ELAP-approved quality assurance program. The test results meet requirements of the current NELAP and DoD-ELAP standards, where applicable, and except as noted in the laboratory case narrative provided. For a specific list of NELAP and DoD-ELAP-accredited analytes, refer to the certifications section at [www.caslab.com](http://www.caslab.com). Results are intended to be considered in their entirety and apply only to the samples analyzed and reported herein.

Columbia Analytical Services, Inc. is certified by the California Department of Health Services, NELAP Laboratory Certificate No. 02115CA; Arizona Department of Health Services, Certificate No. AZ0694; Florida Department of Health, NELAP Certification E871020; New Jersey Department of Environmental Protection, NELAP Laboratory Certification ID #CA009; New York State Department of Health, NELAP NY Lab ID No: 11221; Oregon Environmental Laboratory Accreditation Program, NELAP ID: CA20007; The American Industrial Hygiene Association, Laboratory #101661; United States Department of Defense Environmental Laboratory Accreditation Program (DoD-ELAP), Certificate No. L11-203; Pennsylvania Registration No. 68-03307; TX Commission of Environmental Quality, NELAP ID T104704413-11-2; Minnesota Department of Health, NELAP Certificate No. 362188; Washington State Department of Ecology, ELAP Lab ID: C946, State of Utah Department of Health, NELAP Certificate No. CA015272011-1; Los Angeles Department of Building and Safety, Approval No: TA00001. Each of the certifications listed above have an explicit Scope of Accreditation that applies to specific matrices/methods/analytes; therefore, please contact me for information corresponding to a particular certification.

If you have any questions, please call me at (805) 526-7161.

Respectfully submitted,

**Columbia Analytical Services, Inc.**

Kate Aguilera  
Project Manager

Client: Montana DEQ  
Project: Background Indoor Air Study

CAS Project No: P1200947

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## CASE NARRATIVE

The samples were received intact under chain of custody on March 12, 2012 and were stored in accordance with the analytical method requirements. Please refer to the sample acceptance check form for additional information. The results reported herein are applicable only to the condition of the samples at the time of sample receipt.

### Air-Phase Petroleum Hydrocarbons (APH) Analysis

The samples were analyzed for total aliphatic and aromatic gasoline range hydrocarbons by gas chromatography/mass spectrometry according to the Method for the Determination of Air-Phase Petroleum Hydrocarbons (APH), Massachusetts Department of Environmental Protection, Revision 1, December, 2009.

### Volatile Organic Compound Analysis

The samples were also analyzed in scan and SIM mode for selected volatile organic compounds in accordance with EPA Method TO-15 from the Compendium of Methods for the Determination of Toxic Organic Compounds in Ambient Air, Second Edition (EPA/625/R-96/010b), January, 1999. The analytical system was comprised of a gas chromatograph / mass spectrometer (GC/MS) interfaced to a whole-air preconcentrator.

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*The results of analyses are given in the attached laboratory report. All results are intended to be considered in their entirety, and Columbia Analytical Services, Inc. (CAS) is not responsible for utilization of less than the complete report.*

*Use of Columbia Analytical Services, Inc. (CAS) Name. Client shall not use CAS's name or trademark in any marketing or reporting materials, press releases or in any other manner ("Materials") whatsoever and shall not attribute to CAS any test result, tolerance or specification derived from CAS's data ("Attribution") without CAS's prior written consent, which may be withheld by CAS for any reason in its sole discretion. To request CAS's consent, Client shall provide copies of the proposed Materials or Attribution and describe in writing Client's proposed use of such Materials or Attribution. If CAS has not provided written approval of the Materials or Attribution within ten (10) days of receipt from Client, Client's request to use CAS's name or trademark in any Materials or Attribution shall be deemed denied. CAS may, in its discretion, reasonably charge Client for its time in reviewing Materials or Attribution requests. Client acknowledges and agrees that the unauthorized use of CAS's name or trademark may cause CAS to incur irreparable harm for which the recovery of money damages will be inadequate. Accordingly, Client acknowledges and agrees that a violation shall justify preliminary injunctive relief. For questions contact the laboratory.*

DETAIL SUMMARY REPORT

Client: Montana DEQ  
 Project ID: Background Indoor Air Study

Service Request: P1200947

Date Received: 3/12/2012  
 Time Received: 10:05

Client Sample ID	Lab Code	Matrix	Date Collected	Time Collected	Container ID	Pi1 (psig)	Pf1 (psig)	MA APH 1.0 - MA VOC PH Can	TO-15 - VOC Cans	TO-15 - VOC SIM
01-IF-030512	P1200947-001	Air	3/5/2012	20:30	AC01505	-3.41	3.52	X	X	X
02-IF-030512	P1200947-002	Air	3/5/2012	19:00	AC00687	-4.86	3.52	X	X	X
03-IF-030512	P1200947-003	Air	3/5/2012	19:45	AS00178	-4.80	3.51	X	X	X
05-IF-030712	P1200947-004	Air	3/7/2012	12:45	AC01198	-3.39	3.53	X	X	X
06-IF-030712	P1200947-005	Air	3/7/2012	12:16	AS00137	-2.12	3.74	X	X	X
07-IF-030712	P1200947-006	Air	3/7/2012	15:23	AC01570	-2.36	3.72	X	X	X
08-IF-030712	P1200947-007	Air	3/7/2012	07:17	AS00184	-2.87	3.58	X	X	X
09-IF-030512	P1200947-008	Air	3/5/2012	16:23	AC00553	-4.65	3.56	X	X	X

**Air - Chain of Custody Record & Analytical Service Request**

**Requested Turnaround Time in Business Days (Surcharges) please circle**  
 1 Day (100%) 2 Day (75%) 3 Day (50%) 4 Day (35%) 5 Day (25%) (10 Day-Standard)  
 CAS Project No. P1200947

Company Name & Address (Reporting Information)  
 MTDEG  
 PO Box 200701  
 Helena MT 59620-0701

Project Name: Background Indoor Air Study  
 Project Number: \_\_\_\_\_  
 CAS Contact: Nick Romero  
 Analysis Method: \_\_\_\_\_

Project Manager: Chris Cote  
 Phone: 406-841-5078 Fax: 406-84-5050

Email Address for Result Reporting: cote.c@mt.gov; Marti.d.k.bj@cdm.com

P.O. # / Billing Information: \_\_\_\_\_  
 Reference as: Background Indoor Air Study  
 Sampler (Print & Sign): Bill to Chris Cote of MTDEG

Client Sample ID	Laboratory ID Number	Date Collected	Time Collected	Canister ID (Bar code # - AC, SC, etc.)	Flow Controller ID (Bar code # - FC #)	Canister Start Pressure "Hg	Canister End Pressure "Hg/psig	Sample Volume	TO-15	TO-15 SIM	MA-APH	Comments
<u>01-01E-030512</u>	<u>1-331</u>	<u>3/5/12</u>	<u>8:30pm</u>	<u>AC01505</u>	<u>FC A00377</u>	<u>28.5</u>	<u>5.5</u>		<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<u>Average</u>
<u>02-1E-030512</u>	<u>1-471</u>	<u>3/5/12</u>	<u>19:00</u>	<u>AC00657</u>	<u>FC A00009</u>	<u>28.5</u>	<u>9</u>		<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<u>Project</u>
<u>03-1E-030512</u>	<u>3-467</u>	<u>3/5/12</u>	<u>19:45</u>	<u>AS00178</u>	<u>FC A00201</u>	<u>29</u>	<u>9.5</u>		<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<u>Specific list of compounds</u>
<u>05-1E-030712</u>	<u>4-326</u>	<u>3/7/12</u>	<u>12:45</u>	<u>AC01198</u>	<u>FC A00084</u>	<u>28</u>	<u>5.5</u>		<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<u>to provide</u>
<u>06-1E-030712</u>	<u>5-203</u>	<u>3/7/12</u>	<u>12:16</u>	<u>AS00137</u>	<u>FC A00447</u>	<u>28.5</u>	<u>4.5</u>		<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<u>specific reporting limits</u>
<u>07-1E-030712</u>	<u>6-231</u>	<u>3/7/12</u>	<u>15:23</u>	<u>AC01570</u>	<u>FC A00429</u>	<u>28.0</u>	<u>3</u>		<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	
<u>08-1E-030712</u>	<u>7-276</u>	<u>3/7/12</u>	<u>7:17am</u>	<u>AS00184</u>	<u>FC A00378</u>	<u>28</u>	<u>5.5</u>		<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	
<u>09-1E-030512</u>	<u>8-453</u>	<u>3/5/12</u>	<u>16:23</u>	<u>AC00553</u>	<u>FC A00048</u>	<u>26</u>	<u>6.5</u>		<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	
<u>51-1E-030512</u>		<u>3/5/12</u>	<u>20:30</u>	<u>AC01472</u>	<u>FC A00423</u>	<u>28</u>	<u>7.5</u>		<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	

Report Tier Levels - please select

Tier I - Results (Default if not specified) \_\_\_\_\_  
 Tier II (Results + QC Summaries) \_\_\_\_\_  
 Tier III (Results + QC & Calibration Summaries)   
 Tier IV (Data Validation Package) 10% Surcharge \_\_\_\_\_  
 EDD required  Yes  No  
 Type: CS17 and Excel

Relinquished by: (Signature) [Signature] Date: 3/9/12 Time: 2:00pm  
 Received by: (Signature) [Signature] Date: 3/9/12 Time: 1:00pm  
 Project Requirements (MFLs, QAPP)  
 Cooler / Blank Temperature \_\_\_\_\_ °C

**Sample Acceptance Check Form**

Client: Montana Department of Environmental Quality Work order: P1200947

Project: Background Indoor Air Study

Sample(s) received on: 3/12/12 Date opened: 3/12/12 by: MZAMORA

**Note:** This form is used for all samples received by CAS. The use of this form for custody seals is strictly meant to indicate presence/absence and not as an indication of compliance or nonconformity. Thermal preservation and pH will only be evaluated either at the request of the client and/or as required by the method/SOP.

- |  | <u>Yes</u>                          | <u>No</u>                           | <u>N/A</u>                          |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| 1 Were <b>sample containers</b> properly marked with client sample ID?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 2 Container(s) <b>supplied by CAS</b> ?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 3 Did <b>sample containers</b> arrive in good condition?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 4 Were <b>chain-of-custody</b> papers used and filled out?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 5 Did <b>sample container labels</b> and/or tags agree with custody papers?                                      | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 6 Was <b>sample volume</b> received adequate for analysis?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 7 Are samples within specified holding times?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 8 Was proper <b>temperature</b> (thermal preservation) of cooler at receipt adhered to?                          | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 9 Was a <b>trip blank</b> received?  | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| 10 Were <b>custody seals</b> on outside of cooler/Box?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| Location of seal(s)? _____ Sealing Lid?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were signature and date included?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were seals intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were custody seals on outside of sample container?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| Location of seal(s)? _____ Sealing Lid?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were signature and date included?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were seals intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 11 Do containers have appropriate <b>preservation</b> , according to method/SOP or Client specified information? | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Is there a client indication that the submitted samples are <b>pH</b> preserved?                                 | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were <b>VOA vials</b> checked for presence/absence of air bubbles?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Does the client/method/SOP require that the analyst check the sample pH and <u>if necessary</u> alter it?        | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 12 <b>Tubes:</b> Are the tubes capped and intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Do they contain moisture?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 13 <b>Badges:</b> Are the badges properly capped and intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Are dual bed badges separated and individually capped and intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |

Lab Sample ID	Container Description	Required pH *	Received pH	Adjusted pH	VOA Headspace (Presence/Absence)	Receipt / Preservation Comments
P1200947-001.01	6.0 L Ambient Can					
P1200947-002.01	6.0 L Ambient Can					
P1200947-003.01	6.0 L Silonite Can					
P1200947-004.01	6.0 L Ambient Can					
P1200947-005.01	6.0 L Silonite Can					
P1200947-006.01	6.0 L Ambient Can					
P1200947-007.01	6.0 L Silonite Can					
P1200947-008.01	6.0 L Ambient Can					

Explain any discrepancies: (include lab sample ID numbers): \_\_\_\_\_



RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 01-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-001

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01505

Date Collected: 3/5/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/21/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.41      Final Pressure (psig): 3.52

Canister Dilution Factor: 1.61

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	200	32	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	21	16	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	12	8.1	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 02-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-002

Test Code: Massachusetts APH, Revision 1, December 2009  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00687

Date Collected: 3/5/12  
Date Received: 3/12/12  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.86      Final Pressure (psig): 3.52

Canister Dilution Factor: 1.85

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	350	37	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	57	19	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	41	9.3	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 03-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-003

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00178

Date Collected: 3/5/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/21/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.80      Final Pressure (psig): 3.51

Canister Dilution Factor: 1.84

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	130	37	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	99	18	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	9.2	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 05-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-004

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01198

Date Collected: 3/7/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/21/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.39      Final Pressure (psig): 3.53

Canister Dilution Factor: 1.61

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	500	32	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	100	16	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	70	8.1	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 06-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-005

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00137

Date Collected: 3/7/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/21/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.12      Final Pressure (psig): 3.74

Canister Dilution Factor: 1.47

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	200	29	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	91	15	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	23	7.4	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 07-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-006

Test Code: Massachusetts APH, Revision 1, December 2009  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01570

Date Collected: 3/7/12  
Date Received: 3/12/12  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.36 Final Pressure (psig): 3.72

Canister Dilution Factor: 1.49

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	39	30	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	94	15	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	8.6	7.5	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 08-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-007

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00184

Date Collected: 3/7/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/21/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.87      Final Pressure (psig): 3.58

Canister Dilution Factor: 1.55

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	70	31	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	34	16	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	7.8	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 09-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-008

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00553

Date Collected: 3/5/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/21/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.65      Final Pressure (psig): 3.56

Canister Dilution Factor: 1.82

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	180	36	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	22	18	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	9.1	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P120321-MB

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/21/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	20	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	10	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	5.0	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

LABORATORY DUPLICATE SUMMARY RESULTS

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**Client:** Montana DEQ  
**Client Sample ID:** 02-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-002DUP

Test Code: Massachusetts APH, Revision 1, December 2009  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00687

Date Collected: 3/5/12  
Date Received: 3/12/12  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.86 Final Pressure (psig): 3.52

Canister Dilution Factor: 1.85

Compound	Sample Result	Duplicate Sample Result	Average	% RPD	RPD	Data
	$\mu\text{g}/\text{m}^3$	$\mu\text{g}/\text{m}^3$	$\mu\text{g}/\text{m}^3$		Limit	Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	353	352	352.5	0.3	30	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	57.1	54.2	55.65	5	30	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	41.2	41.1	41.15	0.2	30	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 01-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-001

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01505

Date Collected: 3/5/12  
Date Received: 3/12/12  
Date Analyzed: 3/21 - 3/22/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.050 Liter(s)

Initial Pressure (psig): -3.41 Final Pressure (psig): 3.52

Canister Dilution Factor: 1.61

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	7.4	0.81	4.3	0.47	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.1	0.81	0.43	0.16	
74-87-3	Chloromethane	ND	0.81	ND	0.39	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.81	ND	0.12	
106-99-0	1,3-Butadiene	ND	0.81	ND	0.36	
74-83-9	Bromomethane	ND	0.81	ND	0.21	
75-00-3	Chloroethane	ND	0.81	ND	0.31	
64-17-5	Ethanol	2,200	160	1,200	85	D
75-05-8	Acetonitrile	ND	0.81	ND	0.48	
107-02-8	Acrolein	4.2	3.2	1.8	1.4	
67-64-1	Acetone	76	8.1	32	3.4	
75-69-4	Trichlorofluoromethane	1.2	0.81	0.21	0.14	
67-63-0	2-Propanol (Isopropyl Alcohol)	40	1.6	16	0.66	
107-13-1	Acrylonitrile	ND	0.81	ND	0.37	
75-35-4	1,1-Dichloroethene	ND	0.81	ND	0.20	
75-09-2	Methylene Chloride	0.83	0.81	0.24	0.23	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.81	ND	0.26	
76-13-1	Trichlorotrifluoroethane	ND	0.81	ND	0.11	
75-15-0	Carbon Disulfide	ND	8.1	ND	2.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.81	ND	0.20	
75-34-3	1,1-Dichloroethane	ND	0.81	ND	0.20	
1634-04-4	Methyl tert-Butyl Ether	ND	0.81	ND	0.22	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 01-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-001

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01505

Date Collected: 3/5/12  
Date Received: 3/12/12  
Date Analyzed: 3/21 - 3/22/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.050 Liter(s)

Initial Pressure (psig): -3.41 Final Pressure (psig): 3.52

Canister Dilution Factor: 1.61

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	8.1	ND	2.3	
78-93-3	2-Butanone (MEK)	ND	8.1	ND	2.7	
156-59-2	cis-1,2-Dichloroethene	ND	0.81	ND	0.20	
141-78-6	Ethyl Acetate	<b>25</b>	1.6	<b>7.0</b>	0.45	
110-54-3	n-Hexane	<b>8.0</b>	0.81	<b>2.3</b>	0.23	
67-66-3	Chloroform	ND	0.81	ND	0.16	
109-99-9	Tetrahydrofuran (THF)	ND	0.81	ND	0.27	
71-55-6	1,1,1-Trichloroethane	ND	0.81	ND	0.15	
56-23-5	Carbon Tetrachloride	ND	0.81	ND	0.13	
110-82-7	Cyclohexane	ND	1.6	ND	0.47	
78-87-5	1,2-Dichloropropane	ND	0.81	ND	0.17	
75-27-4	Bromodichloromethane	ND	0.81	ND	0.12	
123-91-1	1,4-Dioxane	ND	0.81	ND	0.22	
80-62-6	Methyl Methacrylate	<b>1.9</b>	1.6	<b>0.47</b>	0.39	
142-82-5	n-Heptane	<b>3.1</b>	0.81	<b>0.76</b>	0.20	
10061-01-5	cis-1,3-Dichloropropene	ND	0.81	ND	0.18	
108-10-1	4-Methyl-2-pentanone	ND	0.81	ND	0.20	
10061-02-6	trans-1,3-Dichloropropene	ND	0.81	ND	0.18	
108-88-3	Toluene	<b>31</b>	0.81	<b>8.2</b>	0.21	
591-78-6	2-Hexanone	<b>0.93</b>	0.81	<b>0.23</b>	0.20	
124-48-1	Dibromochloromethane	ND	0.81	ND	0.095	
106-93-4	1,2-Dibromoethane	ND	0.81	ND	0.10	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 01-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-001

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01505

Date Collected: 3/5/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/21 - 3/22/12  
 Volume(s) Analyzed: 1.00 Liter(s)  
 0.050 Liter(s)

Initial Pressure (psig): -3.41 Final Pressure (psig): 3.52

Canister Dilution Factor: 1.61

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	4.3	0.81	0.90	0.17	
111-65-9	n-Octane	1.4	0.81	0.29	0.17	
108-90-7	Chlorobenzene	ND	0.81	ND	0.17	
179601-23-1	m,p-Xylenes	6.7	1.6	1.5	0.37	
75-25-2	Bromoform	ND	0.81	ND	0.078	
100-42-5	Styrene	2.2	0.81	0.51	0.19	
95-47-6	o-Xylene	2.0	0.81	0.47	0.19	
111-84-2	n-Nonane	ND	0.81	ND	0.15	
98-82-8	Cumene	ND	0.81	ND	0.16	
80-56-8	alpha-Pinene	11	0.81	1.9	0.14	
103-65-1	n-Propylbenzene	ND	0.81	ND	0.16	
622-96-8	4-Ethyltoluene	ND	0.81	ND	0.16	
108-67-8	1,3,5-Trimethylbenzene	ND	0.81	ND	0.16	
95-63-6	1,2,4-Trimethylbenzene	1.5	0.81	0.30	0.16	
100-44-7	Benzyl Chloride	ND	0.81	ND	0.16	
541-73-1	1,3-Dichlorobenzene	ND	0.81	ND	0.13	
106-46-7	1,4-Dichlorobenzene	ND	0.81	ND	0.13	
95-50-1	1,2-Dichlorobenzene	ND	0.81	ND	0.13	
5989-27-5	d-Limonene	170	16	31	2.9	D
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.81	ND	0.083	
120-82-1	1,2,4-Trichlorobenzene	ND	0.81	ND	0.11	
87-68-3	Hexachlorobutadiene	ND	0.81	ND	0.075	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 02-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-002

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00687

Date Collected: 3/5/12  
Date Received: 3/12/12  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.86      Final Pressure (psig): 3.52

Canister Dilution Factor: 1.85

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	3.6	0.93	2.1	0.54	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.1	0.93	0.43	0.19	
74-87-3	Chloromethane	ND	0.93	ND	0.45	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.93	ND	0.13	
106-99-0	1,3-Butadiene	ND	0.93	ND	0.42	
74-83-9	Bromomethane	ND	0.93	ND	0.24	
75-00-3	Chloroethane	ND	0.93	ND	0.35	
64-17-5	Ethanol	580	9.3	310	4.9	
75-05-8	Acetonitrile	ND	0.93	ND	0.55	
107-02-8	Acrolein	ND	3.7	ND	1.6	
67-64-1	Acetone	29	9.3	12	3.9	
75-69-4	Trichlorofluoromethane	1.1	0.93	0.19	0.16	
67-63-0	2-Propanol (Isopropyl Alcohol)	18	1.9	7.2	0.75	
107-13-1	Acrylonitrile	ND	0.93	ND	0.43	
75-35-4	1,1-Dichloroethene	ND	0.93	ND	0.23	
75-09-2	Methylene Chloride	ND	0.93	ND	0.27	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.93	ND	0.30	
76-13-1	Trichlorotrifluoroethane	ND	0.93	ND	0.12	
75-15-0	Carbon Disulfide	ND	9.3	ND	3.0	
156-60-5	trans-1,2-Dichloroethene	ND	0.93	ND	0.23	
75-34-3	1,1-Dichloroethane	ND	0.93	ND	0.23	
1634-04-4	Methyl tert-Butyl Ether	ND	0.93	ND	0.26	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 02-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-002

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00687

Date Collected: 3/5/12  
Date Received: 3/12/12  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.86      Final Pressure (psig): 3.52

Canister Dilution Factor: 1.85

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	9.3	ND	2.6	
78-93-3	2-Butanone (MEK)	ND	9.3	ND	3.1	
156-59-2	cis-1,2-Dichloroethene	ND	0.93	ND	0.23	
141-78-6	Ethyl Acetate	<b>37</b>	1.9	<b>10</b>	0.51	
110-54-3	n-Hexane	<b>14</b>	0.93	<b>3.8</b>	0.26	
67-66-3	Chloroform	ND	0.93	ND	0.19	
109-99-9	Tetrahydrofuran (THF)	ND	0.93	ND	0.31	
71-55-6	1,1,1-Trichloroethane	ND	0.93	ND	0.17	
56-23-5	Carbon Tetrachloride	ND	0.93	ND	0.15	
110-82-7	Cyclohexane	<b>2.4</b>	1.9	<b>0.69</b>	0.54	
78-87-5	1,2-Dichloropropane	ND	0.93	ND	0.20	
75-27-4	Bromodichloromethane	ND	0.93	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.93	ND	0.26	
80-62-6	Methyl Methacrylate	ND	1.9	ND	0.45	
142-82-5	n-Heptane	<b>5.7</b>	0.93	<b>1.4</b>	0.23	
10061-01-5	cis-1,3-Dichloropropene	ND	0.93	ND	0.20	
108-10-1	4-Methyl-2-pentanone	ND	0.93	ND	0.23	
10061-02-6	trans-1,3-Dichloropropene	ND	0.93	ND	0.20	
108-88-3	Toluene	<b>34</b>	0.93	<b>9.1</b>	0.25	
591-78-6	2-Hexanone	ND	0.93	ND	0.23	
124-48-1	Dibromochloromethane	ND	0.93	ND	0.11	
106-93-4	1,2-Dibromoethane	ND	0.93	ND	0.12	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 02-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-002

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00687

Date Collected: 3/5/12  
Date Received: 3/12/12  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.86 Final Pressure (psig): 3.52

Canister Dilution Factor: 1.85

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	5.0	0.93	1.0	0.19	
111-65-9	n-Octane	2.0	0.93	0.43	0.20	
108-90-7	Chlorobenzene	ND	0.93	ND	0.20	
179601-23-1	m,p-Xylenes	22	1.9	5.1	0.43	
75-25-2	Bromoform	ND	0.93	ND	0.089	
100-42-5	Styrene	2.3	0.93	0.54	0.22	
95-47-6	o-Xylene	7.7	0.93	1.8	0.21	
111-84-2	n-Nonane	2.1	0.93	0.40	0.18	
98-82-8	Cumene	ND	0.93	ND	0.19	
80-56-8	alpha-Pinene	11	0.93	2.0	0.17	
103-65-1	n-Propylbenzene	1.7	0.93	0.34	0.19	
622-96-8	4-Ethyltoluene	2.6	0.93	0.52	0.19	
108-67-8	1,3,5-Trimethylbenzene	3.0	0.93	0.61	0.19	
95-63-6	1,2,4-Trimethylbenzene	8.8	0.93	1.8	0.19	
100-44-7	Benzyl Chloride	ND	0.93	ND	0.18	
541-73-1	1,3-Dichlorobenzene	ND	0.93	ND	0.15	
106-46-7	1,4-Dichlorobenzene	ND	0.93	ND	0.15	
95-50-1	1,2-Dichlorobenzene	ND	0.93	ND	0.15	
5989-27-5	d-Limonene	24	0.93	4.3	0.17	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.93	ND	0.096	
120-82-1	1,2,4-Trichlorobenzene	ND	0.93	ND	0.12	
87-68-3	Hexachlorobutadiene	ND	0.93	ND	0.087	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 03-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-003

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00178

Date Collected: 3/5/12  
Date Received: 3/12/12  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.80 Final Pressure (psig): 3.51

Canister Dilution Factor: 1.84

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	78	0.92	45	0.53	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.1	0.92	0.42	0.19	
74-87-3	Chloromethane	ND	0.92	ND	0.45	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.92	ND	0.13	
106-99-0	1,3-Butadiene	ND	0.92	ND	0.42	
74-83-9	Bromomethane	ND	0.92	ND	0.24	
75-00-3	Chloroethane	ND	0.92	ND	0.35	
64-17-5	Ethanol	460	9.2	240	4.9	
75-05-8	Acetonitrile	ND	0.92	ND	0.55	
107-02-8	Acrolein	ND	3.7	ND	1.6	
67-64-1	Acetone	72	9.2	30	3.9	
75-69-4	Trichlorofluoromethane	1.1	0.92	0.20	0.16	
67-63-0	2-Propanol (Isopropyl Alcohol)	27	1.8	11	0.75	
107-13-1	Acrylonitrile	ND	0.92	ND	0.42	
75-35-4	1,1-Dichloroethene	ND	0.92	ND	0.23	
75-09-2	Methylene Chloride	ND	0.92	ND	0.26	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.92	ND	0.29	
76-13-1	Trichlorotrifluoroethane	ND	0.92	ND	0.12	
75-15-0	Carbon Disulfide	ND	9.2	ND	3.0	
156-60-5	trans-1,2-Dichloroethene	ND	0.92	ND	0.23	
75-34-3	1,1-Dichloroethane	ND	0.92	ND	0.23	
1634-04-4	Methyl tert-Butyl Ether	ND	0.92	ND	0.26	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 03-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-003

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AS00178

**Date Collected:** 3/5/12  
**Date Received:** 3/12/12  
**Date Analyzed:** 3/21/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -4.80      Final Pressure (psig): 3.51

Canister Dilution Factor: 1.84

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	9.2	ND	2.6	
78-93-3	2-Butanone (MEK)	ND	9.2	ND	3.1	
156-59-2	cis-1,2-Dichloroethene	ND	0.92	ND	0.23	
141-78-6	Ethyl Acetate	<b>25</b>	1.8	<b>7.0</b>	0.51	
110-54-3	n-Hexane	ND	0.92	ND	0.26	
67-66-3	Chloroform	ND	0.92	ND	0.19	
109-99-9	Tetrahydrofuran (THF)	ND	0.92	ND	0.31	
71-55-6	1,1,1-Trichloroethane	ND	0.92	ND	0.17	
56-23-5	Carbon Tetrachloride	ND	0.92	ND	0.15	
110-82-7	Cyclohexane	ND	1.8	ND	0.53	
78-87-5	1,2-Dichloropropane	ND	0.92	ND	0.20	
75-27-4	Bromodichloromethane	ND	0.92	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.92	ND	0.26	
80-62-6	Methyl Methacrylate	ND	1.8	ND	0.45	
142-82-5	n-Heptane	<b>20</b>	0.92	<b>4.8</b>	0.22	
10061-01-5	cis-1,3-Dichloropropene	ND	0.92	ND	0.20	
108-10-1	4-Methyl-2-pentanone	ND	0.92	ND	0.22	
10061-02-6	trans-1,3-Dichloropropene	ND	0.92	ND	0.20	
108-88-3	Toluene	<b>14</b>	0.92	<b>3.6</b>	0.24	
591-78-6	2-Hexanone	ND	0.92	ND	0.22	
124-48-1	Dibromochloromethane	ND	0.92	ND	0.11	
106-93-4	1,2-Dibromoethane	ND	0.92	ND	0.12	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 03-IF-030512  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1200947  
**CAS Sample ID:** P1200947-003

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AS00178

**Date Collected:** 3/5/12  
**Date Received:** 3/12/12  
**Date Analyzed:** 3/21/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -4.80      Final Pressure (psig): 3.51

Canister Dilution Factor: 1.84

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	7.3	0.92	1.5	0.19	
111-65-9	n-Octane	ND	0.92	ND	0.20	
108-90-7	Chlorobenzene	ND	0.92	ND	0.20	
179601-23-1	m,p-Xylenes	1.9	1.8	0.44	0.42	
75-25-2	Bromoform	ND	0.92	ND	0.089	
100-42-5	Styrene	0.98	0.92	0.23	0.22	
95-47-6	o-Xylene	ND	0.92	ND	0.21	
111-84-2	n-Nonane	3.6	0.92	0.70	0.18	M1
98-82-8	Cumene	ND	0.92	ND	0.19	
80-56-8	alpha-Pinene	11	0.92	1.9	0.17	
103-65-1	n-Propylbenzene	ND	0.92	ND	0.19	
622-96-8	4-Ethyltoluene	ND	0.92	ND	0.19	
108-67-8	1,3,5-Trimethylbenzene	ND	0.92	ND	0.19	
95-63-6	1,2,4-Trimethylbenzene	ND	0.92	ND	0.19	
100-44-7	Benzyl Chloride	ND	0.92	ND	0.18	
541-73-1	1,3-Dichlorobenzene	ND	0.92	ND	0.15	
106-46-7	1,4-Dichlorobenzene	ND	0.92	ND	0.15	
95-50-1	1,2-Dichlorobenzene	ND	0.92	ND	0.15	
5989-27-5	d-Limonene	42	0.92	7.5	0.17	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.92	ND	0.095	
120-82-1	1,2,4-Trichlorobenzene	ND	0.92	ND	0.12	
87-68-3	Hexachlorobutadiene	ND	0.92	ND	0.086	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

M1 = Matrix interference due to coelution with a non-target compound; results may be biased high.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 05-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-004

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01198

Date Collected: 3/7/12  
Date Received: 3/12/12  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.39 Final Pressure (psig): 3.53

Canister Dilution Factor: 1.61

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	8.7	0.81	5.1	0.47	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.1	0.81	0.43	0.16	
74-87-3	Chloromethane	ND	0.81	ND	0.39	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.81	ND	0.12	
106-99-0	1,3-Butadiene	0.99	0.81	0.45	0.36	
74-83-9	Bromomethane	ND	0.81	ND	0.21	
75-00-3	Chloroethane	ND	0.81	ND	0.31	
64-17-5	Ethanol	730	8.1	390	4.3	
75-05-8	Acetonitrile	ND	0.81	ND	0.48	
107-02-8	Acrolein	ND	3.2	ND	1.4	
67-64-1	Acetone	75	8.1	32	3.4	
75-69-4	Trichlorofluoromethane	1.4	0.81	0.24	0.14	
67-63-0	2-Propanol (Isopropyl Alcohol)	22	1.6	9.1	0.66	
107-13-1	Acrylonitrile	ND	0.81	ND	0.37	
75-35-4	1,1-Dichloroethene	ND	0.81	ND	0.20	
75-09-2	Methylene Chloride	2.5	0.81	0.73	0.23	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.81	ND	0.26	
76-13-1	Trichlorotrifluoroethane	ND	0.81	ND	0.11	
75-15-0	Carbon Disulfide	ND	8.1	ND	2.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.81	ND	0.20	
75-34-3	1,1-Dichloroethane	ND	0.81	ND	0.20	
1634-04-4	Methyl tert-Butyl Ether	ND	0.81	ND	0.22	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 05-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-004

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01198

**Date Collected:** 3/7/12  
**Date Received:** 3/12/12  
**Date Analyzed:** 3/21/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -3.39      Final Pressure (psig): 3.53

Canister Dilution Factor: 1.61

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	8.1	ND	2.3	
78-93-3	2-Butanone (MEK)	ND	8.1	ND	2.7	
156-59-2	cis-1,2-Dichloroethene	ND	0.81	ND	0.20	
141-78-6	Ethyl Acetate	6.9	1.6	1.9	0.45	
110-54-3	n-Hexane	19	0.81	5.4	0.23	
67-66-3	Chloroform	5.2	0.81	1.1	0.16	
109-99-9	Tetrahydrofuran (THF)	ND	0.81	ND	0.27	
71-55-6	1,1,1-Trichloroethane	ND	0.81	ND	0.15	
56-23-5	Carbon Tetrachloride	0.97	0.81	0.15	0.13	
110-82-7	Cyclohexane	6.8	1.6	2.0	0.47	
78-87-5	1,2-Dichloropropane	ND	0.81	ND	0.17	
75-27-4	Bromodichloromethane	0.98	0.81	0.15	0.12	
123-91-1	1,4-Dioxane	ND	0.81	ND	0.22	
80-62-6	Methyl Methacrylate	ND	1.6	ND	0.39	
142-82-5	n-Heptane	6.1	0.81	1.5	0.20	
10061-01-5	cis-1,3-Dichloropropene	ND	0.81	ND	0.18	
108-10-1	4-Methyl-2-pentanone	ND	0.81	ND	0.20	
10061-02-6	trans-1,3-Dichloropropene	ND	0.81	ND	0.18	
108-88-3	Toluene	89	0.81	24	0.21	
591-78-6	2-Hexanone	ND	0.81	ND	0.20	
124-48-1	Dibromochloromethane	ND	0.81	ND	0.095	
106-93-4	1,2-Dibromoethane	ND	0.81	ND	0.10	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 05-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-004

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01198

Date Collected: 3/7/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/21/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.39 Final Pressure (psig): 3.53

Canister Dilution Factor: 1.61

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.81	ND	0.17	
111-65-9	n-Octane	2.3	0.81	0.49	0.17	
108-90-7	Chlorobenzene	ND	0.81	ND	0.17	
179601-23-1	m,p-Xylenes	50	1.6	12	0.37	
75-25-2	Bromoform	ND	0.81	ND	0.078	
100-42-5	Styrene	2.4	0.81	0.56	0.19	
95-47-6	o-Xylene	16	0.81	3.6	0.19	
111-84-2	n-Nonane	2.1	0.81	0.40	0.15	
98-82-8	Cumene	ND	0.81	ND	0.16	
80-56-8	alpha-Pinene	1.9	0.81	0.35	0.14	
103-65-1	n-Propylbenzene	2.8	0.81	0.58	0.16	
622-96-8	4-Ethyltoluene	5.4	0.81	1.1	0.16	
108-67-8	1,3,5-Trimethylbenzene	5.5	0.81	1.1	0.16	
95-63-6	1,2,4-Trimethylbenzene	17	0.81	3.4	0.16	
100-44-7	Benzyl Chloride	ND	0.81	ND	0.16	
541-73-1	1,3-Dichlorobenzene	ND	0.81	ND	0.13	
106-46-7	1,4-Dichlorobenzene	ND	0.81	ND	0.13	
95-50-1	1,2-Dichlorobenzene	ND	0.81	ND	0.13	
5989-27-5	d-Limonene	11	0.81	1.9	0.14	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.81	ND	0.083	
120-82-1	1,2,4-Trichlorobenzene	ND	0.81	ND	0.11	
87-68-3	Hexachlorobutadiene	ND	0.81	ND	0.075	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 06-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-005

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AS00137

**Date Collected:** 3/7/12  
**Date Received:** 3/12/12  
**Date Analyzed:** 3/21 - 3/22/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.050 Liter(s)

Initial Pressure (psig): -2.12      Final Pressure (psig): 3.74

Canister Dilution Factor: 1.47

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	7.9	0.74	4.6	0.43	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.1	0.74	0.42	0.15	
74-87-3	Chloromethane	ND	0.74	ND	0.36	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.74	ND	0.11	
106-99-0	1,3-Butadiene	ND	0.74	ND	0.33	
74-83-9	Bromomethane	ND	0.74	ND	0.19	
75-00-3	Chloroethane	ND	0.74	ND	0.28	
64-17-5	Ethanol	970	150	510	78	D
75-05-8	Acetonitrile	ND	0.74	ND	0.44	
107-02-8	Acrolein	ND	2.9	ND	1.3	
67-64-1	Acetone	87	7.4	37	3.1	
75-69-4	Trichlorofluoromethane	2.2	0.74	0.39	0.13	
67-63-0	2-Propanol (Isopropyl Alcohol)	38	1.5	15	0.60	
107-13-1	Acrylonitrile	ND	0.74	ND	0.34	
75-35-4	1,1-Dichloroethene	ND	0.74	ND	0.19	
75-09-2	Methylene Chloride	ND	0.74	ND	0.21	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.74	ND	0.23	
76-13-1	Trichlorotrifluoroethane	ND	0.74	ND	0.096	
75-15-0	Carbon Disulfide	ND	7.4	ND	2.4	
156-60-5	trans-1,2-Dichloroethene	ND	0.74	ND	0.19	
75-34-3	1,1-Dichloroethane	ND	0.74	ND	0.18	
1634-04-4	Methyl tert-Butyl Ether	ND	0.74	ND	0.20	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 06-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-005

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00137

Date Collected: 3/7/12  
Date Received: 3/12/12  
Date Analyzed: 3/21 - 3/22/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.050 Liter(s)

Initial Pressure (psig): -2.12 Final Pressure (psig): 3.74

Canister Dilution Factor: 1.47

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	7.4	ND	2.1	
78-93-3	2-Butanone (MEK)	ND	7.4	ND	2.5	
156-59-2	cis-1,2-Dichloroethene	ND	0.74	ND	0.19	
141-78-6	Ethyl Acetate	<b>40</b>	1.5	<b>11</b>	0.41	
110-54-3	n-Hexane	<b>5.4</b>	0.74	<b>1.5</b>	0.21	
67-66-3	Chloroform	<b>2.1</b>	0.74	<b>0.43</b>	0.15	
109-99-9	Tetrahydrofuran (THF)	ND	0.74	ND	0.25	
71-55-6	1,1,1-Trichloroethane	ND	0.74	ND	0.13	
56-23-5	Carbon Tetrachloride	ND	0.74	ND	0.12	
110-82-7	Cyclohexane	<b>2.3</b>	1.5	<b>0.68</b>	0.43	
78-87-5	1,2-Dichloropropane	ND	0.74	ND	0.16	
75-27-4	Bromodichloromethane	<b>0.76</b>	0.74	<b>0.11</b>	0.11	
123-91-1	1,4-Dioxane	ND	0.74	ND	0.20	
80-62-6	Methyl Methacrylate	ND	1.5	ND	0.36	
142-82-5	n-Heptane	<b>2.4</b>	0.74	<b>0.58</b>	0.18	
10061-01-5	cis-1,3-Dichloropropene	ND	0.74	ND	0.16	
108-10-1	4-Methyl-2-pentanone	ND	0.74	ND	0.18	
10061-02-6	trans-1,3-Dichloropropene	ND	0.74	ND	0.16	
108-88-3	Toluene	<b>20</b>	0.74	<b>5.2</b>	0.20	
591-78-6	2-Hexanone	ND	0.74	ND	0.18	
124-48-1	Dibromochloromethane	ND	0.74	ND	0.086	
106-93-4	1,2-Dibromoethane	ND	0.74	ND	0.096	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 06-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-005

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00137

Date Collected: 3/7/12  
Date Received: 3/12/12  
Date Analyzed: 3/21 - 3/22/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.050 Liter(s)

Initial Pressure (psig): -2.12 Final Pressure (psig): 3.74

Canister Dilution Factor: 1.47

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	14	0.74	2.9	0.15	
111-65-9	n-Octane	1.0	0.74	0.22	0.16	
108-90-7	Chlorobenzene	ND	0.74	ND	0.16	
179601-23-1	m,p-Xylenes	9.8	1.5	2.2	0.34	
75-25-2	Bromoform	ND	0.74	ND	0.071	
100-42-5	Styrene	0.93	0.74	0.22	0.17	
95-47-6	o-Xylene	3.3	0.74	0.76	0.17	
111-84-2	n-Nonane	0.79	0.74	0.15	0.14	
98-82-8	Cumene	ND	0.74	ND	0.15	
80-56-8	alpha-Pinene	14	0.74	2.5	0.13	
103-65-1	n-Propylbenzene	ND	0.74	ND	0.15	
622-96-8	4-Ethyltoluene	1.2	0.74	0.25	0.15	
108-67-8	1,3,5-Trimethylbenzene	1.4	0.74	0.28	0.15	
95-63-6	1,2,4-Trimethylbenzene	4.5	0.74	0.91	0.15	
100-44-7	Benzyl Chloride	ND	0.74	ND	0.14	
541-73-1	1,3-Dichlorobenzene	ND	0.74	ND	0.12	
106-46-7	1,4-Dichlorobenzene	ND	0.74	ND	0.12	
95-50-1	1,2-Dichlorobenzene	ND	0.74	ND	0.12	
5989-27-5	d-Limonene	26	0.74	4.7	0.13	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.74	ND	0.076	
120-82-1	1,2,4-Trichlorobenzene	ND	0.74	ND	0.099	
87-68-3	Hexachlorobutadiene	ND	0.74	ND	0.069	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 07-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-006

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01570

Date Collected: 3/7/12  
Date Received: 3/12/12  
Date Analyzed: 3/21 - 3/22/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.050 Liter(s)

Initial Pressure (psig): -2.36 Final Pressure (psig): 3.72

Canister Dilution Factor: 1.49

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	ND	0.75	ND	0.43	
75-71-8	Dichlorodifluoromethane (CFC 12)	<b>2.1</b>	0.75	<b>0.43</b>	0.15	
74-87-3	Chloromethane	ND	0.75	ND	0.36	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.75	ND	0.11	
106-99-0	1,3-Butadiene	ND	0.75	ND	0.34	
74-83-9	Bromomethane	ND	0.75	ND	0.19	
75-00-3	Chloroethane	ND	0.75	ND	0.28	
64-17-5	Ethanol	<b>1,000</b>	150	<b>530</b>	79	<b>D</b>
75-05-8	Acetonitrile	ND	0.75	ND	0.44	
107-02-8	Acrolein	ND	3.0	ND	1.3	
67-64-1	Acetone	<b>47</b>	7.5	<b>20</b>	3.1	
75-69-4	Trichlorofluoromethane	<b>1.3</b>	0.75	<b>0.22</b>	0.13	
67-63-0	2-Propanol (Isopropyl Alcohol)	<b>5.2</b>	1.5	<b>2.1</b>	0.61	
107-13-1	Acrylonitrile	ND	0.75	ND	0.34	
75-35-4	1,1-Dichloroethene	ND	0.75	ND	0.19	
75-09-2	Methylene Chloride	<b>170</b>	15	<b>50</b>	4.3	<b>D</b>
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.75	ND	0.24	
76-13-1	Trichlorotrifluoroethane	ND	0.75	ND	0.097	
75-15-0	Carbon Disulfide	ND	7.5	ND	2.4	
156-60-5	trans-1,2-Dichloroethene	ND	0.75	ND	0.19	
75-34-3	1,1-Dichloroethane	ND	0.75	ND	0.18	
1634-04-4	Methyl tert-Butyl Ether	ND	0.75	ND	0.21	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 07-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-006

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01570

**Date Collected:** 3/7/12  
**Date Received:** 3/12/12  
**Date Analyzed:** 3/21 - 3/22/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.050 Liter(s)

Initial Pressure (psig): -2.36      Final Pressure (psig): 3.72

Canister Dilution Factor: 1.49

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	7.5	ND	2.1	
78-93-3	2-Butanone (MEK)	ND	7.5	ND	2.5	
156-59-2	cis-1,2-Dichloroethene	ND	0.75	ND	0.19	
141-78-6	Ethyl Acetate	12	1.5	3.4	0.41	
110-54-3	n-Hexane	ND	0.75	ND	0.21	
67-66-3	Chloroform	3.0	0.75	0.61	0.15	
109-99-9	Tetrahydrofuran (THF)	0.89	0.75	0.30	0.25	
71-55-6	1,1,1-Trichloroethane	ND	0.75	ND	0.14	
56-23-5	Carbon Tetrachloride	ND	0.75	ND	0.12	
110-82-7	Cyclohexane	ND	1.5	ND	0.43	
78-87-5	1,2-Dichloropropane	ND	0.75	ND	0.16	
75-27-4	Bromodichloromethane	2.2	0.75	0.32	0.11	
123-91-1	1,4-Dioxane	ND	0.75	ND	0.21	
80-62-6	Methyl Methacrylate	ND	1.5	ND	0.36	
142-82-5	n-Heptane	ND	0.75	ND	0.18	
10061-01-5	cis-1,3-Dichloropropene	ND	0.75	ND	0.16	
108-10-1	4-Methyl-2-pentanone	ND	0.75	ND	0.18	
10061-02-6	trans-1,3-Dichloropropene	ND	0.75	ND	0.16	
108-88-3	Toluene	9.2	0.75	2.5	0.20	
591-78-6	2-Hexanone	ND	0.75	ND	0.18	
124-48-1	Dibromochloromethane	1.3	0.75	0.16	0.087	
106-93-4	1,2-Dibromoethane	ND	0.75	ND	0.097	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 07-IF-030712  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1200947  
**CAS Sample ID:** P1200947-006

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01570

**Date Collected:** 3/7/12  
**Date Received:** 3/12/12  
**Date Analyzed:** 3/21 - 3/22/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.050 Liter(s)

Initial Pressure (psig): -2.36      Final Pressure (psig): 3.72

Canister Dilution Factor: 1.49

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	1.5	0.75	0.31	0.16	
111-65-9	n-Octane	ND	0.75	ND	0.16	
108-90-7	Chlorobenzene	ND	0.75	ND	0.16	
179601-23-1	m,p-Xylenes	1.7	1.5	0.38	0.34	
75-25-2	Bromoform	ND	0.75	ND	0.072	
100-42-5	Styrene	ND	0.75	ND	0.18	
95-47-6	o-Xylene	ND	0.75	ND	0.17	
111-84-2	n-Nonane	ND	0.75	ND	0.14	
98-82-8	Cumene	ND	0.75	ND	0.15	
80-56-8	alpha-Pinene	2.5	0.75	0.44	0.13	
103-65-1	n-Propylbenzene	ND	0.75	ND	0.15	
622-96-8	4-Ethyltoluene	ND	0.75	ND	0.15	
108-67-8	1,3,5-Trimethylbenzene	ND	0.75	ND	0.15	
95-63-6	1,2,4-Trimethylbenzene	1.4	0.75	0.28	0.15	
100-44-7	Benzyl Chloride	ND	0.75	ND	0.14	
541-73-1	1,3-Dichlorobenzene	ND	0.75	ND	0.12	
106-46-7	1,4-Dichlorobenzene	ND	0.75	ND	0.12	
95-50-1	1,2-Dichlorobenzene	ND	0.75	ND	0.12	
5989-27-5	d-Limonene	12	0.75	2.1	0.13	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.75	ND	0.077	
120-82-1	1,2,4-Trichlorobenzene	ND	0.75	ND	0.10	
87-68-3	Hexachlorobutadiene	ND	0.75	ND	0.070	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 08-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-007

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00184

Date Collected: 3/7/12  
Date Received: 3/12/12  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.87 Final Pressure (psig): 3.58

Canister Dilution Factor: 1.55

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	1.2	0.78	0.67	0.45	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.1	0.78	0.42	0.16	
74-87-3	Chloromethane	ND	0.78	ND	0.38	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.78	ND	0.11	
106-99-0	1,3-Butadiene	ND	0.78	ND	0.35	
74-83-9	Bromomethane	ND	0.78	ND	0.20	
75-00-3	Chloroethane	ND	0.78	ND	0.29	
64-17-5	Ethanol	310	7.8	160	4.1	
75-05-8	Acetonitrile	ND	0.78	ND	0.46	
107-02-8	Acrolein	ND	3.1	ND	1.4	
67-64-1	Acetone	30	7.8	13	3.3	
75-69-4	Trichlorofluoromethane	3.6	0.78	0.65	0.14	
67-63-0	2-Propanol (Isopropyl Alcohol)	5.6	1.6	2.3	0.63	
107-13-1	Acrylonitrile	ND	0.78	ND	0.36	
75-35-4	1,1-Dichloroethene	ND	0.78	ND	0.20	
75-09-2	Methylene Chloride	ND	0.78	ND	0.22	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.78	ND	0.25	
76-13-1	Trichlorotrifluoroethane	ND	0.78	ND	0.10	
75-15-0	Carbon Disulfide	ND	7.8	ND	2.5	
156-60-5	trans-1,2-Dichloroethene	ND	0.78	ND	0.20	
75-34-3	1,1-Dichloroethane	ND	0.78	ND	0.19	
1634-04-4	Methyl tert-Butyl Ether	ND	0.78	ND	0.22	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 08-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-007

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AS00184

**Date Collected:** 3/7/12  
**Date Received:** 3/12/12  
**Date Analyzed:** 3/21/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -2.87      Final Pressure (psig): 3.58

Canister Dilution Factor: 1.55

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	7.8	ND	2.2	
78-93-3	2-Butanone (MEK)	ND	7.8	ND	2.6	
156-59-2	cis-1,2-Dichloroethene	ND	0.78	ND	0.20	
141-78-6	Ethyl Acetate	12	1.6	3.4	0.43	
110-54-3	n-Hexane	ND	0.78	ND	0.22	
67-66-3	Chloroform	ND	0.78	ND	0.16	
109-99-9	Tetrahydrofuran (THF)	ND	0.78	ND	0.26	
71-55-6	1,1,1-Trichloroethane	ND	0.78	ND	0.14	
56-23-5	Carbon Tetrachloride	ND	0.78	ND	0.12	
110-82-7	Cyclohexane	2.7	1.6	0.79	0.45	
78-87-5	1,2-Dichloropropane	ND	0.78	ND	0.17	
75-27-4	Bromodichloromethane	ND	0.78	ND	0.12	
123-91-1	1,4-Dioxane	ND	0.78	ND	0.22	
80-62-6	Methyl Methacrylate	ND	1.6	ND	0.38	
142-82-5	n-Heptane	ND	0.78	ND	0.19	
10061-01-5	cis-1,3-Dichloropropene	ND	0.78	ND	0.17	
108-10-1	4-Methyl-2-pentanone	ND	0.78	ND	0.19	
10061-02-6	trans-1,3-Dichloropropene	ND	0.78	ND	0.17	
108-88-3	Toluene	7.1	0.78	1.9	0.21	
591-78-6	2-Hexanone	ND	0.78	ND	0.19	
124-48-1	Dibromochloromethane	ND	0.78	ND	0.091	
106-93-4	1,2-Dibromoethane	ND	0.78	ND	0.10	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 08-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-007

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AS00184

**Date Collected:** 3/7/12  
**Date Received:** 3/12/12  
**Date Analyzed:** 3/21/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -2.87      Final Pressure (psig): 3.58

Canister Dilution Factor: 1.55

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	1.5	0.78	0.32	0.16	
111-65-9	n-Octane	ND	0.78	ND	0.17	
108-90-7	Chlorobenzene	ND	0.78	ND	0.17	
179601-23-1	m,p-Xylenes	ND	1.6	ND	0.36	
75-25-2	Bromoform	ND	0.78	ND	0.075	
100-42-5	Styrene	0.90	0.78	0.21	0.18	
95-47-6	o-Xylene	ND	0.78	ND	0.18	
111-84-2	n-Nonane	ND	0.78	ND	0.15	
98-82-8	Cumene	ND	0.78	ND	0.16	
80-56-8	alpha-Pinene	1.2	0.78	0.21	0.14	
103-65-1	n-Propylbenzene	ND	0.78	ND	0.16	
622-96-8	4-Ethyltoluene	ND	0.78	ND	0.16	
108-67-8	1,3,5-Trimethylbenzene	ND	0.78	ND	0.16	
95-63-6	1,2,4-Trimethylbenzene	ND	0.78	ND	0.16	
100-44-7	Benzyl Chloride	ND	0.78	ND	0.15	
541-73-1	1,3-Dichlorobenzene	ND	0.78	ND	0.13	
106-46-7	1,4-Dichlorobenzene	ND	0.78	ND	0.13	
95-50-1	1,2-Dichlorobenzene	ND	0.78	ND	0.13	
5989-27-5	d-Limonene	64	0.78	11	0.14	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.78	ND	0.080	
120-82-1	1,2,4-Trichlorobenzene	ND	0.78	ND	0.10	
87-68-3	Hexachlorobutadiene	ND	0.78	ND	0.073	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 09-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-008

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00553

Date Collected: 3/5/12  
Date Received: 3/12/12  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.65 Final Pressure (psig): 3.56

Canister Dilution Factor: 1.82

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	1.6	0.91	0.91	0.53	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.7	0.91	0.55	0.18	
74-87-3	Chloromethane	ND	0.91	ND	0.44	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.91	ND	0.13	
106-99-0	1,3-Butadiene	ND	0.91	ND	0.41	
74-83-9	Bromomethane	ND	0.91	ND	0.23	
75-00-3	Chloroethane	ND	0.91	ND	0.34	
64-17-5	Ethanol	120	9.1	62	4.8	
75-05-8	Acetonitrile	ND	0.91	ND	0.54	
107-02-8	Acrolein	ND	3.6	ND	1.6	
67-64-1	Acetone	30	9.1	13	3.8	
75-69-4	Trichlorofluoromethane	2.3	0.91	0.42	0.16	
67-63-0	2-Propanol (Isopropyl Alcohol)	6.7	1.8	2.7	0.74	
107-13-1	Acrylonitrile	ND	0.91	ND	0.42	
75-35-4	1,1-Dichloroethene	ND	0.91	ND	0.23	
75-09-2	Methylene Chloride	1.2	0.91	0.34	0.26	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.91	ND	0.29	
76-13-1	Trichlorotrifluoroethane	ND	0.91	ND	0.12	
75-15-0	Carbon Disulfide	ND	9.1	ND	2.9	
156-60-5	trans-1,2-Dichloroethene	ND	0.91	ND	0.23	
75-34-3	1,1-Dichloroethane	ND	0.91	ND	0.22	
1634-04-4	Methyl tert-Butyl Ether	ND	0.91	ND	0.25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 09-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-008

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00553

Date Collected: 3/5/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/21/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.65      Final Pressure (psig): 3.56

Canister Dilution Factor: 1.82

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	9.1	ND	2.6	
78-93-3	2-Butanone (MEK)	ND	9.1	ND	3.1	
156-59-2	cis-1,2-Dichloroethene	ND	0.91	ND	0.23	
141-78-6	Ethyl Acetate	<b>34</b>	1.8	<b>9.4</b>	0.51	
110-54-3	n-Hexane	ND	0.91	ND	0.26	
67-66-3	Chloroform	ND	0.91	ND	0.19	
109-99-9	Tetrahydrofuran (THF)	ND	0.91	ND	0.31	
71-55-6	1,1,1-Trichloroethane	ND	0.91	ND	0.17	
56-23-5	Carbon Tetrachloride	ND	0.91	ND	0.14	
110-82-7	Cyclohexane	<b>2.7</b>	1.8	<b>0.77</b>	0.53	
78-87-5	1,2-Dichloropropane	ND	0.91	ND	0.20	
75-27-4	Bromodichloromethane	ND	0.91	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.91	ND	0.25	
80-62-6	Methyl Methacrylate	<b>1.9</b>	1.8	<b>0.47</b>	0.44	
142-82-5	n-Heptane	ND	0.91	ND	0.22	
10061-01-5	cis-1,3-Dichloropropene	ND	0.91	ND	0.20	
108-10-1	4-Methyl-2-pentanone	ND	0.91	ND	0.22	
10061-02-6	trans-1,3-Dichloropropene	ND	0.91	ND	0.20	
108-88-3	Toluene	<b>14</b>	0.91	<b>3.8</b>	0.24	
591-78-6	2-Hexanone	ND	0.91	ND	0.22	
124-48-1	Dibromochloromethane	ND	0.91	ND	0.11	
106-93-4	1,2-Dibromoethane	ND	0.91	ND	0.12	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 09-IF-030512  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1200947  
**CAS Sample ID:** P1200947-008

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC00553

**Date Collected:** 3/5/12  
**Date Received:** 3/12/12  
**Date Analyzed:** 3/21/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -4.65      Final Pressure (psig): 3.56

Canister Dilution Factor: 1.82

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	1.2	0.91	0.25	0.19	
111-65-9	n-Octane	ND	0.91	ND	0.19	
108-90-7	Chlorobenzene	ND	0.91	ND	0.20	
179601-23-1	m,p-Xylenes	ND	1.8	ND	0.42	
75-25-2	Bromoform	ND	0.91	ND	0.088	
100-42-5	Styrene	1.3	0.91	0.31	0.21	
95-47-6	o-Xylene	ND	0.91	ND	0.21	
111-84-2	n-Nonane	ND	0.91	ND	0.17	
98-82-8	Cumene	ND	0.91	ND	0.19	
80-56-8	alpha-Pinene	ND	0.91	ND	0.16	
103-65-1	n-Propylbenzene	ND	0.91	ND	0.19	
622-96-8	4-Ethyltoluene	ND	0.91	ND	0.19	
108-67-8	1,3,5-Trimethylbenzene	ND	0.91	ND	0.19	
95-63-6	1,2,4-Trimethylbenzene	ND	0.91	ND	0.19	
100-44-7	Benzyl Chloride	ND	0.91	ND	0.18	
541-73-1	1,3-Dichlorobenzene	ND	0.91	ND	0.15	
106-46-7	1,4-Dichlorobenzene	ND	0.91	ND	0.15	
95-50-1	1,2-Dichlorobenzene	ND	0.91	ND	0.15	
5989-27-5	d-Limonene	3.6	0.91	0.65	0.16	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.91	ND	0.094	
120-82-1	1,2,4-Trichlorobenzene	ND	0.91	ND	0.12	
87-68-3	Hexachlorobutadiene	ND	0.91	ND	0.085	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P120321-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	ND	0.50	ND	0.29	
75-71-8	Dichlorodifluoromethane (CFC 12)	ND	0.50	ND	0.10	
74-87-3	Chloromethane	ND	0.50	ND	0.24	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.50	ND	0.072	
106-99-0	1,3-Butadiene	ND	0.50	ND	0.23	
74-83-9	Bromomethane	ND	0.50	ND	0.13	
75-00-3	Chloroethane	ND	0.50	ND	0.19	
64-17-5	Ethanol	ND	5.0	ND	2.7	
75-05-8	Acetonitrile	ND	0.50	ND	0.30	
107-02-8	Acrolein	ND	2.0	ND	0.87	
67-64-1	Acetone	ND	5.0	ND	2.1	
75-69-4	Trichlorofluoromethane	ND	0.50	ND	0.089	
67-63-0	2-Propanol (Isopropyl Alcohol)	ND	1.0	ND	0.41	
107-13-1	Acrylonitrile	ND	0.50	ND	0.23	
75-35-4	1,1-Dichloroethene	ND	0.50	ND	0.13	
75-09-2	Methylene Chloride	ND	0.50	ND	0.14	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.50	ND	0.16	
76-13-1	Trichlorotrifluoroethane	ND	0.50	ND	0.065	
75-15-0	Carbon Disulfide	ND	5.0	ND	1.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.50	ND	0.13	
75-34-3	1,1-Dichloroethane	ND	0.50	ND	0.12	
1634-04-4	Methyl tert-Butyl Ether	ND	0.50	ND	0.14	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P120321-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	5.0	ND	1.4	
78-93-3	2-Butanone (MEK)	ND	5.0	ND	1.7	
156-59-2	cis-1,2-Dichloroethene	ND	0.50	ND	0.13	
141-78-6	Ethyl Acetate	ND	1.0	ND	0.28	
110-54-3	n-Hexane	ND	0.50	ND	0.14	
67-66-3	Chloroform	ND	0.50	ND	0.10	
109-99-9	Tetrahydrofuran (THF)	ND	0.50	ND	0.17	
71-55-6	1,1,1-Trichloroethane	ND	0.50	ND	0.092	
56-23-5	Carbon Tetrachloride	ND	0.50	ND	0.080	
110-82-7	Cyclohexane	ND	1.0	ND	0.29	
78-87-5	1,2-Dichloropropane	ND	0.50	ND	0.11	
75-27-4	Bromodichloromethane	ND	0.50	ND	0.075	
123-91-1	1,4-Dioxane	ND	0.50	ND	0.14	
80-62-6	Methyl Methacrylate	ND	1.0	ND	0.24	
142-82-5	n-Heptane	ND	0.50	ND	0.12	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-10-1	4-Methyl-2-pentanone	ND	0.50	ND	0.12	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-88-3	Toluene	ND	0.50	ND	0.13	
591-78-6	2-Hexanone	ND	0.50	ND	0.12	
124-48-1	Dibromochloromethane	ND	0.50	ND	0.059	
106-93-4	1,2-Dibromoethane	ND	0.50	ND	0.065	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P120321-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.50	ND	0.11	
111-65-9	n-Octane	ND	0.50	ND	0.11	
108-90-7	Chlorobenzene	ND	0.50	ND	0.11	
179601-23-1	m,p-Xylenes	ND	1.0	ND	0.23	
75-25-2	Bromoform	ND	0.50	ND	0.048	
100-42-5	Styrene	ND	0.50	ND	0.12	
95-47-6	o-Xylene	ND	0.50	ND	0.12	
111-84-2	n-Nonane	ND	0.50	ND	0.095	
98-82-8	Cumene	ND	0.50	ND	0.10	
80-56-8	alpha-Pinene	ND	0.50	ND	0.090	
103-65-1	n-Propylbenzene	ND	0.50	ND	0.10	
622-96-8	4-Ethyltoluene	ND	0.50	ND	0.10	
108-67-8	1,3,5-Trimethylbenzene	ND	0.50	ND	0.10	
95-63-6	1,2,4-Trimethylbenzene	ND	0.50	ND	0.10	
100-44-7	Benzyl Chloride	ND	0.50	ND	0.097	
541-73-1	1,3-Dichlorobenzene	ND	0.50	ND	0.083	
106-46-7	1,4-Dichlorobenzene	ND	0.50	ND	0.083	
95-50-1	1,2-Dichlorobenzene	ND	0.50	ND	0.083	
5989-27-5	d-Limonene	ND	0.50	ND	0.090	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.50	ND	0.052	
120-82-1	1,2,4-Trichlorobenzene	ND	0.50	ND	0.067	
87-68-3	Hexachlorobutadiene	ND	0.50	ND	0.047	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P120322-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/22/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	ND	0.50	ND	0.29	
75-71-8	Dichlorodifluoromethane (CFC 12)	ND	0.50	ND	0.10	
74-87-3	Chloromethane	ND	0.50	ND	0.24	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.50	ND	0.072	
106-99-0	1,3-Butadiene	ND	0.50	ND	0.23	
74-83-9	Bromomethane	ND	0.50	ND	0.13	
75-00-3	Chloroethane	ND	0.50	ND	0.19	
64-17-5	Ethanol	ND	5.0	ND	2.7	
75-05-8	Acetonitrile	ND	0.50	ND	0.30	
107-02-8	Acrolein	ND	2.0	ND	0.87	
67-64-1	Acetone	ND	5.0	ND	2.1	
75-69-4	Trichlorofluoromethane	ND	0.50	ND	0.089	
67-63-0	2-Propanol (Isopropyl Alcohol)	ND	1.0	ND	0.41	
107-13-1	Acrylonitrile	ND	0.50	ND	0.23	
75-35-4	1,1-Dichloroethene	ND	0.50	ND	0.13	
75-09-2	Methylene Chloride	ND	0.50	ND	0.14	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.50	ND	0.16	
76-13-1	Trichlorotrifluoroethane	ND	0.50	ND	0.065	
75-15-0	Carbon Disulfide	ND	5.0	ND	1.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.50	ND	0.13	
75-34-3	1,1-Dichloroethane	ND	0.50	ND	0.12	
1634-04-4	Methyl tert-Butyl Ether	ND	0.50	ND	0.14	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P120322-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/22/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	5.0	ND	1.4	
78-93-3	2-Butanone (MEK)	ND	5.0	ND	1.7	
156-59-2	cis-1,2-Dichloroethene	ND	0.50	ND	0.13	
141-78-6	Ethyl Acetate	ND	1.0	ND	0.28	
110-54-3	n-Hexane	ND	0.50	ND	0.14	
67-66-3	Chloroform	ND	0.50	ND	0.10	
109-99-9	Tetrahydrofuran (THF)	ND	0.50	ND	0.17	
71-55-6	1,1,1-Trichloroethane	ND	0.50	ND	0.092	
56-23-5	Carbon Tetrachloride	ND	0.50	ND	0.080	
110-82-7	Cyclohexane	ND	1.0	ND	0.29	
78-87-5	1,2-Dichloropropane	ND	0.50	ND	0.11	
75-27-4	Bromodichloromethane	ND	0.50	ND	0.075	
123-91-1	1,4-Dioxane	ND	0.50	ND	0.14	
80-62-6	Methyl Methacrylate	ND	1.0	ND	0.24	
142-82-5	n-Heptane	ND	0.50	ND	0.12	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-10-1	4-Methyl-2-pentanone	ND	0.50	ND	0.12	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-88-3	Toluene	ND	0.50	ND	0.13	
591-78-6	2-Hexanone	ND	0.50	ND	0.12	
124-48-1	Dibromochloromethane	ND	0.50	ND	0.059	
106-93-4	1,2-Dibromoethane	ND	0.50	ND	0.065	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P120322-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/22/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.50	ND	0.11	
111-65-9	n-Octane	ND	0.50	ND	0.11	
108-90-7	Chlorobenzene	ND	0.50	ND	0.11	
179601-23-1	m,p-Xylenes	ND	1.0	ND	0.23	
75-25-2	Bromoform	ND	0.50	ND	0.048	
100-42-5	Styrene	ND	0.50	ND	0.12	
95-47-6	o-Xylene	ND	0.50	ND	0.12	
111-84-2	n-Nonane	ND	0.50	ND	0.095	
98-82-8	Cumene	ND	0.50	ND	0.10	
80-56-8	alpha-Pinene	ND	0.50	ND	0.090	
103-65-1	n-Propylbenzene	ND	0.50	ND	0.10	
622-96-8	4-Ethyltoluene	ND	0.50	ND	0.10	
108-67-8	1,3,5-Trimethylbenzene	ND	0.50	ND	0.10	
95-63-6	1,2,4-Trimethylbenzene	ND	0.50	ND	0.10	
100-44-7	Benzyl Chloride	ND	0.50	ND	0.097	
541-73-1	1,3-Dichlorobenzene	ND	0.50	ND	0.083	
106-46-7	1,4-Dichlorobenzene	ND	0.50	ND	0.083	
95-50-1	1,2-Dichlorobenzene	ND	0.50	ND	0.083	
5989-27-5	d-Limonene	ND	0.50	ND	0.090	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.50	ND	0.052	
120-82-1	1,2,4-Trichlorobenzene	ND	0.50	ND	0.067	
87-68-3	Hexachlorobutadiene	ND	0.50	ND	0.047	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister(s)  
 Test Notes:

Date(s) Collected: 3/5 - 3/7/12  
 Date(s) Received: 3/12/12  
 Date(s) Analyzed: 3/21 - 3/22/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		Percent Recovered	Percent Recovered	Percent Recovered		
Method Blank	P120321-MB	104	102	94	70-130	
Method Blank	P120322-MB	105	100	97	70-130	
Lab Control Sample	P120321-LCS	107	98	98	70-130	
Lab Control Sample	P120322-LCS	105	99	98	70-130	
01-IF-030512	P1200947-001	106	100	93	70-130	
02-IF-030512	P1200947-002	106	98	95	70-130	
02-IF-030512	P1200947-002DUP	106	98	96	70-130	
03-IF-030512	P1200947-003	105	98	93	70-130	
05-IF-030712	P1200947-004	104	99	96	70-130	
06-IF-030712	P1200947-005	106	99	97	70-130	
07-IF-030712	P1200947-006	102	100	96	70-130	
08-IF-030712	P1200947-007	104	98	95	70-130	
09-IF-030512	P1200947-008	105	99	97	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.

Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P120321-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/21/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
115-07-1	Propene	198	202	102	54-137	
75-71-8	Dichlorodifluoromethane (CFC 12)	196	167	85	55-112	
74-87-3	Chloromethane	190	165	87	66-122	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	200	167	84	64-113	
106-99-0	1,3-Butadiene	204	197	97	74-142	
74-83-9	Bromomethane	194	171	88	72-124	
75-00-3	Chloroethane	196	168	86	69-115	
64-17-5	Ethanol	928	860	93	67-127	
75-05-8	Acetonitrile	194	190	98	63-126	
107-02-8	Acrolein	198	175	88	62-127	
67-64-1	Acetone	1,010	880	87	67-106	
75-69-4	Trichlorofluoromethane	202	173	86	66-121	
67-63-0	2-Propanol (Isopropyl Alcohol)	382	308	81	56-112	
107-13-1	Acrylonitrile	198	194	98	78-128	
75-35-4	1,1-Dichloroethene	212	190	90	74-116	
75-09-2	Methylene Chloride	206	177	86	69-103	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	208	211	101	76-142	
76-13-1	Trichlorotrifluoroethane	206	184	89	69-118	
75-15-0	Carbon Disulfide	208	180	87	71-112	
156-60-5	trans-1,2-Dichloroethene	196	182	93	73-121	
75-34-3	1,1-Dichloroethane	200	187	94	71-118	
1634-04-4	Methyl tert-Butyl Ether	198	188	95	72-115	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P120321-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/21/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
108-05-4	Vinyl Acetate	952	937	98	51-150	
78-93-3	2-Butanone (MEK)	206	196	95	80-130	
156-59-2	cis-1,2-Dichloroethene	206	190	92	73-117	
141-78-6	Ethyl Acetate	398	375	94	79-126	
110-54-3	n-Hexane	198	175	88	68-109	
67-66-3	Chloroform	214	186	87	67-118	
109-99-9	Tetrahydrofuran (THF)	202	182	90	57-130	
71-55-6	1,1,1-Trichloroethane	198	178	90	70-116	
56-23-5	Carbon Tetrachloride	202	187	93	68-123	
110-82-7	Cyclohexane	390	341	87	73-111	
78-87-5	1,2-Dichloropropane	198	175	88	74-114	
75-27-4	Bromodichloromethane	198	182	92	75-120	
123-91-1	1,4-Dioxane	200	180	90	74-120	
80-62-6	Methyl Methacrylate	400	373	93	80-124	
142-82-5	n-Heptane	196	178	91	75-114	
10061-01-5	cis-1,3-Dichloropropene	188	177	94	79-120	
108-10-1	4-Methyl-2-pentanone	204	196	96	79-128	
10061-02-6	trans-1,3-Dichloropropene	210	204	97	83-131	
108-88-3	Toluene	202	170	84	64-115	
591-78-6	2-Hexanone	222	203	91	73-120	
124-48-1	Dibromochloromethane	206	187	91	72-137	
106-93-4	1,2-Dibromoethane	200	171	86	70-126	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P120321-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
123-86-4	n-Butyl Acetate	220	201	91	65-127	
111-65-9	n-Octane	200	184	92	72-118	
108-90-7	Chlorobenzene	202	167	83	65-117	
179601-23-1	m,p-Xylenes	392	342	87	68-119	
75-25-2	Bromoform	208	180	87	79-150	
100-42-5	Styrene	200	178	89	74-127	
95-47-6	o-Xylene	194	167	86	68-118	
111-84-2	n-Nonane	196	175	89	72-116	
98-82-8	Cumene	190	163	86	68-119	
80-56-8	alpha-Pinene	186	165	89	70-123	
103-65-1	n-Propylbenzene	192	163	85	69-119	
622-96-8	4-Ethyltoluene	198	167	84	68-121	
108-67-8	1,3,5-Trimethylbenzene	200	172	86	67-118	
95-63-6	1,2,4-Trimethylbenzene	194	172	89	66-122	
100-44-7	Benzyl Chloride	200	193	97	73-144	
541-73-1	1,3-Dichlorobenzene	200	157	79	64-122	
106-46-7	1,4-Dichlorobenzene	206	153	74	65-125	
95-50-1	1,2-Dichlorobenzene	198	156	79	63-128	
5989-27-5	d-Limonene	200	189	95	72-126	
96-12-8	1,2-Dibromo-3-chloropropane	196	173	88	72-139	
120-82-1	1,2,4-Trichlorobenzene	196	158	81	65-139	
87-68-3	Hexachlorobutadiene	202	161	80	58-137	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P120322-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/22/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
115-07-1	Propene	198	197	99	54-137	
75-71-8	Dichlorodifluoromethane (CFC 12)	196	159	81	55-112	
74-87-3	Chloromethane	190	159	84	66-122	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	200	159	80	64-113	
106-99-0	1,3-Butadiene	204	184	90	74-142	
74-83-9	Bromomethane	194	162	84	72-124	
75-00-3	Chloroethane	196	160	82	69-115	
64-17-5	Ethanol	928	797	86	67-127	
75-05-8	Acetonitrile	194	183	94	63-126	
107-02-8	Acrolein	198	169	85	62-127	
67-64-1	Acetone	1,010	851	84	67-106	
75-69-4	Trichlorofluoromethane	202	165	82	66-121	
67-63-0	2-Propanol (Isopropyl Alcohol)	382	290	76	56-112	
107-13-1	Acrylonitrile	198	190	96	78-128	
75-35-4	1,1-Dichloroethene	212	186	88	74-116	
75-09-2	Methylene Chloride	206	172	83	69-103	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	208	205	99	76-142	
76-13-1	Trichlorotrifluoroethane	206	177	86	69-118	
75-15-0	Carbon Disulfide	208	173	83	71-112	
156-60-5	trans-1,2-Dichloroethene	196	177	90	73-121	
75-34-3	1,1-Dichloroethane	200	182	91	71-118	
1634-04-4	Methyl tert-Butyl Ether	198	181	91	72-115	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result. Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P120322-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/22/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
108-05-4	Vinyl Acetate	952	916	96	51-150	
78-93-3	2-Butanone (MEK)	206	191	93	80-130	
156-59-2	cis-1,2-Dichloroethene	206	185	90	73-117	
141-78-6	Ethyl Acetate	398	369	93	79-126	
110-54-3	n-Hexane	198	167	84	68-109	
67-66-3	Chloroform	214	181	85	67-118	
109-99-9	Tetrahydrofuran (THF)	202	178	88	57-130	
71-55-6	1,1,1-Trichloroethane	198	177	89	70-116	
56-23-5	Carbon Tetrachloride	202	184	91	68-123	
110-82-7	Cyclohexane	390	335	86	73-111	
78-87-5	1,2-Dichloropropane	198	173	87	74-114	
75-27-4	Bromodichloromethane	198	180	91	75-120	
123-91-1	1,4-Dioxane	200	178	89	74-120	
80-62-6	Methyl Methacrylate	400	371	93	80-124	
142-82-5	n-Heptane	196	178	91	75-114	
10061-01-5	cis-1,3-Dichloropropene	188	173	92	79-120	
108-10-1	4-Methyl-2-pentanone	204	194	95	79-128	
10061-02-6	trans-1,3-Dichloropropene	210	200	95	83-131	
108-88-3	Toluene	202	170	84	64-115	
591-78-6	2-Hexanone	222	202	91	73-120	
124-48-1	Dibromochloromethane	206	188	91	72-137	
106-93-4	1,2-Dibromoethane	200	172	86	70-126	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result. Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P120322-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/22/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
123-86-4	n-Butyl Acetate	220	199	90	65-127	
111-65-9	n-Octane	200	184	92	72-118	
108-90-7	Chlorobenzene	202	167	83	65-117	
179601-23-1	m,p-Xylenes	392	342	87	68-119	
75-25-2	Bromoform	208	181	87	79-150	
100-42-5	Styrene	200	178	89	74-127	
95-47-6	o-Xylene	194	167	86	68-118	
111-84-2	n-Nonane	196	174	89	72-116	
98-82-8	Cumene	190	164	86	68-119	
80-56-8	alpha-Pinene	186	166	89	70-123	
103-65-1	n-Propylbenzene	192	164	85	69-119	
622-96-8	4-Ethyltoluene	198	168	85	68-121	
108-67-8	1,3,5-Trimethylbenzene	200	173	87	67-118	
95-63-6	1,2,4-Trimethylbenzene	194	172	89	66-122	
100-44-7	Benzyl Chloride	200	194	97	73-144	
541-73-1	1,3-Dichlorobenzene	200	158	79	64-122	
106-46-7	1,4-Dichlorobenzene	206	155	75	65-125	
95-50-1	1,2-Dichlorobenzene	198	155	78	63-128	
5989-27-5	d-Limonene	200	187	94	72-126	
96-12-8	1,2-Dibromo-3-chloropropane	196	172	88	72-139	
120-82-1	1,2,4-Trichlorobenzene	196	157	80	65-139	
87-68-3	Hexachlorobutadiene	202	160	79	58-137	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY DUPLICATE SUMMARY RESULTS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 02-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-002DUP

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00687

Date Collected: 3/5/12  
Date Received: 3/12/12  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.86

Final Pressure (psig): 3.52

Canister Dilution Factor: 1.85

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
Propene	3.60	2.09	3.43	1.99	3.515	<b>5</b>	25	
Dichlorodifluoromethane (CFC 12)	2.14	0.433	2.09	0.423	2.115	<b>2</b>	25	
Chloromethane	ND	ND	ND	ND	-	-	25	
1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	ND	ND	ND	-	-	25	
1,3-Butadiene	ND	ND	ND	ND	-	-	25	
Bromomethane	ND	ND	ND	ND	-	-	25	
Chloroethane	ND	ND	ND	ND	-	-	25	
Ethanol	579	308	562	298	570.5	<b>3</b>	25	
Acetonitrile	ND	ND	ND	ND	-	-	25	
Acrolein	ND	ND	ND	ND	-	-	25	
Acetone	28.9	12.2	28.7	12.1	28.8	<b>0.7</b>	25	
Trichlorofluoromethane	1.08	0.192	1.12	0.200	1.1	<b>4</b>	25	
2-Propanol (Isopropyl Alcohol)	17.6	7.17	17.7	7.22	17.65	<b>0.6</b>	25	
Acrylonitrile	ND	ND	ND	ND	-	-	25	
1,1-Dichloroethene	ND	ND	ND	ND	-	-	25	
Methylene Chloride	ND	ND	ND	ND	-	-	25	
3-Chloro-1-propene (Allyl Chloride)	ND	ND	ND	ND	-	-	25	
Trichlorotrifluoroethane	ND	ND	ND	ND	-	-	25	
Carbon Disulfide	ND	ND	ND	ND	-	-	25	
trans-1,2-Dichloroethene	ND	ND	ND	ND	-	-	25	
1,1-Dichloroethane	ND	ND	ND	ND	-	-	25	
Methyl tert-Butyl Ether	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.



LABORATORY DUPLICATE SUMMARY RESULTS

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**Client:** Montana DEQ  
**Client Sample ID:** 02-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-002DUP

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00687

Date Collected: 3/5/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/21/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.86

Final Pressure (psig): 3.52

Canister Dilution Factor: 1.85

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
Vinyl Acetate	ND	ND	ND	ND	-	-	25	
2-Butanone (MEK)	ND	ND	ND	ND	-	-	25	
cis-1,2-Dichloroethene	ND	ND	ND	ND	-	-	25	
Ethyl Acetate	37.1	10.3	37.3	10.4	37.2	<b>0.5</b>	25	
n-Hexane	13.5	3.84	13.5	3.83	13.5	<b>0</b>	25	
Chloroform	ND	ND	ND	ND	-	-	25	
Tetrahydrofuran (THF)	ND	ND	ND	ND	-	-	25	
1,1,1-Trichloroethane	ND	ND	ND	ND	-	-	25	
Carbon Tetrachloride	ND	ND	ND	ND	-	-	25	
Cyclohexane	2.36	0.686	2.33	0.677	2.345	<b>1</b>	25	
1,2-Dichloropropane	ND	ND	ND	ND	-	-	25	
Bromodichloromethane	ND	ND	ND	ND	-	-	25	
1,4-Dioxane	ND	ND	ND	ND	-	-	25	
Methyl Methacrylate	ND	ND	ND	ND	-	-	25	
n-Heptane	5.70	1.39	5.81	1.42	5.755	<b>2</b>	25	
cis-1,3-Dichloropropene	ND	ND	ND	ND	-	-	25	
4-Methyl-2-pentanone	ND	ND	ND	ND	-	-	25	
trans-1,3-Dichloropropene	ND	ND	ND	ND	-	-	25	
Toluene	34.4	9.12	34.3	9.10	34.35	<b>0.3</b>	25	
2-Hexanone	ND	ND	ND	ND	-	-	25	
Dibromochloromethane	ND	ND	ND	ND	-	-	25	
1,2-Dibromoethane	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

LABORATORY DUPLICATE SUMMARY RESULTS

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**Client:** Montana DEQ  
**Client Sample ID:** 02-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-002DUP

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00687

Date Collected: 3/5/12  
Date Received: 3/12/12  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.86

Final Pressure (psig): 3.52

Canister Dilution Factor: 1.85

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
n-Butyl Acetate	4.95	1.04	4.87	1.03	4.91	2	25	
n-Octane	1.99	0.425	1.94	0.415	1.965	3	25	
Chlorobenzene	ND	ND	ND	ND	-	-	25	
m,p-Xylenes	22.3	5.13	22.0	5.07	22.15	1	25	
Bromoform	ND	ND	ND	ND	-	-	25	
Styrene	2.31	0.543	2.26	0.532	2.285	2	25	
o-Xylene	7.71	1.78	7.63	1.76	7.67	1	25	
n-Nonane	2.09	0.399	2.14	0.408	2.115	2	25	
Cumene	ND	ND	ND	ND	-	-	25	
alpha-Pinene	11.0	1.97	10.9	1.96	10.95	0.9	25	
n-Propylbenzene	1.65	0.336	1.62	0.330	1.635	2	25	
4-Ethyltoluene	2.55	0.519	2.53	0.515	2.54	0.8	25	
1,3,5-Trimethylbenzene	3.02	0.615	3.00	0.610	3.01	0.7	25	
1,2,4-Trimethylbenzene	8.80	1.79	8.72	1.77	8.76	0.9	25	
Benzyl Chloride	ND	ND	ND	ND	-	-	25	
1,3-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
1,4-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
1,2-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
d-Limonene	24.2	4.34	23.8	4.28	24	2	25	
1,2-Dibromo-3-chloropropane	ND	ND	ND	ND	-	-	25	
1,2,4-Trichlorobenzene	ND	ND	ND	ND	-	-	25	
Hexachlorobutadiene	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 01-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P1200947-001

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01505

Date Collected: 3/5/12  
Date Received: 3/12/12  
Date Analyzed: 3/22/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.41 Final Pressure (psig): 3.52

Canister Dilution Factor: 1.61

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	<b>0.048</b>	0.040	<b>0.019</b>	0.016	
107-06-2	1,2-Dichloroethane	<b>0.23</b>	0.040	<b>0.056</b>	0.0099	
71-43-2	Benzene	<b>4.0</b>	0.12	<b>1.2</b>	0.038	
79-01-6	Trichloroethene	<b>0.29</b>	0.040	<b>0.054</b>	0.0075	
79-00-5	1,1,2-Trichloroethane	ND	0.16	ND	0.030	
127-18-4	Tetrachloroethene	<b>0.17</b>	0.040	<b>0.026</b>	0.0059	
100-41-4	Ethylbenzene	<b>2.3</b>	0.16	<b>0.52</b>	0.037	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.040	ND	0.0059	
91-20-3	Naphthalene	<b>0.36</b>	0.16	<b>0.068</b>	0.031	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 02-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-002

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00687

Date Collected: 3/5/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/22/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.86      Final Pressure (psig): 3.52

Canister Dilution Factor: 1.85

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.046	ND	0.018	
107-06-2	1,2-Dichloroethane	<b>0.69</b>	0.046	<b>0.17</b>	0.011	
71-43-2	Benzene	<b>8.2</b>	0.14	<b>2.6</b>	0.043	
79-01-6	Trichloroethene	<b>0.26</b>	0.046	<b>0.048</b>	0.0086	
79-00-5	1,1,2-Trichloroethane	ND	0.19	ND	0.034	
127-18-4	Tetrachloroethene	<b>0.23</b>	0.046	<b>0.033</b>	0.0068	
100-41-4	Ethylbenzene	<b>5.8</b>	0.19	<b>1.3</b>	0.043	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.046	ND	0.0067	
91-20-3	Naphthalene	<b>1.1</b>	0.19	<b>0.21</b>	0.035	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 03-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-003

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00178

Date Collected: 3/5/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/22/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.80      Final Pressure (psig): 3.51

Canister Dilution Factor: 1.84

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.046	ND	0.018	
107-06-2	1,2-Dichloroethane	<b>0.11</b>	0.046	<b>0.027</b>	0.011	
71-43-2	Benzene	ND	0.14	ND	0.043	
79-01-6	Trichloroethene	ND	0.046	ND	0.0086	
79-00-5	1,1,2-Trichloroethane	ND	0.18	ND	0.034	
127-18-4	Tetrachloroethene	<b>3.9</b>	0.046	<b>0.57</b>	0.0068	
100-41-4	Ethylbenzene	ND	0.18	ND	0.042	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.046	ND	0.0067	
91-20-3	Naphthalene	ND	0.18	ND	0.035	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 05-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-004

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Wida Ang  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01198

Date Collected: 3/7/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/23/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.39      Final Pressure (psig): 3.53

Canister Dilution Factor: 1.61

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.040	ND	0.016	
107-06-2	1,2-Dichloroethane	1.2	0.040	0.31	0.0099	
71-43-2	Benzene	20	0.12	6.2	0.038	
79-01-6	Trichloroethene	0.24	0.040	0.044	0.0075	
79-00-5	1,1,2-Trichloroethane	ND	0.16	ND	0.030	
127-18-4	Tetrachloroethene	0.060	0.040	0.0088	0.0059	
100-41-4	Ethylbenzene	11	0.16	2.6	0.037	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.040	ND	0.0059	
91-20-3	Naphthalene	1.5	0.16	0.28	0.031	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 06-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-005

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Wida Ang  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00137

Date Collected: 3/7/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/23/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.12      Final Pressure (psig): 3.74

Canister Dilution Factor: 1.47

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.037	ND	0.014	
107-06-2	1,2-Dichloroethane	<b>0.57</b>	0.037	<b>0.14</b>	0.0091	
71-43-2	Benzene	<b>4.6</b>	0.11	<b>1.4</b>	0.035	
79-01-6	Trichloroethene	<b>0.27</b>	0.037	<b>0.051</b>	0.0068	
79-00-5	1,1,2-Trichloroethane	ND	0.15	ND	0.027	
127-18-4	Tetrachloroethene	<b>0.19</b>	0.037	<b>0.028</b>	0.0054	
100-41-4	Ethylbenzene	<b>2.6</b>	0.15	<b>0.59</b>	0.034	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.037	ND	0.0054	
91-20-3	Naphthalene	<b>0.65</b>	0.15	<b>0.12</b>	0.028	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 07-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-006

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Wida Ang  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01570

Date Collected: 3/7/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/23/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.36      Final Pressure (psig): 3.72

Canister Dilution Factor: 1.49

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.037	ND	0.015	
107-06-2	1,2-Dichloroethane	<b>0.38</b>	0.037	<b>0.095</b>	0.0092	
71-43-2	Benzene	<b>0.50</b>	0.11	<b>0.16</b>	0.035	
79-01-6	Trichloroethene	<b>0.22</b>	0.037	<b>0.041</b>	0.0069	
79-00-5	1,1,2-Trichloroethane	ND	0.15	ND	0.027	
127-18-4	Tetrachloroethene	<b>0.11</b>	0.037	<b>0.017</b>	0.0055	
100-41-4	Ethylbenzene	<b>0.68</b>	0.15	<b>0.16</b>	0.034	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.037	ND	0.0054	
91-20-3	Naphthalene	<b>0.37</b>	0.15	<b>0.070</b>	0.028	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 08-IF-030712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-007

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Wida Ang  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00184

Date Collected: 3/7/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/23/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.87      Final Pressure (psig): 3.58

Canister Dilution Factor: 1.55

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.039	ND	0.015	
107-06-2	1,2-Dichloroethane	<b>0.48</b>	0.039	<b>0.12</b>	0.0096	
71-43-2	Benzene	<b>0.52</b>	0.12	<b>0.16</b>	0.036	
79-01-6	Trichloroethene	<b>0.96</b>	0.039	<b>0.18</b>	0.0072	
79-00-5	1,1,2-Trichloroethane	ND	0.16	ND	0.028	
127-18-4	Tetrachloroethene	<b>0.14</b>	0.039	<b>0.021</b>	0.0057	
100-41-4	Ethylbenzene	<b>0.65</b>	0.16	<b>0.15</b>	0.036	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.039	ND	0.0056	
91-20-3	Naphthalene	ND	0.16	ND	0.030	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 09-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-008

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Wida Ang  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00553

Date Collected: 3/5/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/23/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.65      Final Pressure (psig): 3.56

Canister Dilution Factor: 1.82

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.046	ND	0.018	
107-06-2	1,2-Dichloroethane	<b>0.47</b>	0.046	<b>0.12</b>	0.011	
71-43-2	Benzene	<b>0.64</b>	0.14	<b>0.20</b>	0.043	
79-01-6	Trichloroethene	<b>5.2</b>	0.046	<b>0.96</b>	0.0085	
79-00-5	1,1,2-Trichloroethane	ND	0.18	ND	0.033	
127-18-4	Tetrachloroethene	<b>0.097</b>	0.046	<b>0.014</b>	0.0067	
100-41-4	Ethylbenzene	<b>1.6</b>	0.18	<b>0.37</b>	0.042	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.046	ND	0.0066	
91-20-3	Naphthalene	ND	0.18	ND	0.035	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P120322-MB

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/22/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result		MRL		Data Qualifier
		$\mu\text{g}/\text{m}^3$	$\mu\text{g}/\text{m}^3$	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.025	ND	0.0098	
107-06-2	1,2-Dichloroethane	ND	0.025	ND	0.0062	
71-43-2	Benzene	ND	0.075	ND	0.023	
79-01-6	Trichloroethene	ND	0.025	ND	0.0047	
79-00-5	1,1,2-Trichloroethane	ND	0.10	ND	0.018	
127-18-4	Tetrachloroethene	ND	0.025	ND	0.0037	
100-41-4	Ethylbenzene	ND	0.10	ND	0.023	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.025	ND	0.0036	
91-20-3	Naphthalene	ND	0.10	ND	0.019	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P120323-MB

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/23/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result		MRL		Data Qualifier
		$\mu\text{g}/\text{m}^3$	$\mu\text{g}/\text{m}^3$	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.025	ND	0.0098	
107-06-2	1,2-Dichloroethane	ND	0.025	ND	0.0062	
71-43-2	Benzene	ND	0.075	ND	0.023	
79-01-6	Trichloroethene	ND	0.025	ND	0.0047	
79-00-5	1,1,2-Trichloroethane	ND	0.10	ND	0.018	
127-18-4	Tetrachloroethene	ND	0.025	ND	0.0037	
100-41-4	Ethylbenzene	ND	0.10	ND	0.023	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.025	ND	0.0036	
91-20-3	Naphthalene	ND	0.10	ND	0.019	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister(s)  
 Test Notes:

Date(s) Collected: 3/5 - 3/7/12  
 Date(s) Received: 3/12/12  
 Date(s) Analyzed: 3/22 - 3/23/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		% Recovered	% Recovered	% Recovered		
Method Blank	P120322-MB	102	116	110	70-130	
Method Blank	P120323-MB	100	112	114	70-130	
Lab Control Sample	P120322-LCS	103	115	114	70-130	
Lab Control Sample	P120323-LCS	98	112	110	70-130	
01-IF-030512	P1200947-001	98	115	119	70-130	
02-IF-030512	P1200947-002	96	114	115	70-130	
03-IF-030512	P1200947-003	98	114	120	70-130	
05-IF-030712	P1200947-004	100	113	99	70-130	
06-IF-030712	P1200947-005	100	112	109	70-130	
07-IF-030712	P1200947-006	98	114	114	70-130	
08-IF-030712	P1200947-007	97	113	115	70-130	
09-IF-030512	P1200947-008	97	113	114	70-130	
09-IF-030512	P1200947-008DUP	98	111	113	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.

Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P120322-LCS

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/22/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
75-01-4	Vinyl Chloride	3.88	3.45	89	56-127	
107-06-2	1,2-Dichloroethane	4.00	3.48	87	51-140	
71-43-2	Benzene	3.96	3.89	98	56-125	
79-01-6	Trichloroethene	3.88	3.29	85	51-127	
79-00-5	1,1,2-Trichloroethane	3.92	3.39	86	49-137	
127-18-4	Tetrachloroethene	3.68	3.22	88	58-134	
100-41-4	Ethylbenzene	3.96	3.23	82	56-126	
79-34-5	1,1,2,2-Tetrachloroethane	3.84	3.33	87	53-148	
91-20-3	Naphthalene	3.44	3.35	97	25-149	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
CAS Sample ID: P120323-LCS

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/23/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
75-01-4	Vinyl Chloride	3.88	3.38	87	56-127	
107-06-2	1,2-Dichloroethane	4.00	3.33	83	51-140	
71-43-2	Benzene	3.96	3.93	99	56-125	
79-01-6	Trichloroethene	3.88	3.22	83	51-127	
79-00-5	1,1,2-Trichloroethane	3.92	3.26	83	49-137	
127-18-4	Tetrachloroethene	3.68	3.18	86	58-134	
100-41-4	Ethylbenzene	3.96	3.39	86	56-126	
79-34-5	1,1,2,2-Tetrachloroethane	3.84	3.33	87	53-148	
91-20-3	Naphthalene	3.44	3.47	101	25-149	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY DUPLICATE SUMMARY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 09-IF-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1200947  
 CAS Sample ID: P1200947-008DUP

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Wida Ang  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00553

Date Collected: 3/5/12  
 Date Received: 3/12/12  
 Date Analyzed: 3/23/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.65      Final Pressure (psig): 3.56

Canister Dilution Factor: 1.82

CAS #	Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
		µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
75-01-4	Vinyl Chloride	ND	ND	ND	ND	-	-	25	
107-06-2	1,2-Dichloroethane	0.466	0.115	0.469	0.116	0.4675	<b>0.6</b>	25	
71-43-2	Benzene	0.638	0.200	0.631	0.198	0.6345	<b>1</b>	25	
79-01-6	Trichloroethene	5.18	0.964	5.09	0.947	5.135	<b>2</b>	25	
79-00-5	1,1,2-Trichloroethane	ND	ND	ND	ND	-	-	25	
127-18-4	Tetrachloroethene	0.0969	0.0143	0.0953	0.0141	0.0961	<b>2</b>	25	
100-41-4	Ethylbenzene	1.63	0.375	1.62	0.374	1.625	<b>0.6</b>	25	
79-34-5	1,1,2,2-Tetrachloroethane	ND	ND	ND	ND	-	-	25	
91-20-3	Naphthalene	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.



Massachusetts APH

Hydrocarbon Ranges

ICAL Method: M8011412A.M

ICAL Date: 1/14/12

Instrument ID: MS08

	areas				masses							
	0.5	1	5	25	50	100	0.5	1	5	25	50	100
<b>C5-C8 Aliphatics</b>												
Isopentane	60478	112606	488243	2480105	4710733	9222808	0.520	1.04	5.20	26.0	52.0	104
n-Hexane	59953	110631	507126	2577865	4931102	9678199	0.490	0.98	4.90	24.5	49.0	98
Cyclohexane	63203	115230	548910	2874726	5524402	10890911	0.500	1.00	5.00	25.0	50.0	100
2,3-Dimethylpentane	68648	127222	585183	2998017	5742104	11319597	0.520	1.04	5.20	26.0	52.0	104
n-Heptane	66143	120286	558074	2884817	5509801	10873046	0.490	0.98	4.90	24.5	49.0	98
n-Octane	76169	137829	632891	3286017	6205035	12092514	0.505	1.01	5.05	25.3	50.5	101
<b>area sum:</b>	<b>394594</b>	<b>723804</b>	<b>3320427</b>	<b>17101547</b>	<b>32623177</b>	<b>64077075</b>	<b>3.025</b>	<b>6.050</b>	<b>30.25</b>	<b>151.30</b>	<b>302.50</b>	<b>605.00</b>

	areas				masses							
	0.5	1	5	25	50	100	0.5	1	5	25	50	100
<b>C9-C12 Aliphatics</b>												
2,3-Dimethylheptane	78780	147724	682518	3501487	6691428	13148840	0.485	0.97	4.85	24.3	48.5	97
n-Nonane	77464	140680	662221	3417128	6555165	12940625	0.485	0.97	4.85	24.3	48.5	97
n-Decane	80593	147724	705259	3651028	7021424	13870530	0.490	0.98	4.90	24.5	49.0	98
2-Butylcyclohexane	88915	166782	776903	4026931	7749936	15415340	0.495	0.99	4.95	24.8	49.5	99
n-Undecane	80867	152557	719618	3772331	7263077	14385881	0.480	0.96	4.80	24.0	48.0	96
n-Dodecane	70045	143746	687255	3673524	7115012	14041583	0.500	1.00	5.00	25.0	50.0	100
<b>area sum:</b>	<b>476664</b>	<b>899213</b>	<b>4233774</b>	<b>22042429</b>	<b>42396042</b>	<b>83802799</b>	<b>2.935</b>	<b>5.870</b>	<b>29.35</b>	<b>146.90</b>	<b>293.50</b>	<b>587.00</b>

	areas				masses							
	0.5	1	5	25	50	100	0.5	1	5	25	50	100
<b>C9-C10 Aromatics</b>												
Isopropylbenzene	10658	19169	87439	448287	849124	1695737	0.485	0.97	4.85	24.3	48.5	97
3-Ethyltoluene	11805	20564	97823	504510	977020	1973178	0.470	0.94	4.70	23.5	47.0	94
1,3,5-Trimethylbenzene	15759	29260	134073	700133	1353305	2726630	0.490	0.98	4.90	24.5	49.0	98
1,2,3-Trimethylbenzene	17195	31328	143473	757946	1493373	3063059	0.495	0.99	4.95	24.8	49.5	99
p-Isopropyltoluene	9490	18116	83401	440962	872799	1814314	0.460	0.92	4.60	23.0	46.0	92
<b>area sum:</b>	<b>64907</b>	<b>118437</b>	<b>546209</b>	<b>2851838</b>	<b>5545621</b>	<b>11272918</b>	<b>2.400</b>	<b>4.800</b>	<b>24.00</b>	<b>120.10</b>	<b>240.00</b>	<b>480.00</b>

Massachusetts APH  
Hydrocarbon Ranges

ICAL: M8011412A.M

ICAL Date: 1/14/12

Instrument ID: MS08

areas

<u>Internal Standards (TIC)</u>	0.5	1	5	25	50	100
Bromochloromethane (IS1)	1312793	1307309	1293586	1289283	1284194	1283817
1,4-Difluorobenzene (IS2)	2547079	2541159	2512498	2475255	2422119	2352159
Chlorobenzene-d5 (IS3)	2929189	2913374	2880258	2895080	2878314	2908396

Internal Standards (EIC)

Bromochloromethane (IS1)	249696	247899	245985	245585	244327	243561
1,4-Difluorobenzene (IS2)	1153657	1149947	1133233	1136223	1127203	1138730
Chlorobenzene-d5 (IS3)	1007702	1004638	991212	992925	982812	986950

Surrogates (TIC)

	0.5	1	5	25	50	100
1,2-Dichloroethane-d4	1162497	1154617	1105630	1126201	1103958	1095945
m-Toluene-d8	3508018	3490493	3455382	3466688	3437986	3457323
o-p-Bromofluorobenzene	2298220	2284797	2197239	2227396	2192159	2197999

C5-C8 Aliphatics

	<u>RRFs</u>					
	0.5	1	5	25	50	100
	2.8268	2.6009	2.4215	2.4870	2.3919	2.3252
						<u>RRF<sub>avg</sub></u> 2.509
						<u>%RSD</u> 7.24

C9-C12 Aliphatics

	<u>RRFs</u>					
	0.5	1	5	25	50	100
	4.0291	3.8120	3.6383	3.7780	3.6744	3.6163
						<u>RRF<sub>avg</sub></u> 3.758
						<u>%RSD</u> 4.09

C9-C10 Aromatics

	<u>RRFs</u>					
	0.5	1	5	25	50	100
	0.6709	0.6140	0.5740	0.5979	0.5878	0.5949
						<u>RRF<sub>avg</sub></u> 0.607
						<u>%RSD</u> 5.63

**Massachusetts APH**  
**Continuing Calibration Verification Check Sheet**

Data File Name: 03211203.D  
 Data File Path: J:\MS08\Data\2012\_03\21\  
 Operator: EM  
 Date Acquired: 3/21/12 5:39  
 Acq. Method File: TO15.M  
 Sample Name: 25ng TO-15/MAPH CCV STD  
 Misc Info: S25-03151201/S25-03161202  
 Instrument Name: MS08

*Enter RRFs from current ICAL!*

Internal Standards	RT	Area						
7) 1,4-Difluorobenzene (IS2)	15.89	1102904						
16) Chlorobenzene-d5 (IS3)	21.69	947969						
<b>C5-C8 Aliphatics</b>								
	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
3) Isopentane	8.19	2464516	2.569	154.9	2.39	-30	30	Pass
4) n-Hexane	13.08	2578864						
9) Cyclohexane	15.81	2862780						
10) 2,3-Dimethylpentane	16.17	3020819	Spike	ICAL				
11) n-Heptane	17.35	2903399	Amt (ng)	RRF				
14) n-Octane	20.69	3316542	151.30	2.509				
		<b>17146920</b>						
<b>C9-C12 Aliphatics</b>								
	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
18) 2,3-Dimethylheptane	22.31	3515879	3.955	154.6	5.23	-30	30	Pass
19) n-Nonane	23.30	3434658						
25) n-Decane	25.26	3663558						
28) Butylcyclohexane	25.99	4002852	Spike	ICAL				
29) n-Undecane	26.76	3775128	Amt (ng)	RRF				
30) n-Dodecane	28.00	3635722	146.90	3.758				
		<b>22027797</b>						
<b>C9-C10 Aromatics</b>								
	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
22) Isopropylbenzene	23.78	446789	0.612	121.1	0.80	-30	30	Pass
23) 1-Methyl-3-ethylbenzene	24.52	492493						
24) 1,3,5-Trimethylbenzene	24.66	681321						
26) p-Isopropyltoluene	25.68	426000	Spike	ICAL				
27) 1,2,3-Trimethylbenzene	25.68	739855	Amt (ng)	RRF				
		<b>2786458</b>	120.1	0.607				

*EM 3/21/12*

Response Factor Report MS08

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Mon Dec 12 09:45:45 2011  
 Response Via : Initial Calibration

Calibration Files

0.1 =12091118.D 0.2 =12091119.D 0.5 =12091120.D 1.0 =12091121.D 5.0 =12091122.D 25 =12091123.D  
 50 =12091124.D 100 =12091125.D

Compound	0.1	0.2	0.5	1.0	5.0	25	50	100	Avg	%RSD
1) IR Bromochloromethane...										
2) T Propene	2.487	1.965	1.952	1.692	2.194	2.303	2.185	2.036	2.102	11.62
3) T Dichlorodifluo...	3.545	2.805	2.940	2.748	2.432	2.361	2.278	2.199	2.664	16.73
4) T Chloromethane	3.290	2.811	2.875	2.512	2.176	2.286	1.977	1.664	2.449	21.62
5) T 1,2-Dichloro-1...	1.901	1.661	1.575	1.458	1.359	1.333	1.200	1.190	1.460	16.68
6) T Vinyl Chloride	2.906	2.590	2.631	2.374	2.145	2.129	1.950	1.939	2.333	15.07
7) T 1,3-Butadiene	2.145	1.788	1.790	1.718	1.581	1.689	1.621	1.616	1.743	10.33
8) T Bromomethane	1.687	1.328	1.349	1.212	1.094	1.142	1.111	1.092	1.252	16.21
9) T Chloroethane	1.671	1.259	1.333	1.193	1.101	1.091	1.046	1.044	1.217	17.29
10) T Ethanol	2.040	1.345	1.186	1.085	1.022	1.096	1.076	1.064	1.239	27.34
11) T Acetonitrile	4.365	2.973	2.723	2.546	2.480	2.562	2.482	2.468	2.825	22.85
12) T Acrolein	1.328	0.931	0.907	0.861	0.772	0.814	0.789	0.779	0.897	20.48
13) T Acetone	1.645	1.316	1.249	1.129	0.994	1.015	0.981	0.964	1.161	20.26
14) T Trichlorofluor...	2.776	2.385	2.496	2.272	2.097	2.135	2.042	2.011	2.277	11.59
15) T 2-Propanol (Is...		4.463	4.536	4.168	2.920	2.763	2.577	2.672	3.443	26.09
16) T Acrylonitrile	2.701	1.907	1.878	1.783	1.745	1.813	1.765	1.738	1.916	16.85
17) T 1,1-Dichloroet...	1.458	1.418	1.452	1.263	1.206	1.223	1.172	1.152	1.293	9.95
18) T 2-Methyl-2-Pro...	4.464	3.710	3.939	3.616	3.558	3.584			3.812	9.14
19) T Methylene Chlo...			1.958	1.578	1.294	1.273	1.220	1.201	1.421	20.88
20) T 3-Chloro-1-pro...	2.347	2.260	2.345	2.220	2.150	2.296	2.230	2.191	2.255	3.15
21) T Trichlorotrifl...	1.348	1.244	1.299	1.193	1.089	1.098	1.062	1.035	1.171	10.01
22) T Carbon Disulfide		6.210	5.678	5.057	4.654	4.699	4.488	4.424	5.030	13.40
23) T trans-1,2-Dich...	2.436	2.237	2.206	2.142	1.951	1.953	1.883	1.835	2.080	10.00
24) T 1,1-Dichloroet...	2.896	2.757	2.747	2.500	2.307	2.379	2.288	2.236	2.514	10.07
25) T Methyl tert-Bu...	5.128	4.611	4.650	4.307	4.119	4.185	4.042	3.984	4.378	8.93
26) T Vinyl Acetate	0.430	0.368	0.423	0.399	0.393	0.410	0.402	0.398	0.403	4.70
27) T 2-Butanone (MEK)	1.265	1.003	1.038	0.972	0.927	0.964	0.928	0.794	0.986	13.56
28) T cis-1,2-Dichlo...	2.442	2.056	2.101	1.938	1.826	1.836	1.764	1.720	1.960	12.05
29) T Diisopropyl Ether	1.446	1.223	1.323	1.262	1.171	1.188	1.153	1.148	1.239	8.26
30) T Ethyl Acetate	0.611	0.554	0.573	0.545	0.517	0.535	0.520	0.516	0.546	6.02
31) T n-Hexane	3.541	3.006	3.181	2.805	2.582	2.633	2.545	2.543	2.855	12.70

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Response Factor Report MS08

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M

Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

32) T	Chloroform	2.679	2.408	2.409	2.204	1.983	2.054	1.974	1.937	2.206	12.18
33) S	1,2-Dichloroet...	1.586	1.577	1.582	1.590	1.535	1.533	1.520	1.504	1.553	2.19
34) T	Tetrahydrofura...	1.274	0.932	1.100	0.938	0.910	0.879	0.830	0.821	0.961	15.96
35) T	Ethyl tert-But...	1.999	1.832	2.033	1.818	1.732	1.757	1.708	1.687	1.821	7.18
36) T	1,2-Dichloroet...	2.099	1.992	1.976	1.806	1.683	1.700	1.599	1.531	1.798	11.39
37) IR	1,4-Difluorobenzen...	-----ISTD-----									
38) T	1,1,1-Trichlor...	0.531	0.465	0.486	0.453	0.419	0.434	0.413	0.403	0.451	9.46
39) T	Isopropyl Acetate	0.227	0.205	0.220	0.206	0.194	0.201	0.192	0.188	0.204	6.71
40) T	1-Butanol		0.348	0.319	0.300	0.320	0.335	0.324	0.318	0.323	4.66
41) T	Benzene	1.633	1.385	1.359	1.227	1.136	1.158	1.101	1.070	1.258	15.11
42) T	Carbon Tetrach...	0.394	0.356	0.369	0.340	0.334	0.352	0.337	0.333	0.352	6.03
43) T	Cyclohexane	0.666	0.576	0.588	0.539	0.492	0.500	0.479	0.469	0.539	12.59
44) T	tert-Amyl Meth...	1.092	0.926	0.998	0.937	0.900	0.918	0.881	0.858	0.939	7.95
45) T	1,2-Dichloropr...	0.397	0.361	0.368	0.324	0.312	0.316	0.300	0.293	0.334	11.12
46) T	Bromodichlorom...	0.458	0.387	0.392	0.378	0.362	0.378	0.362	0.353	0.384	8.57
47) T	Trichloroethene	0.466	0.375	0.379	0.344	0.323	0.328	0.316	0.309	0.355	14.62
48) T	1,4-Dioxane	0.305	0.282	0.288	0.248	0.245	0.246	0.237	0.230	0.260	10.55
49) T	2,2,4-Trimethy...	1.964	1.629	1.686	1.546	1.447	1.461	1.387	1.337	1.557	12.98
50) T	Methyl Methacr...	0.156	0.139	0.150	0.135	0.136	0.142	0.136	0.135	0.141	5.53
51) T	n-Heptane	0.427	0.373	0.385	0.364	0.341	0.344	0.332	0.324	0.361	9.35
52) T	cis-1,3-Dichlo...	0.571	0.502	0.521	0.491	0.482	0.500	0.481	0.471	0.502	6.33
53) T	4-Methyl-2-pen...	0.359	0.295	0.308	0.303	0.297	0.303	0.289	0.284	0.305	7.68
54) T	trans-1,3-Dich...	0.476	0.413	0.443	0.421	0.429	0.453	0.437	0.427	0.437	4.59
55) T	1,1,2-Trichlor...	0.340	0.301	0.310	0.284	0.276	0.285	0.271	0.265	0.292	8.44
56) IR	Chlorobenzene-d5 (...	-----ISTD-----									
57) S	Toluene-d8 (SS2)	2.305	2.306	2.305	2.309	2.306	2.300	2.310	2.304	2.306	0.14
58) T	Toluene	3.623	2.920	3.025	2.738	2.571	2.574	2.473	2.421	2.793	14.19
59) T	2-Hexanone	2.049	1.429	1.515	1.403	1.395	1.422	1.367	1.337	1.490	15.57
60) T	Dibromochlorom...	0.696	0.618	0.636	0.636	0.612	0.648	0.629	0.625	0.638	4.14
61) T	1,2-Dibromoethane	0.831	0.730	0.718	0.658	0.641	0.662	0.638	0.628	0.688	9.97
62) T	n-Butyl Acetate	2.338	1.850	1.857	1.721	1.704	1.765	1.737	1.737	1.839	11.40
63) T	n-Octane	0.785	0.652	0.714	0.643	0.602	0.611	0.580	0.572	0.645	11.28
64) T	Tetrachloroethene	1.090	0.904	0.904	0.826	0.767	0.779	0.756	0.755	0.848	13.63
65) T	Chlorobenzene	2.375	1.868	1.886	1.744	1.610	1.637	1.574	1.548	1.780	15.30
66) T	Ethylbenzene	3.846	3.171	3.188	2.956	2.803	2.848	2.743	2.680	3.030	12.51
67) T	m- & p-Xylenes	3.081	2.499	2.611	2.392	2.216	2.273	2.170	2.112	2.419	13.08
68) T	Bromoform	0.622	0.494	0.556	0.541	0.551	0.598	0.589	0.594	0.568	7.19
69) T	Styrene	2.478	1.895	1.937	1.800	1.735	1.803	1.726	1.675	1.881	13.64
70) T	o-Xylene	3.279	2.708	2.726	2.479	2.331	2.373	2.283	2.224	2.550	13.65

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

71)	T	n-Nonane	2.005	1.753	1.847	1.675	1.581	1.597	1.518	1.470	1.681	10.69
72)	T	1,1,2,2-Tetrac...	1.348	1.164	1.262	1.150	1.083	1.121	1.085	1.066	1.160	8.48
73)	S	Bromofluoroben...	0.774	0.771	0.779	0.781	0.775	0.778	0.782	0.780	0.777	0.50
74)	T	Cumene	4.522	3.495	3.524	3.169	2.962	3.002	2.884	2.831	3.299	16.98
75)	T	alpha-Pinene	1.782	1.636	1.734	1.601	1.537	1.578	1.518	1.498	1.611	6.35
76)	T	n-Propylbenzene	5.441	4.229	4.306	3.924	3.710	3.754	3.553	3.394	4.039	16.01
77)	T	3-Ethyltoluene	4.091	3.373	3.461	3.038	2.964	2.981	2.893	2.845	3.206	13.14
78)	T	4-Ethyltoluene	4.003	3.202	3.301	3.151	2.837	2.983	2.810	2.776	3.133	12.84
79)	T	1,3,5-Trimethy...	3.324	2.662	2.723	2.537	2.378	2.426	2.344	2.403	2.600	12.43
80)	T	alpha-Methylst...	1.742	1.386	1.449	1.371	1.390	1.402	1.354	1.399	1.436	8.80
81)	T	2-Ethyltoluene	4.143	3.362	3.430	3.169	2.975	3.048	2.919	2.985	3.254	12.43
82)	T	1,2,4-Trimethy...	3.356	2.663	2.791	2.551	2.433	2.513	2.454	2.548	2.664	11.37
83)	T	n-Decane	1.969	1.702	1.844	1.705	1.582	1.626	1.554	1.609	1.699	8.39
84)	T	Benzyl Chloride	2.373	1.648	1.772	1.688	1.799	2.058	2.054	2.190	1.948	13.35
85)	T	1,3-Dichlorobe...	2.498	1.658	1.649	1.455	1.377	1.409	1.374	1.436	1.607	23.49
86)	T	1,4-Dichlorobe...	2.655	1.785	1.618	1.492	1.380	1.428	1.388	1.449	1.649	25.99
87)	T	sec-Butylbenzene	4.527	3.773	3.816	3.471	3.264	3.328	3.214	3.176	3.571	12.80
88)	T	4-Isopropyltol...	4.426	3.674	3.683	3.416	3.254	3.356	3.253	3.366	3.554	10.98
89)	T	1,2,3-Trimethy...	3.266	2.650	2.852	2.588	2.471	2.573	2.506	2.619	2.691	9.63
90)	T	1,2-Dichlorobe...	2.173	1.589	1.545	1.435	1.332	1.387	1.370	1.415	1.531	17.89
91)	T	d-Limonene	1.124	0.998	1.106	1.069	1.064	1.072	1.040	1.065	1.067	3.58
92)	T	1,2-Dibromo-3-...	0.563	0.488	0.501	0.486	0.494	0.525	0.532	0.542	0.517	5.50
93)	T	n-Undecane	1.922	1.637	1.755	1.671	1.608	1.630	1.608	1.592	1.678	6.63
94)	T	1,2,4-Trichlor...		1.335	1.165	1.062	1.017	1.071	1.086	1.118	1.122	9.32
95)	T	Naphthalene		4.848	3.923	3.638	3.497	3.721	3.755	3.783	3.881	11.50
96)	T	n-Dodecane	1.866	1.599	1.669	1.586	1.545	1.575	1.553	1.509	1.613	6.99
97)	T	Hexachlorobuta...	0.841	0.749	0.730	0.673	0.635	0.664	0.663	0.676	0.704	9.50
98)	T	Cyclohexanone	1.172	0.979	1.021	0.969	0.952	0.962	0.931	0.903	0.986	8.37
99)	T	tert-Butylbenzene	3.486	2.799	2.880	2.658	2.449	2.533	2.453	2.528	2.723	12.71
100)	T	n-Butylbenzene	3.596	2.904	2.958	2.670	2.572	2.625	2.613	2.593	2.817	12.33

(#) = Out of Range

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\21\  
 Data File : 03211201.D  
 Acq On : 21 Mar 2012 4:13  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 21 07:38:45 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 IR	Bromochloromethane (IS1)	1.000	1.000	0.0	83	-0.02
2 T	Propene	2.102	2.230	-6.1	80	0.00
3 T	Dichlorodifluoromethane (CF	2.664	2.371	11.0	83	0.00
4 T	Chloromethane	2.449	2.229	9.0	81	0.00
5 T	1,2-Dichloro-1,1,2,2-tetra	1.460	1.253	14.2	78	0.00
6 T	Vinyl Chloride	2.333	2.088	10.5	82	0.00
7 T	1,3-Butadiene	1.743	1.698	2.6	84	-0.01
8 T	Bromomethane	1.252	1.103	11.9	80	0.00
9 T	Chloroethane	1.217	1.086	10.8	83	0.00
10 T	Ethanol	1.239	1.182	4.6	90	-0.07
11 T	Acetonitrile	2.825	2.652	6.1	86	-0.05
12 T	Acrolein	0.897	0.782	12.8	80	-0.02
13 T	Acetone	1.161	1.046	9.9	86	-0.05
14 T	Trichlorofluoromethane	2.277	2.123	6.8	83	0.00
15 T	2-Propanol (Isopropanol)	3.443	2.957	14.1	89	-0.05
16 T	Acrylonitrile	1.916	1.850	3.4	85	-0.03
17 T	1,1-Dichloroethene	1.293	1.194	7.7	81	0.00
18 T	2-Methyl-2-Propanol (tert-B	3.812	3.371	11.6	78	-0.03
19 T	Methylene Chloride	1.421	1.258	11.5	82	-0.01
20 T	3-Chloro-1-propene (Allyl C	2.255	2.357	-4.5	85	-0.02
21 T	Trichlorotrifluoroethane	1.171	1.062	9.3	80	0.00
22 T	Carbon Disulfide	5.030	4.631	7.9	82	0.00
23 T	trans-1,2-Dichloroethene	2.080	1.980	4.8	84	-0.01
24 T	1,1-Dichloroethane	2.514	2.403	4.4	84	-0.02
25 T	Methyl tert-Butyl Ether	4.378	4.257	2.8	85	0.00
26 T	Vinyl Acetate	0.403	0.410	-1.7	83	-0.03
27 T	2-Butanone (MEK)	0.986	0.978	0.8	84	-0.03
28 T	cis-1,2-Dichloroethene	1.960	1.861	5.1	84	-0.02
29 T	Diisopropyl Ether	1.239	1.185	4.4	83	-0.01
30 T	Ethyl Acetate	0.546	0.540	1.1	84	-0.03
31 T	n-Hexane	2.855	2.629	7.9	83	-0.01
32 T	Chloroform	2.206	2.043	7.4	83	-0.03
33 S	1,2-Dichloroethane-d4 (SS1)	1.553	1.659	-6.8	90	-0.01
34 T	Tetrahydrofuran (THF)	0.961	0.885	7.9	84	-0.01
35 T	Ethyl tert-Butyl Ether	1.821	1.752	3.8	83	-0.01
36 T	1,2-Dichloroethane	1.798	1.746	2.9	85	-0.02
37 IR	1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	87	0.00
38 T	1,1,1-Trichloroethane	0.451	0.419	7.1	84	0.00

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\21\  
 Data File : 03211201.D  
 Acq On : 21 Mar 2012 4:13  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 21 07:38:45 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
39 T	Isopropyl Acetate	0.204	0.195	4.4	84	-0.02
40 T	1-Butanol	0.323	0.351	-8.7	91	-0.05
41 T	Benzene	1.258	1.108	11.9	83	-0.02
42 T	Carbon Tetrachloride	0.352	0.330	6.2	81	-0.01
43 T	Cyclohexane	0.539	0.483	10.4	84	-0.02
44 T	tert-Amyl Methyl Ether	0.939	0.885	5.8	84	-0.02
45 T	1,2-Dichloropropane	0.334	0.306	8.4	84	-0.01
46 T	Bromodichloromethane	0.384	0.360	6.3	83	-0.01
47 T	Trichloroethene	0.355	0.306	13.8	81	-0.01
48 T	1,4-Dioxane	0.260	0.240	7.7	85	-0.02
49 T	2,2,4-Trimethylpentane (Iso	1.557	1.440	7.5	85	-0.01
50 T	Methyl Methacrylate	0.141	0.136	3.5	83	-0.02
51 T	n-Heptane	0.361	0.337	6.6	85	-0.01
52 T	cis-1,3-Dichloropropene	0.502	0.482	4.0	84	0.00
53 T	4-Methyl-2-pentanone	0.305	0.303	0.7	87	-0.02
54 T	trans-1,3-Dichloropropene	0.437	0.437	0.0	83	-0.01
55 T	1,1,2-Trichloroethane	0.292	0.269	7.9	82	-0.01
56 IR	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	87	0.00
57 S	Toluene-d8 (SS2)	2.306	2.280	1.1	86	0.00
58 T	Toluene	2.793	2.457	12.0	83	-0.01
59 T	2-Hexanone	1.490	1.438	3.5	88	-0.02
60 T	Dibromochloromethane	0.638	0.606	5.0	81	0.00
61 T	1,2-Dibromoethane	0.688	0.621	9.7	81	-0.01
62 T	n-Butyl Acetate	1.839	1.765	4.0	87	-0.01
63 T	n-Octane	0.645	0.625	3.1	89	-0.01
64 T	Tetrachloroethene	0.848	0.719	15.2	80	0.00
65 T	Chlorobenzene	1.780	1.542	13.4	82	0.00
66 T	Ethylbenzene	3.030	2.707	10.7	83	0.00
67 T	m- & p-Xylenes	2.419	2.215	8.4	85	-0.02
68 T	Bromoform	0.568	0.556	2.1	81	-0.01
69 T	Styrene	1.881	1.754	6.8	84	-0.01
70 T	o-Xylene	2.550	2.308	9.5	84	-0.01
71 T	n-Nonane	1.681	1.571	6.5	85	-0.01
72 T	1,1,2,2-Tetrachloroethane	1.160	1.066	8.1	83	-0.02
73 S	Bromofluorobenzene (SS3)	0.777	0.752	3.2	84	0.00
74 T	Cumene	3.299	2.988	9.4	86	-0.01
75 T	alpha-Pinene	1.611	1.480	8.1	81	-0.01
76 T	n-Propylbenzene	4.039	3.645	9.8	84	0.00

*Em 3/21/12*



Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\21\  
 Data File : 03211201.D  
 Acq On : 21 Mar 2012 4:13  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 21 07:38:45 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
77 T	3-Ethyltoluene	3.206	2.881	10.1	84	-0.01
78 T	4-Ethyltoluene	3.133	2.806	10.4	82	-0.02
79 T	1,3,5-Trimethylbenzene	2.600	2.325	10.6	83	-0.01
80 T	alpha-Methylstyrene	1.436	1.326	7.7	82	-0.01
81 T	2-Ethyltoluene	3.254	2.901	10.8	83	-0.01
82 T	1,2,4-Trimethylbenzene	2.664	2.417	9.3	83	-0.02
83 T	n-Decane	1.699	1.581	6.9	84	-0.01
84 T	Benzyl Chloride	1.948	2.028	-4.1	86	-0.02
85 T	1,3-Dichlorobenzene	1.607	1.306	18.7	80	-0.01
86 T	1,4-Dichlorobenzene	1.649	1.324	19.7	80	-0.02
87 T	sec-Butylbenzene	3.571	3.214	10.0	84	-0.01
88 T	4-Isopropyltoluene (p-Cymen	3.554	3.220	9.4	83	-0.01
89 T	1,2,3-Trimethylbenzene	2.691	2.461	8.5	83	-0.01
90 T	1,2-Dichlorobenzene	1.531	1.292	15.6	81	-0.01
91 T	d-Limonene	1.067	1.038	2.7	84	-0.01
92 T	1,2-Dibromo-3-Chloropropane	0.517	0.488	5.6	81	-0.01
93 T	n-Undecane	1.678	1.592	5.1	85	0.00
94 T	1,2,4-Trichlorobenzene	1.122	0.982	12.5	80	-0.01
95 T	Naphthalene	3.881	3.509	9.6	82	-0.01
96 T	n-Dodecane	1.613	1.592	1.3	88	0.00
97 T	Hexachlorobutadiene	0.704	0.608	13.6	80	0.00
98 T	Cyclohexanone	0.986	1.003	-1.7	91	-0.02
99 T	tert-Butylbenzene	2.723	2.453	9.9	84	-0.02
100 T	n-Butylbenzene	2.817	2.521	10.5	83	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*EM 3/21/12*

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\22\  
 Data File : 03221201.D  
 Acq On : 22 Mar 2012 2:49  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 22 08:55:06 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
1	IR Bromochloromethane (IS1)	1.000	1.000	0.0	86	-0.01
2	T Propene	2.102	2.177	-3.6	81	0.00
3	T Dichlorodifluoromethane (CF	2.664	2.284	14.3	83	0.00
4	T Chloromethane	2.449	2.142	12.5	80	0.00
5	T 1,2-Dichloro-1,1,2,2-tetra	1.460	1.223	16.2	79	0.00
6	T Vinyl Chloride	2.333	2.030	13.0	82	0.00
7	T 1,3-Butadiene	1.743	1.690	3.0	86	0.00
8	T Bromomethane	1.252	1.088	13.1	82	0.00
9	T Chloroethane	1.217	1.072	11.9	84	0.00
10	T Ethanol	1.239	1.208	2.5	94	-0.07
11	T Acetonitrile	2.825	2.612	7.5	87	-0.05
12	T Acrolein	0.897	0.768	14.4	81	-0.02
13	T Acetone	1.161	1.021	12.1	86	-0.05
14	T Trichlorofluoromethane	2.277	2.085	8.4	84	0.00
15	T 2-Propanol (Isopropanol)	3.443	2.917	15.3	90	-0.05
16	T Acrylonitrile	1.916	1.802	5.9	85	-0.03
17	T 1,1-Dichloroethene	1.293	1.158	10.4	81	0.00
18	T 2-Methyl-2-Propanol (tert-B	3.812	3.266	14.3	78	-0.03
19	T Methylene Chloride	1.421	1.230	13.4	83	-0.01
20	T 3-Chloro-1-propene (Allyl C	2.255	2.332	-3.4	87	-0.01
21	T Trichlorotrifluoroethane	1.171	1.050	10.3	82	0.00
22	T Carbon Disulfide	5.030	4.560	9.3	83	0.00
23	T trans-1,2-Dichloroethene	2.080	1.936	6.9	85	-0.01
24	T 1,1-Dichloroethane	2.514	2.368	5.8	85	-0.02
25	T Methyl tert-Butyl Ether	4.378	4.176	4.6	85	0.00
26	T Vinyl Acetate	0.403	0.404	-0.2	84	-0.03
27	T 2-Butanone (MEK)	0.986	0.952	3.4	85	-0.03
28	T cis-1,2-Dichloroethene	1.960	1.817	7.3	85	-0.01
29	T Diisopropyl Ether	1.239	1.163	6.1	84	0.00
30	T Ethyl Acetate	0.546	0.535	2.0	86	-0.02
31	T n-Hexane	2.855	2.576	9.8	84	0.00
32	T Chloroform	2.206	2.014	8.7	84	-0.03
33	S 1,2-Dichloroethane-d4 (SS1)	1.553	1.633	-5.2	91	-0.01
34	T Tetrahydrofuran (THF)	0.961	0.871	9.4	85	-0.01
35	T Ethyl tert-Butyl Ether	1.821	1.717	5.7	84	-0.01
36	T 1,2-Dichloroethane	1.798	1.691	6.0	85	-0.01
37	IR 1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	88	0.00
38	T 1,1,1-Trichloroethane	0.451	0.413	8.4	84	0.00

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\22\  
 Data File : 03221201.D  
 Acq On : 22 Mar 2012 2:49  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 22 08:55:06 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
39 T	Isopropyl Acetate	0.204	0.193	5.4	85	-0.02
40 T	1-Butanol	0.323	0.355	-9.9	94	-0.05
41 T	Benzene	1.258	1.095	13.0	84	-0.01
42 T	Carbon Tetrachloride	0.352	0.327	7.1	82	-0.01
43 T	Cyclohexane	0.539	0.478	11.3	84	-0.01
44 T	tert-Amyl Methyl Ether	0.939	0.875	6.8	84	-0.01
45 T	1,2-Dichloropropane	0.334	0.301	9.9	84	-0.01
46 T	Bromodichloromethane	0.384	0.359	6.5	84	0.00
47 T	Trichloroethene	0.355	0.302	14.9	81	-0.01
48 T	1,4-Dioxane	0.260	0.239	8.1	86	-0.02
49 T	2,2,4-Trimethylpentane (Iso	1.557	1.417	9.0	86	-0.01
50 T	Methyl Methacrylate	0.141	0.134	5.0	84	-0.02
51 T	n-Heptane	0.361	0.333	7.8	85	-0.01
52 T	cis-1,3-Dichloropropene	0.502	0.480	4.4	85	0.00
53 T	4-Methyl-2-pentanone	0.305	0.298	2.3	87	-0.02
54 T	trans-1,3-Dichloropropene	0.437	0.432	1.1	84	-0.01
55 T	1,1,2-Trichloroethane	0.292	0.265	9.2	82	-0.01
56 IR	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	90	0.00
57 S	Toluene-d8 (SS2)	2.306	2.283	1.0	89	0.00
58 T	Toluene	2.793	2.406	13.9	84	0.00
59 T	2-Hexanone	1.490	1.411	5.3	89	-0.02
60 T	Dibromochloromethane	0.638	0.592	7.2	82	0.00
61 T	1,2-Dibromoethane	0.688	0.603	12.4	82	-0.01
62 T	n-Butyl Acetate	1.839	1.728	6.0	88	-0.01
63 T	n-Octane	0.645	0.608	5.7	89	-0.01
64 T	Tetrachloroethene	0.848	0.698	17.7	80	0.00
65 T	Chlorobenzene	1.780	1.521	14.6	83	0.00
66 T	Ethylbenzene	3.030	2.626	13.3	83	0.00
67 T	m- & p-Xylenes	2.419	2.167	10.4	86	-0.02
68 T	Bromoform	0.568	0.544	4.2	82	0.00
69 T	Styrene	1.881	1.704	9.4	85	-0.01
70 T	o-Xylene	2.550	2.242	12.1	85	-0.01
71 T	n-Nonane	1.681	1.532	8.9	86	-0.01
72 T	1,1,2,2-Tetrachloroethane	1.160	1.040	10.3	83	-0.02
73 S	Bromofluorobenzene (SS3)	0.777	0.770	0.9	89	0.00
74 T	Cumene	3.299	2.905	11.9	87	0.00
75 T	alpha-Pinene	1.611	1.440	10.6	82	0.00
76 T	n-Propylbenzene	4.039	3.532	12.6	84	0.00

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\22\  
 Data File : 03221201.D  
 Acq On : 22 Mar 2012 2:49  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 22 08:55:06 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
77 T	3-Ethyltoluene	3.206	2.780	13.3	84	-0.01
78 T	4-Ethyltoluene	3.133	2.755	12.1	83	-0.01
79 T	1,3,5-Trimethylbenzene	2.600	2.260	13.1	84	-0.01
80 T	alpha-Methylstyrene	1.436	1.281	10.8	82	-0.01
81 T	2-Ethyltoluene	3.254	2.816	13.5	83	-0.01
82 T	1,2,4-Trimethylbenzene	2.664	2.332	12.5	83	-0.01
83 T	n-Decane	1.699	1.533	9.8	85	-0.01
84 T	Benzyl Chloride	1.948	2.018	-3.6	88	-0.02
85 T	1,3-Dichlorobenzene	1.607	1.274	20.7	81	-0.01
86 T	1,4-Dichlorobenzene	1.649	1.279	22.4	80	-0.01
87 T	sec-Butylbenzene	3.571	3.095	13.3	83	-0.01
88 T	4-Isopropyltoluene (p-Cymen	3.554	3.108	12.5	83	-0.01
89 T	1,2,3-Trimethylbenzene	2.691	2.390	11.2	83	-0.01
90 T	1,2-Dichlorobenzene	1.531	1.254	18.1	81	-0.01
91 T	d-Limonene	1.067	1.005	5.8	84	-0.01
92 T	1,2-Dibromo-3-Chloropropane	0.517	0.476	7.9	81	-0.01
93 T	n-Undecane	1.678	1.540	8.2	85	0.00
94 T	1,2,4-Trichlorobenzene	1.122	0.940	16.2	79	-0.01
95 T	Naphthalene	3.881	3.420	11.9	82	-0.01
96 T	n-Dodecane	1.613	1.535	4.8	87	0.00
97 T	Hexachlorobutadiene	0.704	0.582	17.3	79	0.00
98 T	Cyclohexanone	0.986	0.974	1.2	91	-0.02
99 T	tert-Butylbenzene	2.723	2.366	13.1	84	-0.01
100 T	n-Butylbenzene	2.817	2.442	13.3	83	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*EM 3/22/12*

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_03\22\  
 Data File : 03221204.D  
 Acq On : 22 Mar 2012 5:21 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00069  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Mar 22 06:57:51 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane (IS1)	1.000	1.000	0.0	97	0.00
2 T	Dichlorodifluoromethane (CF	3.313	3.019	8.9	89	0.00
3 T	Chloromethane	0.949	0.890	6.2	95	0.00
4 T	Vinyl Chloride	2.615	2.487	4.9	92	0.00
5 T	Bromomethane	1.407	1.417	-0.7	102	0.00
6 T	Chloroethane	1.340	1.295	3.4	97	0.00
7 T	Acetone	1.208	1.294	-7.1	103	0.00
8 T	Trichlorofluoromethane	2.761	2.600	5.8	93	0.00
9 T	1,1-Dichloroethene	1.376	1.351	1.8	99	0.00
10 T	Methylene Chloride	1.633	1.511	7.5	95	0.00
11 T	Trichlorotrifluoroethane	1.220	1.203	1.4	99	0.00
12 T	trans-1,2-Dichloroethene	1.588	1.483	6.6	95	0.00
13 T	1,1-Dichloroethane	2.980	2.876	3.5	95	0.00
14 T	Methyl tert-Butyl Ether	4.613	4.232	8.3	94	0.00
15 T	cis-1,2-Dichloroethene	1.591	1.557	2.1	99	0.00
16 T	Chloroform	2.636	2.617	0.7	98	0.00
17 S	1,2-Dichloroethane-d4 (SS1)	1.983	1.977	0.3	96	0.00
18 T	1,2-Dichloroethane	2.293	2.073	9.6	90	0.00
19 T	1,1,1-Trichloroethane	2.291	2.199	4.0	97	0.00
20 T	Benzene	6.222	6.262	-0.6	100	0.00
21 T	Carbon Tetrachloride	1.790	1.688	5.7	97	0.00
22 I	1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	106	0.00
23 T	1,2-Dichloropropane	0.366	0.322	12.0	96	0.00
24 T	Bromodichloromethane	0.439	0.383	12.8	96	0.00
25 T	Trichloroethene	0.343	0.303	11.7	99	0.00
26 T	1,4-Dioxane	0.263	0.233	11.4	102	0.00
27 T	cis-1,3-Dichloropropene	0.521	0.474	9.0	100	0.00
28 T	trans-1,3-Dichloropropene	0.444	0.403	9.2	101	0.00
29 T	1,1,2-Trichloroethane	0.276	0.252	8.7	101	0.00
30 S	Toluene-d8 (SS2)	1.056	1.191	-12.8	120	0.00
31 T	Toluene	1.374	1.230	10.5	102	0.00
32 T	1,2-Dibromoethane	0.340	0.316	7.1	103	0.00
33 T	Tetrachloroethene	0.359	0.336	6.4	104	0.00
34 I	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	105	0.00
35 T	Chlorobenzene	3.844	3.647	5.1	104	0.00
36 T	Ethylbenzene	6.608	5.864	11.3	98	0.00
37 T	m,p-Xylene	5.125	4.630	9.7	99	0.00

*Handwritten signature/initials*

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_03\22\  
 Data File : 03221204.D  
 Acq On : 22 Mar 2012 5:21 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00069  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Mar 22 06:57:51 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
38 T	o-Xylene	5.490	5.028	8.4	102	0.00
39 T	1,1,2,2-Tetrachloroethane	2.561	2.469	3.6	107	0.00
40 S	Bromofluorobenzene (SS3)	1.675	1.947	-16.2	122	0.00
41 T	1,3-Dichlorobenzene	3.015	3.092	-2.6	115	0.00
42 T	1,4-Dichlorobenzene	3.034	3.052	-0.6	115	0.00
43 T	1,2-Dichlorobenzene	2.886	2.957	-2.5	117	0.00
44 T	1,2,4-Trichlorobenzene	1.907	2.112	-10.7	135	0.00
45 T	Naphthalene	6.232	7.427	-19.2	153	0.00
46 T	Hexachlorobutadiene	1.216	1.227	-0.9	122	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*KR 3/22/12*

Response Factor Report MS07

Method : J:\Ms07\METHODS\X7021712B.M (RTE Integrator)  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
 1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	Avg	%RSD
1) I Bromochloromethan	3.784	3.527	3.467	3.822	3.279	2.986	2.875	3.155	2.920	3.313	10.83
2) T Dichlorodifluorom		1.470	0.956	1.190	0.904	0.778	0.569	0.897	0.827	0.949	28.81
3) T Chloromethane	2.802	2.749	2.588	3.044	2.597	2.406	2.314	2.588	2.446	2.615	8.57
4) T Vinyl Chloride	1.894	1.496	1.211	1.613	1.335	1.208	1.105	1.433	1.367	1.407	17.11
5) T Bromomethane	1.511	1.416	1.357	1.517	1.295	1.237	1.195	1.312	1.224	1.340	8.92
6) T Chloroethane		1.548	1.214	1.122	1.080	1.193	1.090			1.208	14.52
7) T Acetone	3.051	2.887	2.831	3.145	2.692	2.565	2.447	2.691	2.536	2.761	8.59
8) T Trichlorofluorome	1.591	1.440	1.366	1.547	1.322	1.229	1.214	1.385	1.295	1.376	9.52
9) T 1,1-Dichloroethen		2.134	1.795	1.866	1.536	1.426	1.404	1.494	1.411	1.633	16.43
10) T Methylene Chlorid	1.408	1.325	1.295	1.376	1.169	1.081	1.062	1.170	1.097	1.220	10.87
11) T Trichlorotrifluor	1.972	1.710	1.589	1.762	1.505	1.406	1.393	1.512	1.439	1.588	12.18
12) T trans-1,2-Dichlor	3.265	3.020	3.016	3.383	2.925	2.756	2.681	2.937	2.841	2.980	7.60
13) T 1,1-Dichloroethan	5.168	4.740	4.762	5.045	4.335	4.137	4.166	4.690	4.471	4.613	7.90
14) T Methyl tert-Butyl	1.885	1.626	1.630	1.776	1.524	1.434	1.388	1.575	1.482	1.591	10.09
15) T cis-1,2-Dichloroe		3.041	3.257	2.586	2.390	2.280	2.524	2.374		2.636	14.03
16) T Chloroform	2.000	2.002	2.032	2.018	1.989	1.973	1.960	1.938	1.932	1.983	1.75
17) S 1,2-Dichloroethan	2.582	2.460	2.375	2.604	2.222	2.090	1.988	2.222	2.094	2.293	9.73
18) T 1,2-Dichloroethan	2.587	2.431	2.340	2.573	2.191	2.059	2.007	2.278	2.150	2.291	9.18
19) T 1,1,1-Trichloroet		7.536	7.534	6.055	5.665	5.665	5.395	5.903	5.468	6.222	14.88
20) T Benzene	2.280	1.842	1.769	1.950	1.676	1.579	1.549	1.771	1.692	1.790	12.41
21) T Carbon Tetrachlor											
22) I 1,4-Difluorobenze	0.417	0.397	0.373	0.409	0.354	0.331	0.319	0.358	0.336	0.366	9.68
23) T 1,2-Dichloropropa	0.491	0.464	0.436	0.489	0.420	0.398	0.394	0.447	0.416	0.439	8.26
24) T Bromodichlorometh	0.418	0.379	0.349	0.390	0.324	0.305	0.287	0.326	0.307	0.343	12.96
25) T Trichloroethene	0.324	0.290	0.270	0.278	0.242	0.231	0.227	0.258	0.243	0.263	11.99
26) T 1,4-Dioxane	0.567	0.523	0.508	0.540	0.499	0.484	0.488	0.559	0.522	0.521	5.67
27) T cis-1,3-Dichlorop	0.457	0.434	0.431	0.444	0.419	0.409	0.417	0.505	0.482	0.444	7.22
28) T trans-1,3-Dichlor											

(#) = Out of Range ## Number of calibration levels exceeded format ##  
 X7021712B.M Thu Mar 22 06:58:21 2012

Response Factor Report MS07

Method : J:\Ms07\METHODS\X7021712B.M (RTE Integrator)  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
 1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	Avg	%RSD
29) T 1,1,2-Trichloroet	0.315	0.296	0.286	0.305	0.263	0.248	0.242	0.274	0.257	0.276	9.32
30) S Toluene-d8 (SS2)	1.053	1.071	1.054	1.047	1.051	1.052	1.071	1.060	1.049	1.056	0.84
31) T Toluene	1.782	1.541	1.420	1.485	1.269	1.198	1.177	1.307	1.186	1.374	14.76
32) T 1,2-Dibromoethane	0.379	0.362	0.345	0.362	0.323	0.307	0.304	0.351	0.331	0.340	7.65
33) T Tetrachloroethene	0.423	0.394	0.371	0.400	0.341	0.320	0.311	0.347	0.326	0.359	11.03
-----ISTD-----											
34) I Chlorobenzene-d5											
35) T Chlorobenzene	4.581	4.172	4.017	4.308	3.688	3.469	3.311	3.705	3.345	3.844	11.68
36) T Ethylbenzene	7.604	7.082	6.707	7.235	6.297	6.070	5.915	6.682	5.878	6.608	9.30
37) T m,p-Xylene	5.861	5.347	5.132	5.536	4.930	4.777	4.719	5.268	4.558	5.125	8.24
38) T o-Xylene	6.505	5.826	5.526	5.889	5.188	4.995	4.950	5.594	4.932	5.490	9.67
39) T 1,1,2,2-Tetrachlo	2.655	2.674	2.710	2.646	2.437	2.371	2.298	2.761	2.498	2.561	6.40
40) S Bromofluorobenze	1.712	1.701	1.695	1.697	1.683	1.680	1.670	1.642	1.593	1.675	2.20
41) T 1,3-Dichlorobenze	3.304	3.261	3.222	3.204	2.835	2.719	2.608	3.138	2.846	3.015	8.71
42) T 1,4-Dichlorobenze	3.470	3.303	3.208	3.208	2.803	2.671	2.579	3.174	2.891	3.034	10.13
43) T 1,2-Dichlorobenze	3.271	3.123	3.112	3.026	2.662	2.536	2.418	3.053	2.772	2.886	10.31
44) T 1,2,4-Trichlorobe	2.184	2.118	2.169	1.753	1.641	1.608	1.575	2.139	1.978	1.907	13.66
45) T Naphthalene	6.857	6.355	6.704	5.143	5.107	5.172	5.297	8.126	7.328	6.232	17.82
46) T Hexachlorobutadie	1.486	1.423	1.403	1.217	1.055	1.013	0.975	1.227	1.145	1.216	15.40

(#) = Out of Range ### Number of calibration levels exceeded format ###  
 X7021712B.M Thu Mar 22 06:58:21 2012



Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_03\23\  
 Data File : 03231204.D  
 Acq On : 23 Mar 2012 5:21  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00069  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Mar 23 06:42:58 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane (IS1)	1.000	1.000	0.0	98	0.00
2 T	Dichlorodifluoromethane (CF)	3.313	2.906	12.3	87	0.00
3 T	Chloromethane	0.949	0.826	13.0	89	0.00
4 T	Vinyl Chloride	2.615	2.307	11.8	87	0.00
5 T	Bromomethane	1.407	1.342	4.6	98	0.00
6 T	Chloroethane	1.340	1.216	9.3	92	0.00
7 T	Acetone	1.208	1.242	-2.8	100	0.00
8 T	Trichlorofluoromethane	2.761	2.447	11.4	89	0.00
9 T	1,1-Dichloroethene	1.376	1.302	5.4	96	0.00
10 T	Methylene Chloride	1.633	1.454	11.0	93	0.00
11 T	Trichlorotrifluoroethane	1.220	1.190	2.5	100	0.00
12 T	trans-1,2-Dichloroethene	1.588	1.441	9.3	94	0.00
13 T	1,1-Dichloroethane	2.980	2.739	8.1	92	0.00
14 T	Methyl tert-Butyl Ether	4.613	4.149	10.1	94	0.00
15 T	cis-1,2-Dichloroethene	1.591	1.514	4.8	97	0.00
16 T	Chloroform	2.636	2.525	4.2	95	0.00
17 S	1,2-Dichloroethane-d4 (SS1)	1.983	1.948	1.8	96	0.00
18 T	1,2-Dichloroethane	2.293	1.984	13.5	87	0.00
19 T	1,1,1-Trichloroethane	2.291	2.107	8.0	94	0.00
20 T	Benzene	6.222	6.119	1.7	99	0.00
21 T	Carbon Tetrachloride	1.790	1.600	10.6	93	0.00
22 I	1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	106	0.00
23 T	1,2-Dichloropropane	0.366	0.315	13.9	94	0.00
24 T	Bromodichloromethane	0.439	0.369	15.9	93	0.00
25 T	Trichloroethene	0.343	0.295	14.0	97	0.00
26 T	1,4-Dioxane	0.263	0.226	14.1	99	0.00
27 T	cis-1,3-Dichloropropene	0.521	0.455	12.7	97	0.00
28 T	trans-1,3-Dichloropropene	0.444	0.391	11.9	99	0.00
29 T	1,1,2-Trichloroethane	0.276	0.239	13.4	96	0.00
30 S	Toluene-d8 (SS2)	1.056	1.185	-12.2	120	0.00
31 T	Toluene	1.374	1.185	13.8	99	0.00
32 T	1,2-Dibromoethane	0.340	0.302	11.2	99	0.00
33 T	Tetrachloroethene	0.359	0.322	10.3	100	0.00
34 I	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	105	0.00
35 T	Chlorobenzene	3.844	3.553	7.6	101	0.00
36 T	Ethylbenzene	6.608	5.962	9.8	99	0.00
37 T	m,p-Xylene	5.125	4.745	7.4	101	0.00

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_03\23\  
 Data File : 03231204.D  
 Acq On : 23 Mar 2012 5:21  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00069  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Mar 23 06:42:58 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
38 T	o-Xylene	5.490	5.051	8.0	102	0.00
39 T	1,1,2,2-Tetrachloroethane	2.561	2.386	6.8	102	0.00
40 S	Bromofluorobenzene (SS3)	1.675	2.041	-21.9	127	0.00
41 T	1,3-Dichlorobenzene	3.015	3.016	-0.0	111	0.00
42 T	1,4-Dichlorobenzene	3.034	2.996	1.3	112	0.00
43 T	1,2-Dichlorobenzene	2.886	2.910	-0.8	114	0.00
44 T	1,2,4-Trichlorobenzene	1.907	2.116	-11.0	135	0.00
45 T	Naphthalene	6.232	7.511	-20.5	154	0.00
46 T	Hexachlorobutadiene	1.216	1.242	-2.1	123	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

## LABORATORY REPORT

March 28, 2012

Christoher Cote  
Montana DEQ  
1100 N. Last Chance Gulch  
Helena, MT 59601

### **RE: Background Indoor Air Study**

Dear Christoher:

Enclosed are the results of the sample submitted to our laboratory on March 14, 2012. For your reference, these analyses have been assigned our service request number P1201004.

All analyses were performed according to our laboratory's NELAP and DoD-ELAP-approved quality assurance program. The test results meet requirements of the current NELAP and DoD-ELAP standards, where applicable, and except as noted in the laboratory case narrative provided. For a specific list of NELAP and DoD-ELAP-accredited analytes, refer to the certifications section at [www.caslab.com](http://www.caslab.com). Results are intended to be considered in their entirety and apply only to the samples analyzed and reported herein.

Columbia Analytical Services, Inc. is certified by the California Department of Health Services, NELAP Laboratory Certificate No. 02115CA; Arizona Department of Health Services, Certificate No. AZ0694; Florida Department of Health, NELAP Certification E871020; New Jersey Department of Environmental Protection, NELAP Laboratory Certification ID #CA009; New York State Department of Health, NELAP NY Lab ID No: 11221; Oregon Environmental Laboratory Accreditation Program, NELAP ID: CA20007; The American Industrial Hygiene Association, Laboratory #101661; United States Department of Defense Environmental Laboratory Accreditation Program (DoD-ELAP), Certificate No. L11-203; Pennsylvania Registration No. 68-03307; TX Commission of Environmental Quality, NELAP ID T104704413-11-2; Minnesota Department of Health, NELAP Certificate No. 362188; Washington State Department of Ecology, ELAP Lab ID: C946, State of Utah Department of Health, NELAP Certificate No. CA015272011-1; Los Angeles Department of Building and Safety, Approval No: TA00001. Each of the certifications listed above have an explicit Scope of Accreditation that applies to specific matrices/methods/analytes; therefore, please contact me for information corresponding to a particular certification.

If you have any questions, please call me at (805) 526-7161.

Respectfully submitted,

**Columbia Analytical Services, Inc.**

Kate Aguilera  
Project Manager

Client: Montana DEQ  
Project: Background Indoor Air Study

CAS Project No: P1201004

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## CASE NARRATIVE

The sample was received intact under chain of custody on March 14, 2012 and was stored in accordance with the analytical method requirements. Please refer to the sample acceptance check form for additional information. The results reported herein are applicable only to the condition of the sample at the time of sample receipt.

### Air-Phase Petroleum Hydrocarbons (APH) Analysis

The samples were analyzed for total aliphatic and aromatic gasoline range hydrocarbons by gas chromatography/mass spectrometry according to the Method for the Determination of Air-Phase Petroleum Hydrocarbons (APH), Massachusetts Department of Environmental Protection, Revision 1, December, 2009.

### Volatile Organic Compound Analysis

The samples were also analyzed in SIM and Scan mode for selected volatile organic compounds in accordance with EPA Method TO-15 from the Compendium of Methods for the Determination of Toxic Organic Compounds in Ambient Air, Second Edition (EPA/625/R-96/010b), January, 1999. The analytical system was comprised of a gas chromatograph / mass spectrometer (GC/MS) interfaced to a whole-air preconcentrator.

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*The results of analyses are given in the attached laboratory report. All results are intended to be considered in their entirety, and Columbia Analytical Services, Inc. (CAS) is not responsible for utilization of less than the complete report.*

*Use of Columbia Analytical Services, Inc. (CAS) Name. Client shall not use CAS's name or trademark in any marketing or reporting materials, press releases or in any other manner ("Materials") whatsoever and shall not attribute to CAS any test result, tolerance or specification derived from CAS's data ("Attribution") without CAS's prior written consent, which may be withheld by CAS for any reason in its sole discretion. To request CAS's consent, Client shall provide copies of the proposed Materials or Attribution and describe in writing Client's proposed use of such Materials or Attribution. If CAS has not provided written approval of the Materials or Attribution within ten (10) days of receipt from Client, Client's request to use CAS's name or trademark in any Materials or Attribution shall be deemed denied. CAS may, in its discretion, reasonably charge Client for its time in reviewing Materials or Attribution requests. Client acknowledges and agrees that the unauthorized use of CAS's name or trademark may cause CAS to incur irreparable harm for which the recovery of money damages will be inadequate. Accordingly, Client acknowledges and agrees that a violation shall justify preliminary injunctive relief. For questions contact the laboratory.*

**DETAIL SUMMARY REPORT**

Client: Montana DEQ  
 Project ID: Background Indoor Air Study

Service Request: P1201004

Date Received: 3/14/2012  
 Time Received: 10:15

MA APH 1.0 - MA VOC PH Can
TO-15 - VOC Cans
TO-15 - VOC SIM

Client Sample ID	Lab Code	Matrix	Date Collected	Time Collected	Container ID	Pi1 (psig)	Pf1 (psig)	MA APH 1.0 - MA VOC PH Can	TO-15 - VOC Cans	TO-15 - VOC SIM
18-1F-031212	P1201004-001	Air	3/13/2012	11:02	AC01843	-2.66	3.59	X	X	X

2655 Park Center Drive, Suite A  
 Simi Valley, California 93065  
 Phone (805) 526-7161  
 Fax (805) 526-7270

Requested Turnaround Time in Business Days (Surcharges) please circle  
 1 Day (100%) 2 Day (75%) 3 Day (50%) 4 Day (35%) 5 Day (25%) 10 Day-Standard

CAS Project No. **P1201004**

Company Name & Address (Reporting Information) Mountain Dept of Env Quality PO Box 20090 Helena, MT 59620-0901							Project Name <b>Background Indoor Air Study</b>		CAS Contact:	Analysis Method				
Project Manager <b>Chris Cole</b>		P.O. # / Billing Information Reference as: <b>Background Indoor Air Study</b>		Sampler (Print & Sign) <b>Aimee Reynolds</b>		Flow Controller ID (Bar code # - FC #)	Canister Start Pressure (H <sub>g</sub> )	Canister End Pressure (H <sub>g</sub> /psig)	Sample Volume	Received by (Signature)	Date: <b>8/13/12</b>	Time:		
Phone <b>406-844-5078</b>	Fax <b>406-844-5050</b>	Client Sample ID <b>18-1F-03121Z</b>	Laboratory ID Number <del>18-1F-03121Z</del>	Date Collected <b>8/13/12</b>	Time Collected <b>11DZ</b>	Canister ID (Bar code # - AC, SC, etc.) <b>ACD1843</b>	Flow Controller ID (Bar code # - FC #) <b>FGA0031</b>	Canister Start Pressure (H <sub>g</sub> ) <b>AVG02193</b>	Canister End Pressure (H <sub>g</sub> /psig) <b>-6.5"</b>	Sample Volume <b>6L</b>	Received by (Signature)	Date: <b>8/13/12</b>	Time:	
Email Address for Result Reporting <b>ccolez@mt.gov</b>							Flow Controller ID (Bar code # - FC #)		Canister Start Pressure (H <sub>g</sub> )	Canister End Pressure (H <sub>g</sub> /psig)	Sample Volume	CAS Contact:		
Tier I - Results (Default if not specified) _____							Tier II (Results + QC Summaries) _____		Tier III (Results + QC & Calibration Summaries) _____		Tier IV (Data Validation Package) 10% Surcharge _____		Project Requirements (MRLs, GAPP)	
Relinquished by: (Signature) <b>Aimee Reynolds</b>							Date: <b>8/13/12</b>		Time:		Received by: (Signature)		Date: <b>8/13/12</b>	Time:

COC AIR REV 3-11

**Sample Acceptance Check Form**

Client: Montana DEQ Work order: P1201004

Project: Background Indoor Air Study

Sample(s) received on: 3/14/12 Date opened: 3/14/12 by: MZAMORA

**Note:** This form is used for all samples received by CAS. The use of this form for custody seals is strictly meant to indicate presence/absence and not as an indication of compliance or nonconformity. Thermal preservation and pH will only be evaluated either at the request of the client and/or as required by the method/SOP.

- |    |   | Yes                                 | No                                  | N/A                                 |
|----|---|-------------------------------------|-------------------------------------|-------------------------------------|
| 1  | Were <b>sample containers</b> properly marked with client sample ID?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 2  | Container(s) <b>supplied by CAS</b> ?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 3  | Did <b>sample containers</b> arrive in good condition?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 4  | Were <b>chain-of-custody</b> papers used and filled out?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 5  | Did <b>sample container labels</b> and/or tags agree with custody papers?                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 6  | Was <b>sample volume</b> received adequate for analysis?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 7  | Are samples within specified holding times?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 8  | Was proper <b>temperature</b> (thermal preservation) of cooler at receipt adhered to?                         | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 9  | Was a <b>trip blank</b> received?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| 10 | Were <b>custody seals</b> on outside of cooler/Box?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
|    | Location of seal(s)? _____ Sealing Lid?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Were signature and date included?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Were seals intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Were custody seals on outside of sample container?  | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
|    | Location of seal(s)? _____ Sealing Lid?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Were signature and date included?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Were seals intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 11 | Do containers have appropriate <b>preservation</b> , according to method/SOP or Client specified information? | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Is there a client indication that the submitted samples are <b>pH</b> preserved?                              | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Were <b>VOA vials</b> checked for presence/absence of air bubbles?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Does the client/method/SOP require that the analyst check the sample pH and <u>if necessary</u> alter it?     | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 12 | <b>Tubes:</b> Are the tubes capped and intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Do they contain moisture?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 13 | <b>Badges:</b> Are the badges properly capped and intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Are dual bed badges separated and individually capped and intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |

Lab Sample ID	Container Description	Required pH *	Received pH	Adjusted pH	VOA Headspace (Presence/Absence)	Receipt / Preservation Comments
P1201004-001.01	6.0 L Ambient Can					

Explain any discrepancies: (include lab sample ID numbers): \_\_\_\_\_

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 18-1F-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004  
CAS Sample ID: P1201004-001

Test Code: Massachusetts APH, Revision 1, December 2009  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01843

Date Collected: 3/13/12  
Date Received: 3/14/12  
Date Analyzed: 3/22/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.66      Final Pressure (psig): 3.59

Canister Dilution Factor: 1.52

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	450	30	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	62	15	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	23	7.6	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004  
 CAS Sample ID: P120321-MB

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/21/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	20	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	10	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	5.0	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004  
 CAS Sample ID: P120321-LCS

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/21/12  
 Volume(s) Analyzed: NA Liter(s)

Compound	Spike Amount ng	Result ng	% Recovery	CAS	Data Qualifier
				Acceptance Limits	
C5 - C8 Aliphatic Hydrocarbons	24.5	24.2	99	70-130	
C9 - C12 Aliphatic Hydrocarbons	25.0	24.5	98	70-130	
C9 - C10 Aromatic Hydrocarbons	48.5	43.8	90	70-130	

## RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 18-1F-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004  
 CAS Sample ID: P1201004-001

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01843

**Date Collected:** 3/13/12  
**Date Received:** 3/14/12  
**Date Analyzed:** 3/22/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -2.66      Final Pressure (psig): 3.59

Canister Dilution Factor: 1.52

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	12	0.76	7.0	0.44	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.0	0.76	0.41	0.15	
74-87-3	Chloromethane	ND	0.76	ND	0.37	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.76	ND	0.11	
75-01-4	Vinyl Chloride	ND	0.76	ND	0.30	
106-99-0	1,3-Butadiene	ND	0.76	ND	0.34	
74-83-9	Bromomethane	ND	0.76	ND	0.20	
75-00-3	Chloroethane	ND	0.76	ND	0.29	
64-17-5	Ethanol	430	7.6	230	4.0	
75-05-8	Acetonitrile	ND	0.76	ND	0.45	
107-02-8	Acrolein	ND	3.0	ND	1.3	
67-64-1	Acetone	56	7.6	24	3.2	
75-69-4	Trichlorofluoromethane	1.2	0.76	0.22	0.14	
67-63-0	2-Propanol (Isopropyl Alcohol)	76	1.5	31	0.62	
107-13-1	Acrylonitrile	ND	0.76	ND	0.35	
75-35-4	1,1-Dichloroethene	ND	0.76	ND	0.19	
75-09-2	Methylene Chloride	1.5	0.76	0.42	0.22	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.76	ND	0.24	
76-13-1	Trichlorotrifluoroethane	ND	0.76	ND	0.099	
75-15-0	Carbon Disulfide	ND	7.6	ND	2.4	
156-60-5	trans-1,2-Dichloroethene	ND	0.76	ND	0.19	
75-34-3	1,1-Dichloroethane	ND	0.76	ND	0.19	
1634-04-4	Methyl tert-Butyl Ether	ND	0.76	ND	0.21	
108-05-4	Vinyl Acetate	ND	7.6	ND	2.2	
78-93-3	2-Butanone (MEK)	ND	7.6	ND	2.6	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 18-1F-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004  
 CAS Sample ID: P1201004-001

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01843

**Date Collected:** 3/13/12  
**Date Received:** 3/14/12  
**Date Analyzed:** 3/22/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -2.66      Final Pressure (psig): 3.59

Canister Dilution Factor: 1.52

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
156-59-2	cis-1,2-Dichloroethene	ND	0.76	ND	0.19	
141-78-6	Ethyl Acetate	15	1.5	4.2	0.42	
110-54-3	n-Hexane	15	0.76	4.1	0.22	
67-66-3	Chloroform	1.8	0.76	0.36	0.16	
109-99-9	Tetrahydrofuran (THF)	ND	0.76	ND	0.26	
107-06-2	1,2-Dichloroethane	ND	0.76	ND	0.19	
71-55-6	1,1,1-Trichloroethane	11	0.76	2.0	0.14	
71-43-2	Benzene	11	0.76	3.3	0.24	
56-23-5	Carbon Tetrachloride	ND	0.76	ND	0.12	
110-82-7	Cyclohexane	3.9	1.5	1.1	0.44	
78-87-5	1,2-Dichloropropane	ND	0.76	ND	0.16	
75-27-4	Bromodichloromethane	ND	0.76	ND	0.11	
79-01-6	Trichloroethene	ND	0.76	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.76	ND	0.21	
80-62-6	Methyl Methacrylate	ND	1.5	ND	0.37	
142-82-5	n-Heptane	5.8	0.76	1.4	0.19	
10061-01-5	cis-1,3-Dichloropropene	ND	0.76	ND	0.17	
108-10-1	4-Methyl-2-pentanone	0.80	0.76	0.20	0.19	
10061-02-6	trans-1,3-Dichloropropene	ND	0.76	ND	0.17	
79-00-5	1,1,2-Trichloroethane	ND	0.76	ND	0.14	
108-88-3	Toluene	63	0.76	17	0.20	
591-78-6	2-Hexanone	ND	0.76	ND	0.19	
124-48-1	Dibromochloromethane	ND	0.76	ND	0.089	
106-93-4	1,2-Dibromoethane	ND	0.76	ND	0.099	
123-86-4	n-Butyl Acetate	1.3	0.76	0.27	0.16	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 18-1F-031212  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201004  
**CAS Sample ID:** P1201004-001

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01843

**Date Collected:** 3/13/12  
**Date Received:** 3/14/12  
**Date Analyzed:** 3/22/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -2.66      Final Pressure (psig): 3.59

Canister Dilution Factor: 1.52

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
111-65-9	n-Octane	2.3	0.76	0.50	0.16	
127-18-4	Tetrachloroethene	2.2	0.76	0.33	0.11	
108-90-7	Chlorobenzene	ND	0.76	ND	0.17	
100-41-4	Ethylbenzene	4.5	0.76	1.0	0.18	
179601-23-1	m,p-Xylenes	16	1.5	3.7	0.35	
75-25-2	Bromoform	ND	0.76	ND	0.074	
100-42-5	Styrene	ND	0.76	ND	0.18	
95-47-6	o-Xylene	5.1	0.76	1.2	0.18	
111-84-2	n-Nonane	1.5	0.76	0.29	0.14	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.76	ND	0.11	
98-82-8	Cumene	ND	0.76	ND	0.15	
80-56-8	alpha-Pinene	2.0	0.76	0.35	0.14	
103-65-1	n-Propylbenzene	0.91	0.76	0.19	0.15	
622-96-8	4-Ethyltoluene	1.5	0.76	0.31	0.15	
108-67-8	1,3,5-Trimethylbenzene	1.7	0.76	0.35	0.15	
95-63-6	1,2,4-Trimethylbenzene	5.0	0.76	1.0	0.15	
100-44-7	Benzyl Chloride	ND	0.76	ND	0.15	
541-73-1	1,3-Dichlorobenzene	ND	0.76	ND	0.13	
106-46-7	1,4-Dichlorobenzene	ND	0.76	ND	0.13	
95-50-1	1,2-Dichlorobenzene	ND	0.76	ND	0.13	
5989-27-5	d-Limonene	16	0.76	2.8	0.14	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.76	ND	0.079	
120-82-1	1,2,4-Trichlorobenzene	ND	0.76	ND	0.10	
91-20-3	Naphthalene	ND	0.76	ND	0.15	
87-68-3	Hexachlorobutadiene	ND	0.76	ND	0.071	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004  
CAS Sample ID: P120321-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	ND	0.50	ND	0.29	
75-71-8	Dichlorodifluoromethane (CFC 12)	ND	0.50	ND	0.10	
74-87-3	Chloromethane	ND	0.50	ND	0.24	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.50	ND	0.072	
75-01-4	Vinyl Chloride	ND	0.50	ND	0.20	
106-99-0	1,3-Butadiene	ND	0.50	ND	0.23	
74-83-9	Bromomethane	ND	0.50	ND	0.13	
75-00-3	Chloroethane	ND	0.50	ND	0.19	
64-17-5	Ethanol	ND	5.0	ND	2.7	
75-05-8	Acetonitrile	ND	0.50	ND	0.30	
107-02-8	Acrolein	ND	2.0	ND	0.87	
67-64-1	Acetone	ND	5.0	ND	2.1	
75-69-4	Trichlorofluoromethane	ND	0.50	ND	0.089	
67-63-0	2-Propanol (Isopropyl Alcohol)	ND	1.0	ND	0.41	
107-13-1	Acrylonitrile	ND	0.50	ND	0.23	
75-35-4	1,1-Dichloroethene	ND	0.50	ND	0.13	
75-09-2	Methylene Chloride	ND	0.50	ND	0.14	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.50	ND	0.16	
76-13-1	Trichlorotrifluoroethane	ND	0.50	ND	0.065	
75-15-0	Carbon Disulfide	ND	5.0	ND	1.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.50	ND	0.13	
75-34-3	1,1-Dichloroethane	ND	0.50	ND	0.12	
1634-04-4	Methyl tert-Butyl Ether	ND	0.50	ND	0.14	
108-05-4	Vinyl Acetate	ND	5.0	ND	1.4	
78-93-3	2-Butanone (MEK)	ND	5.0	ND	1.7	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004  
CAS Sample ID: P120321-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
156-59-2	cis-1,2-Dichloroethene	ND	0.50	ND	0.13	
141-78-6	Ethyl Acetate	ND	1.0	ND	0.28	
110-54-3	n-Hexane	ND	0.50	ND	0.14	
67-66-3	Chloroform	ND	0.50	ND	0.10	
109-99-9	Tetrahydrofuran (THF)	ND	0.50	ND	0.17	
107-06-2	1,2-Dichloroethane	ND	0.50	ND	0.12	
71-55-6	1,1,1-Trichloroethane	ND	0.50	ND	0.092	
71-43-2	Benzene	ND	0.50	ND	0.16	
56-23-5	Carbon Tetrachloride	ND	0.50	ND	0.080	
110-82-7	Cyclohexane	ND	1.0	ND	0.29	
78-87-5	1,2-Dichloropropane	ND	0.50	ND	0.11	
75-27-4	Bromodichloromethane	ND	0.50	ND	0.075	
79-01-6	Trichloroethene	ND	0.50	ND	0.093	
123-91-1	1,4-Dioxane	ND	0.50	ND	0.14	
80-62-6	Methyl Methacrylate	ND	1.0	ND	0.24	
142-82-5	n-Heptane	ND	0.50	ND	0.12	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-10-1	4-Methyl-2-pentanone	ND	0.50	ND	0.12	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	ND	0.11	
79-00-5	1,1,2-Trichloroethane	ND	0.50	ND	0.092	
108-88-3	Toluene	ND	0.50	ND	0.13	
591-78-6	2-Hexanone	ND	0.50	ND	0.12	
124-48-1	Dibromochloromethane	ND	0.50	ND	0.059	
106-93-4	1,2-Dibromoethane	ND	0.50	ND	0.065	
123-86-4	n-Butyl Acetate	ND	0.50	ND	0.11	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004  
 CAS Sample ID: P120321-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/21/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
111-65-9	n-Octane	ND	0.50	ND	0.11	
127-18-4	Tetrachloroethene	ND	0.50	ND	0.074	
108-90-7	Chlorobenzene	ND	0.50	ND	0.11	
100-41-4	Ethylbenzene	ND	0.50	ND	0.12	
179601-23-1	m,p-Xylenes	ND	1.0	ND	0.23	
75-25-2	Bromoform	ND	0.50	ND	0.048	
100-42-5	Styrene	ND	0.50	ND	0.12	
95-47-6	o-Xylene	ND	0.50	ND	0.12	
111-84-2	n-Nonane	ND	0.50	ND	0.095	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	ND	0.073	
98-82-8	Cumene	ND	0.50	ND	0.10	
80-56-8	alpha-Pinene	ND	0.50	ND	0.090	
103-65-1	n-Propylbenzene	ND	0.50	ND	0.10	
622-96-8	4-Ethyltoluene	ND	0.50	ND	0.10	
108-67-8	1,3,5-Trimethylbenzene	ND	0.50	ND	0.10	
95-63-6	1,2,4-Trimethylbenzene	ND	0.50	ND	0.10	
100-44-7	Benzyl Chloride	ND	0.50	ND	0.097	
541-73-1	1,3-Dichlorobenzene	ND	0.50	ND	0.083	
106-46-7	1,4-Dichlorobenzene	ND	0.50	ND	0.083	
95-50-1	1,2-Dichlorobenzene	ND	0.50	ND	0.083	
5989-27-5	d-Limonene	ND	0.50	ND	0.090	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.50	ND	0.052	
120-82-1	1,2,4-Trichlorobenzene	ND	0.50	ND	0.067	
91-20-3	Naphthalene	ND	0.50	ND	0.095	
87-68-3	Hexachlorobutadiene	ND	0.50	ND	0.047	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister(s)  
 Test Notes:

Date(s) Collected: 3/13/12  
 Date(s) Received: 3/14/12  
 Date(s) Analyzed: 3/21 - 3/22/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		Percent Recovered	Percent Recovered	Percent Recovered		
Method Blank	P120321-MB	104	102	94	70-130	
Lab Control Sample	P120321-LCS	107	98	98	70-130	
18-1F-031212	P1201004-001	105	98	97	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004  
CAS Sample ID: P120321-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
115-07-1	Propene	198	202	102	54-137	
75-71-8	Dichlorodifluoromethane (CFC 12)	196	167	85	55-112	
74-87-3	Chloromethane	190	165	87	66-122	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	200	167	84	64-113	
75-01-4	Vinyl Chloride	194	167	86	68-115	
106-99-0	1,3-Butadiene	204	197	97	74-142	
74-83-9	Bromomethane	194	171	88	72-124	
75-00-3	Chloroethane	196	168	86	69-115	
64-17-5	Ethanol	928	860	93	67-127	
75-05-8	Acetonitrile	194	190	98	63-126	
107-02-8	Acrolein	198	175	88	62-127	
67-64-1	Acetone	1,010	880	87	67-106	
75-69-4	Trichlorofluoromethane	202	173	86	66-121	
67-63-0	2-Propanol (Isopropyl Alcohol)	382	308	81	56-112	
107-13-1	Acrylonitrile	198	194	98	78-128	
75-35-4	1,1-Dichloroethene	212	190	90	74-116	
75-09-2	Methylene Chloride	206	177	86	69-103	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	208	211	101	76-142	
76-13-1	Trichlorotrifluoroethane	206	184	89	69-118	
75-15-0	Carbon Disulfide	208	180	87	71-112	
156-60-5	trans-1,2-Dichloroethene	196	182	93	73-121	
75-34-3	1,1-Dichloroethane	200	187	94	71-118	
1634-04-4	Methyl tert-Butyl Ether	198	188	95	72-115	
108-05-4	Vinyl Acetate	952	937	98	51-150	
78-93-3	2-Butanone (MEK)	206	196	95	80-130	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result. Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004  
 CAS Sample ID: P120321-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/21/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
156-59-2	cis-1,2-Dichloroethene	206	190	92	73-117	
141-78-6	Ethyl Acetate	398	375	94	79-126	
110-54-3	n-Hexane	198	175	88	68-109	
67-66-3	Chloroform	214	186	87	67-118	
109-99-9	Tetrahydrofuran (THF)	202	182	90	57-130	
107-06-2	1,2-Dichloroethane	200	189	95	62-121	
71-55-6	1,1,1-Trichloroethane	198	178	90	70-116	
71-43-2	Benzene	198	178	90	66-103	
56-23-5	Carbon Tetrachloride	202	187	93	68-123	
110-82-7	Cyclohexane	390	341	87	73-111	
78-87-5	1,2-Dichloropropane	198	175	88	74-114	
75-27-4	Bromodichloromethane	198	182	92	75-120	
79-01-6	Trichloroethene	194	164	85	65-109	
123-91-1	1,4-Dioxane	200	180	90	74-120	
80-62-6	Methyl Methacrylate	400	373	93	80-124	
142-82-5	n-Heptane	196	178	91	75-114	
10061-01-5	cis-1,3-Dichloropropene	188	177	94	79-120	
108-10-1	4-Methyl-2-pentanone	204	196	96	79-128	
10061-02-6	trans-1,3-Dichloropropene	210	204	97	83-131	
79-00-5	1,1,2-Trichloroethane	196	176	90	76-116	
108-88-3	Toluene	202	170	84	64-115	
591-78-6	2-Hexanone	222	203	91	73-120	
124-48-1	Dibromochloromethane	206	187	91	72-137	
106-93-4	1,2-Dibromoethane	200	171	86	70-126	
123-86-4	n-Butyl Acetate	220	201	91	65-127	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004  
CAS Sample ID: P120321-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/21/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
111-65-9	n-Octane	200	184	92	72-118	
127-18-4	Tetrachloroethene	184	149	81	63-123	
108-90-7	Chlorobenzene	202	167	83	65-117	
100-41-4	Ethylbenzene	198	168	85	69-118	
179601-23-1	m,p-Xylenes	392	342	87	68-119	
75-25-2	Bromoform	208	180	87	79-150	
100-42-5	Styrene	200	178	89	74-127	
95-47-6	o-Xylene	194	167	86	68-118	
111-84-2	n-Nonane	196	175	89	72-116	
79-34-5	1,1,2,2-Tetrachloroethane	192	163	85	72-135	
98-82-8	Cumene	190	163	86	68-119	
80-56-8	alpha-Pinene	186	165	89	70-123	
103-65-1	n-Propylbenzene	192	163	85	69-119	
622-96-8	4-Ethyltoluene	198	167	84	68-121	
108-67-8	1,3,5-Trimethylbenzene	200	172	86	67-118	
95-63-6	1,2,4-Trimethylbenzene	194	172	89	66-122	
100-44-7	Benzyl Chloride	200	193	97	73-144	
541-73-1	1,3-Dichlorobenzene	200	157	79	64-122	
106-46-7	1,4-Dichlorobenzene	206	153	74	65-125	
95-50-1	1,2-Dichlorobenzene	198	156	79	63-128	
5989-27-5	d-Limonene	200	189	95	72-126	
96-12-8	1,2-Dibromo-3-chloropropane	196	173	88	72-139	
120-82-1	1,2,4-Trichlorobenzene	196	158	81	65-139	
91-20-3	Naphthalene	172	142	83	60-142	
87-68-3	Hexachlorobutadiene	202	161	80	58-137	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result. Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 18-1F-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004  
CAS Sample ID: P1201004-001

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01843

Date Collected: 3/13/12  
Date Received: 3/14/12  
Date Analyzed: 3/22/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.66      Final Pressure (psig): 3.59

Canister Dilution Factor: 1.52

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.038	ND	0.015	
107-06-2	1,2-Dichloroethane	<b>0.38</b>	0.038	<b>0.094</b>	0.0094	
71-43-2	Benzene	<b>11</b>	0.11	<b>3.3</b>	0.036	
79-01-6	Trichloroethene	<b>0.087</b>	0.038	<b>0.016</b>	0.0071	
79-00-5	1,1,2-Trichloroethane	ND	0.15	ND	0.028	
127-18-4	Tetrachloroethene	<b>2.3</b>	0.038	<b>0.34</b>	0.0056	
100-41-4	Ethylbenzene	<b>3.9</b>	0.15	<b>0.89</b>	0.035	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.038	ND	0.0055	
91-20-3	Naphthalene	<b>0.57</b>	0.15	<b>0.11</b>	0.029	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004  
CAS Sample ID: P120322-MB

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/22/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result		MRL		Data Qualifier
		$\mu\text{g}/\text{m}^3$	$\mu\text{g}/\text{m}^3$	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.025	ND	0.0098	
107-06-2	1,2-Dichloroethane	ND	0.025	ND	0.0062	
71-43-2	Benzene	ND	0.075	ND	0.023	
79-01-6	Trichloroethene	ND	0.025	ND	0.0047	
79-00-5	1,1,2-Trichloroethane	ND	0.10	ND	0.018	
127-18-4	Tetrachloroethene	ND	0.025	ND	0.0037	
100-41-4	Ethylbenzene	ND	0.10	ND	0.023	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.025	ND	0.0036	
91-20-3	Naphthalene	ND	0.10	ND	0.019	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister(s)  
 Test Notes:

Date(s) Collected: 3/13/12  
 Date(s) Received: 3/14/12  
 Date(s) Analyzed: 3/22/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		% Recovered	% Recovered	% Recovered		
Method Blank	P120322-MB	102	116	110	70-130	
Lab Control Sample	P120322-LCS	103	115	114	70-130	
18-1F-031212	P1201004-001	98	112	97	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.

Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201004  
CAS Sample ID: P120322-LCS

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/22/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
75-01-4	Vinyl Chloride	3.88	3.45	89	56-127	
107-06-2	1,2-Dichloroethane	4.00	3.48	87	51-140	
71-43-2	Benzene	3.96	3.89	98	56-125	
79-01-6	Trichloroethene	3.88	3.29	85	51-127	
79-00-5	1,1,2-Trichloroethane	3.92	3.39	86	49-137	
127-18-4	Tetrachloroethene	3.68	3.22	88	58-134	
100-41-4	Ethylbenzene	3.96	3.23	82	56-126	
79-34-5	1,1,2,2-Tetrachloroethane	3.84	3.33	87	53-148	
91-20-3	Naphthalene	3.44	3.35	97	25-149	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.



Massachusetts APH  
Hydrocarbon Ranges

ICAL Method: M8011412A.M

ICAL Date: 1/14/12

Instrument ID: MS08

areas

masses

	areas				masses							
	0.5	1	5	25	50	100	0.5	1	5	25	50	100
<b>C5-C8 Aliphatics</b>												
Isopentane	60478	112606	488243	2480105	4710733	9222808	0.520	1.04	5.20	26.0	52.0	104
n-Hexane	59953	110631	507126	2577865	4931102	9678199	0.490	0.98	4.90	24.5	49.0	98
Cyclohexane	63203	115230	548910	2874726	5524402	10890911	0.500	1.00	5.00	25.0	50.0	100
2,3-Dimethylpentane	68648	127222	585183	2998017	5742104	11319597	0.520	1.04	5.20	26.0	52.0	104
n-Heptane	66143	120286	558074	2884817	5509801	10873046	0.490	0.98	4.90	24.5	49.0	98
n-Octane	76169	137829	632891	3286017	6205035	12092514	0.505	1.01	5.05	25.3	50.5	101
<b>area sum:</b>	<b>394594</b>	<b>723804</b>	<b>3320427</b>	<b>17101547</b>	<b>32623177</b>	<b>64077075</b>	<b>3.025</b>	<b>6.050</b>	<b>30.25</b>	<b>151.30</b>	<b>302.50</b>	<b>605.00</b>

**C9-C12 Aliphatics**

	0.5	1	5	25	50	100	0.5	1	5	25	50	100
2,3-Dimethylheptane	78780	147724	682518	3501487	6691428	13148840	0.485	0.97	4.85	24.3	48.5	97
n-Nonane	77464	140680	662221	3417128	6555165	12940625	0.485	0.97	4.85	24.3	48.5	97
n-Decane	80593	147724	705259	3651028	7021424	13870530	0.490	0.98	4.90	24.5	49.0	98
2-Ethylcyclohexane	88915	166782	776903	4026931	7749936	15415340	0.495	0.99	4.95	24.8	49.5	99
n-Undecane	80867	152557	719618	3772331	7263077	14385881	0.480	0.96	4.80	24.0	48.0	96
n-Dodecane	70045	143746	687255	3673524	7115012	14041583	0.500	1.00	5.00	25.0	50.0	100
<b>area sum:</b>	<b>476664</b>	<b>899213</b>	<b>4233774</b>	<b>22042429</b>	<b>42396042</b>	<b>83802799</b>	<b>2.935</b>	<b>5.870</b>	<b>29.35</b>	<b>146.90</b>	<b>293.50</b>	<b>587.00</b>

**C9-C10 Aromatics**

	0.5	1	5	25	50	100	0.5	1	5	25	50	100
Isopropylbenzene	10658	19169	87439	448287	849124	1695737	0.485	0.97	4.85	24.3	48.5	97
3-Ethyltoluene	11805	20564	97823	504510	977020	1973178	0.470	0.94	4.70	23.5	47.0	94
1,3,5-Trimethylbenzene	15759	29260	134073	700133	1353305	2726630	0.490	0.98	4.90	24.5	49.0	98
1,2,3-Trimethylbenzene	17195	31328	143473	757946	1493373	3063059	0.495	0.99	4.95	24.8	49.5	99
p-Isopropyltoluene	9490	18116	83401	440962	872799	1814314	0.460	0.92	4.60	23.0	46.0	92
<b>area sum:</b>	<b>64907</b>	<b>118437</b>	<b>546209</b>	<b>2851838</b>	<b>5545621</b>	<b>11272918</b>	<b>2.400</b>	<b>4.800</b>	<b>24.00</b>	<b>120.10</b>	<b>240.00</b>	<b>480.00</b>

**Massachusetts APH**  
Hydrocarbon Ranges

ICAL: M8011412A.M

ICAL Date: 1/14/12

Instrument ID: MS08

areas

<u>Internal Standards (TIC)</u>	0.5	1	5	25	50	100
Bromochloromethane (IS1)	1312793	1307309	1293586	1289283	1284194	1283817
1,4-Difluorobenzene (IS2)	2547079	2541159	2512498	2475255	2422119	2352159
Chlorobenzene-d5 (IS3)	2929189	2913374	2880258	2895080	2878314	2908396

Internal Standards (EIC)

Bromochloromethane (IS1)	249696	247899	245985	245585	244327	243561
1,4-Difluorobenzene (IS2)	1153657	1149947	1133233	1136223	1127203	1138730
Chlorobenzene-d5 (IS3)	1007702	1004638	991212	992925	982812	986950

Surrogates (TIC)

1,2-Dichloroethane-d4	0.5	1	5	25	50	100
1,2-Dichloroethane-d4	1162497	1154617	1105630	1126201	1103958	1095945
m-Toluene-d8	3508018	3490493	3455382	3466688	3437986	3457323
p-Bromofluorobenzene	2298220	2284797	2197239	2227396	2192159	2197999

RRFs

<u>C5-C8 Aliphatics</u>	0.5	1	5	25	50	100	<u>RRF<sub>avg</sub></u>	<u>%RSD</u>
	2.8268	2.6009	2.4215	2.4870	2.3919	2.3252	2.509	7.24
<u>C9-C12 Aliphatics</u>	0.5	1	5	25	50	100	<u>RRF<sub>avg</sub></u>	<u>%RSD</u>
	4.0291	3.8120	3.6383	3.7780	3.6744	3.6163	3.758	4.09
<u>C9-C10 Aromatics</u>	0.5	1	5	25	50	100	<u>RRF<sub>avg</sub></u>	<u>%RSD</u>
	0.6709	0.6140	0.5740	0.5979	0.5878	0.5949	0.607	5.63

**Massachusetts APH**  
**Continuing Calibration Verification Check Sheet**

Data File Name: 03211203.D  
 Data File Path: J:\MS08\Data\2012\_03\21\  
 Operator: EM  
 Date Acquired: 3/21/12 5:39  
 Acq. Method File: TO15.M  
 Sample Name: 25ng TO-15/MAPH CCV STD  
 Misc Info: S25-03151201/S25-03161202  
 Instrument Name: MS08

Enter RRFs from current ICAL!

Internal Standards	RT	Area
7) 1,4-Difluorobenzene (IS2)	15.89	1102904
16) Chlorobenzene-d5 (IS3)	21.69	947969

C5-C8 Aliphatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
3) Isopentane	8.19	2464516	2.569	154.9	2.39	-30	30	Pass
4) n-Hexane	13.08	2578864						
9) Cyclohexane	15.81	2862780						
10) 2,3-Dimethylpentane	16.17	3020819	Spike	ICAL				
11) n-Heptane	17.35	2903399	Amt (ng)	RRF				
14) n-Octane	20.69	3316542	151.30	2.509				
		<b>17146920</b>						

C9-C12 Aliphatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
18) 2,3-Dimethylheptane	22.31	3515879	3.955	154.6	5.23	-30	30	Pass
19) n-Nonane	23.30	3434658						
25) n-Decane	25.26	3663558						
28) Butylcyclohexane	25.99	4002852	Spike	ICAL				
29) n-Undecane	26.76	3775128	Amt (ng)	RRF				
30) n-Dodecane	28.00	3635722	146.90	3.758				
		<b>22027797</b>						

C9-C10 Aromatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
22) Isopropylbenzene	23.78	446789	0.612	121.1	0.80	-30	30	Pass
23) 1-Methyl-3-ethylbenzene	24.52	492493						
24) 1,3,5-Trimethylbenzene	24.66	681321						
26) p-Isopropyltoluene	25.68	426000	Spike	ICAL				
27) 1,2,3-Trimethylbenzene	25.68	739855	Amt (ng)	RRF				
		<b>2786458</b>	120.1	0.607				

*EM 3/21/12*

Response Factor Report MS08

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Mon Dec 12 09:45:45 2011  
 Response Via : Initial Calibration

Calibration Files

0.1 =12091118.D 0.2 =12091119.D 0.5 =12091120.D 1.0 =12091121.D 5.0 =12091122.D 25 =12091123.D  
 50 =12091124.D 100 =12091125.D

Compound	0.1	0.2	0.5	1.0	5.0	25	50	100	Avg	%RSD
1) IR Bromochloromethane...				ISTD						
2) T Propene	2.487	1.965	1.952	1.692	2.194	2.303	2.185	2.036	2.102	11.62
3) T Dichlorodifluo...	3.545	2.805	2.940	2.748	2.432	2.361	2.278	2.199	2.664	16.73
4) T Chloromethane	3.290	2.811	2.875	2.512	2.176	2.286	1.977	1.664	2.449	21.62
5) T 1,2-Dichloro-1...	1.901	1.661	1.575	1.458	1.359	1.333	1.200	1.190	1.460	16.68
6) T Vinyl Chloride	2.906	2.590	2.631	2.374	2.145	2.129	1.950	1.939	2.333	15.07
7) T 1,3-Butadiene	2.145	1.788	1.790	1.718	1.581	1.689	1.621	1.616	1.743	10.33
8) T Bromomethane	1.687	1.328	1.349	1.212	1.094	1.142	1.111	1.092	1.252	16.21
9) T Chloroethane	1.671	1.259	1.333	1.193	1.101	1.091	1.046	1.044	1.217	17.29
10) T Ethanol	2.040	1.345	1.186	1.085	1.022	1.096	1.076	1.064	1.239	27.34
11) T Acetonitrile	4.365	2.973	2.723	2.546	2.480	2.562	2.482	2.468	2.825	22.85
12) T Acrolein	1.328	0.931	0.907	0.861	0.772	0.814	0.789	0.779	0.897	20.48
13) T Acetone	1.645	1.316	1.249	1.129	0.994	1.015	0.981	0.964	1.161	20.26
14) T Trichlorofluor...	2.776	2.385	2.496	2.272	2.097	2.135	2.042	2.011	2.277	11.59
15) T 2-Propanol (Is...	4.463	4.536	4.168	2.920	2.763	2.577	2.672	3.443	26.09	26.09
16) T Acrylonitrile	2.701	1.907	1.878	1.783	1.745	1.813	1.765	1.738	1.916	16.85
17) T 1,1-Dichloroet...	1.458	1.418	1.452	1.263	1.206	1.223	1.172	1.152	1.293	9.95
18) T 2-Methyl-2-Pro...	4.464	3.710	3.939	3.616	3.558	3.584		3.812	9.14	9.14
19) T Methylene Chlo...			1.958	1.578	1.294	1.273	1.220	1.201	1.421	20.88
20) T 3-Chloro-1-pro...	2.347	2.260	2.345	2.220	2.150	2.296	2.230	2.191	2.255	3.15
21) T Trichlorotrifl...	1.348	1.244	1.299	1.193	1.089	1.098	1.062	1.035	1.171	10.01
22) T Carbon Disulfide	6.210	5.678	5.057	4.654	4.699	4.488	4.424	5.030	13.40	13.40
23) T trans-1,2-Dich...	2.436	2.237	2.206	2.142	1.951	1.953	1.883	1.835	2.080	10.00
24) T 1,1-Dichloroet...	2.896	2.757	2.747	2.500	2.307	2.379	2.288	2.236	2.514	10.07
25) T Methyl tert-Bu...	5.128	4.611	4.650	4.307	4.119	4.185	4.042	3.984	4.378	8.93
26) T Vinyl Acetate	0.430	0.368	0.423	0.399	0.393	0.410	0.402	0.398	0.403	4.70
27) T 2-Butanone (MEK)	1.265	1.003	1.038	0.972	0.927	0.964	0.928	0.794	0.986	13.56
28) T cis-1,2-Dichlo...	2.442	2.056	2.101	1.938	1.826	1.836	1.764	1.720	1.960	12.05
29) T Diisopropyl Ether	1.446	1.223	1.323	1.262	1.171	1.188	1.153	1.148	1.239	8.26
30) T Ethyl Acetate	0.611	0.554	0.573	0.545	0.517	0.535	0.520	0.516	0.546	6.02
31) T n-Hexane	3.541	3.006	3.181	2.805	2.582	2.633	2.545	2.543	2.855	12.70

Response Factor Report MS08

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

32) T	Chloroform	2.679	2.408	2.409	2.204	1.983	2.054	1.974	1.937	2.206	12.18
33) S	1,2-Dichloroet...	1.586	1.577	1.582	1.590	1.535	1.533	1.520	1.504	1.553	2.19
34) T	Tetrahydrofura...	1.274	0.932	1.100	0.938	0.910	0.879	0.830	0.821	0.961	15.96
35) T	Ethyl tert-But...	1.999	1.832	2.033	1.818	1.732	1.757	1.708	1.687	1.821	7.18
36) T	1,2-Dichloroet...	2.099	1.992	1.976	1.806	1.683	1.700	1.599	1.531	1.798	11.39
37) IR	1,4-Difluorobenzen...	-----ISTD-----									
38) T	1,1,1-Trichlor...	0.531	0.465	0.486	0.453	0.419	0.434	0.413	0.403	0.451	9.46
39) T	Isopropyl Acetate	0.227	0.205	0.220	0.206	0.194	0.201	0.192	0.188	0.204	6.71
40) T	1-Butanol		0.348	0.319	0.300	0.320	0.335	0.324	0.318	0.323	4.66
41) T	Benzene	1.633	1.385	1.359	1.227	1.136	1.158	1.101	1.070	1.258	15.11
42) T	Carbon Tetrach...	0.394	0.356	0.369	0.340	0.334	0.352	0.337	0.333	0.352	6.03
43) T	Cyclohexane	0.666	0.576	0.588	0.539	0.492	0.500	0.479	0.469	0.539	12.59
44) T	tert-Amyl Meth...	1.092	0.926	0.998	0.937	0.900	0.918	0.881	0.858	0.939	7.95
45) T	1,2-Dichloropr...	0.397	0.361	0.368	0.324	0.312	0.316	0.300	0.293	0.334	11.12
46) T	Bromodichlorom...	0.458	0.387	0.392	0.378	0.362	0.378	0.362	0.353	0.384	8.57
47) T	Trichloroethene	0.466	0.375	0.379	0.344	0.323	0.328	0.316	0.309	0.355	14.62
48) T	1,4-Dioxane	0.305	0.282	0.288	0.248	0.245	0.246	0.237	0.230	0.260	10.55
49) T	2,2,4-Trimethy...	1.964	1.629	1.686	1.546	1.447	1.461	1.387	1.337	1.557	12.98
50) T	Methyl Methacr...	0.156	0.139	0.150	0.135	0.136	0.142	0.136	0.135	0.141	5.53
51) T	n-Heptane	0.427	0.373	0.385	0.364	0.341	0.344	0.332	0.324	0.361	9.35
52) T	cis-1,3-Dichlo...	0.571	0.502	0.521	0.491	0.482	0.500	0.481	0.471	0.502	6.33
53) T	4-Methyl-2-pen...	0.359	0.295	0.308	0.303	0.297	0.303	0.289	0.284	0.305	7.68
54) T	trans-1,3-Dich...	0.476	0.413	0.443	0.421	0.429	0.453	0.437	0.427	0.437	4.59
55) T	1,1,2-Trichlor...	0.340	0.301	0.310	0.284	0.276	0.285	0.271	0.265	0.292	8.44
56) IR	Chlorobenzene-d5 (...)	-----ISTD-----									
57) S	Toluene-d8 (SS2)	2.305	2.306	2.305	2.309	2.306	2.300	2.310	2.304	2.306	0.14
58) T	Toluene	3.623	2.920	3.025	2.738	2.571	2.574	2.473	2.421	2.793	14.19
59) T	2-Hexanone	2.049	1.429	1.515	1.403	1.395	1.422	1.367	1.337	1.490	15.57
60) T	Dibromochlorom...	0.696	0.618	0.636	0.636	0.612	0.648	0.629	0.625	0.638	4.14
61) T	1,2-Dibromoethane	0.831	0.730	0.718	0.658	0.641	0.662	0.638	0.628	0.688	9.97
62) T	n-Butyl Acetate	2.338	1.850	1.857	1.721	1.704	1.765	1.737	1.737	1.839	11.40
63) T	n-Octane	0.785	0.652	0.714	0.643	0.602	0.611	0.580	0.572	0.645	11.28
64) T	Tetrachloroethene	1.090	0.904	0.904	0.826	0.767	0.779	0.756	0.755	0.848	13.63
65) T	Chlorobenzene	2.375	1.868	1.886	1.744	1.610	1.637	1.574	1.548	1.780	15.30
66) T	Ethylbenzene	3.846	3.171	3.188	2.956	2.803	2.848	2.743	2.680	3.030	12.51
67) T	m- & p-Xylenes	3.081	2.499	2.611	2.392	2.216	2.273	2.170	2.112	2.419	13.08
68) T	Bromoform	0.622	0.494	0.556	0.541	0.551	0.598	0.589	0.594	0.568	7.19
69) T	Styrene	2.478	1.895	1.937	1.800	1.735	1.803	1.726	1.675	1.881	13.64
70) T	o-Xylene	3.279	2.708	2.726	2.479	2.331	2.373	2.283	2.224	2.550	13.65

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M

Title	SOP	VOA-TO15	CASS	TO-15/GC-MS	1.518	1.470	1.681	10.69
71) T n-Nonane	2.005	1.753	1.847	1.675	1.581	1.597	1.470	1.681
72) T 1,1,2,2-Tetrac...	1.348	1.164	1.262	1.150	1.083	1.121	1.085	1.160
73) S Bromofluoroben...	0.774	0.771	0.779	0.781	0.775	0.778	0.782	0.777
74) T Cumene	4.522	3.495	3.524	3.169	2.962	3.002	2.884	2.831
75) T alpha-Pinene	1.782	1.636	1.734	1.601	1.537	1.578	1.518	1.611
76) T n-Propylbenzene	5.441	4.229	4.306	3.924	3.710	3.754	3.394	4.039
77) T 3-Ethyltoluene	4.091	3.373	3.461	3.038	2.964	2.981	2.893	3.206
78) T 4-Ethyltoluene	4.003	3.202	3.301	3.151	2.837	2.983	2.810	2.776
79) T 1,3,5-Trimethy...	3.324	2.662	2.723	2.537	2.378	2.426	2.344	2.403
80) T alpha-Methylst...	1.742	1.386	1.449	1.371	1.390	1.402	1.354	1.436
81) T 2-Ethyltoluene	4.143	3.362	3.430	3.169	2.975	3.048	2.919	2.985
82) T 1,2,4-Trimethy...	3.356	2.663	2.791	2.551	2.433	2.513	2.454	2.664
83) T n-Decane	1.969	1.702	1.844	1.705	1.582	1.626	1.554	1.609
84) T Benzyl Chloride	2.373	1.648	1.772	1.688	1.799	2.058	2.054	2.190
85) T 1,3-Dichlorobe...	2.498	1.658	1.649	1.455	1.377	1.409	1.374	1.436
86) T 1,4-Dichlorobe...	2.655	1.785	1.618	1.492	1.380	1.428	1.388	1.449
87) T sec-Butylbenzene	4.527	3.773	3.816	3.471	3.264	3.328	3.214	3.176
88) T 4-Isopropyltol...	4.426	3.674	3.683	3.416	3.254	3.356	3.253	3.366
889) T 1,2,3-Trimethy...	3.266	2.650	2.852	2.588	2.471	2.573	2.506	2.619
890) T 1,2-Dichlorobe...	2.173	1.589	1.545	1.435	1.332	1.387	1.370	1.415
891) T d-Limonene	1.124	0.998	1.106	1.069	1.064	1.072	1.040	1.065
92) T 1,2-Dibromo-3-...	0.563	0.488	0.501	0.486	0.494	0.525	0.532	0.517
93) T n-Undecane	1.922	1.637	1.755	1.671	1.608	1.630	1.608	1.678
94) T 1,2,4-Trichlor...	1.335	1.165	1.062	1.017	1.071	1.071	1.086	1.118
95) T Naphthalene	4.848	3.923	3.638	3.497	3.721	3.721	3.755	3.881
96) T n-Dodecane	1.866	1.599	1.669	1.586	1.545	1.575	1.553	1.613
97) T Hexachlorobuta...	0.841	0.749	0.730	0.673	0.635	0.664	0.663	0.704
98) T Cyclohexanone	1.172	0.979	1.021	0.969	0.952	0.962	0.931	0.986
99) T tert-Butylbenzene	3.486	2.799	2.880	2.658	2.449	2.533	2.453	2.723
100) T n-Butylbenzene	3.596	2.904	2.958	2.670	2.572	2.625	2.613	2.817

(#) = Out of Range

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\21\  
 Data File : 03211201.D  
 Acq On : 21 Mar 2012 4:13  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 21 07:38:45 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	IR Bromochloromethane (IS1)	1.000	1.000	0.0	83	-0.02
2	T Propene	2.102	2.230	-6.1	80	0.00
3	T Dichlorodifluoromethane (CF	2.664	2.371	11.0	83	0.00
4	T Chloromethane	2.449	2.229	9.0	81	0.00
5	T 1,2-Dichloro-1,1,2,2-tetra	1.460	1.253	14.2	78	0.00
6	T Vinyl Chloride	2.333	2.088	10.5	82	0.00
7	T 1,3-Butadiene	1.743	1.698	2.6	84	-0.01
8	T Bromomethane	1.252	1.103	11.9	80	0.00
9	T Chloroethane	1.217	1.086	10.8	83	0.00
10	T Ethanol	1.239	1.182	4.6	90	-0.07
11	T Acetonitrile	2.825	2.652	6.1	86	-0.05
12	T Acrolein	0.897	0.782	12.8	80	-0.02
13	T Acetone	1.161	1.046	9.9	86	-0.05
14	T Trichlorofluoromethane	2.277	2.123	6.8	83	0.00
15	T 2-Propanol (Isopropanol)	3.443	2.957	14.1	89	-0.05
16	T Acrylonitrile	1.916	1.850	3.4	85	-0.03
17	T 1,1-Dichloroethene	1.293	1.194	7.7	81	0.00
18	T 2-Methyl-2-Propanol (tert-B	3.812	3.371	11.6	78	-0.03
19	T Methylene Chloride	1.421	1.258	11.5	82	-0.01
20	T 3-Chloro-1-propene (Allyl C	2.255	2.357	-4.5	85	-0.02
21	T Trichlorotrifluoroethane	1.171	1.062	9.3	80	0.00
22	T Carbon Disulfide	5.030	4.631	7.9	82	0.00
23	T trans-1,2-Dichloroethene	2.080	1.980	4.8	84	-0.01
24	T 1,1-Dichloroethane	2.514	2.403	4.4	84	-0.02
25	T Methyl tert-Butyl Ether	4.378	4.257	2.8	85	0.00
26	T Vinyl Acetate	0.403	0.410	-1.7	83	-0.03
27	T 2-Butanone (MEK)	0.986	0.978	0.8	84	-0.03
28	T cis-1,2-Dichloroethene	1.960	1.861	5.1	84	-0.02
29	T Diisopropyl Ether	1.239	1.185	4.4	83	-0.01
30	T Ethyl Acetate	0.546	0.540	1.1	84	-0.03
31	T n-Hexane	2.855	2.629	7.9	83	-0.01
32	T Chloroform	2.206	2.043	7.4	83	-0.03
33	S 1,2-Dichloroethane-d4 (SS1)	1.553	1.659	-6.8	90	-0.01
34	T Tetrahydrofuran (THF)	0.961	0.885	7.9	84	-0.01
35	T Ethyl tert-Butyl Ether	1.821	1.752	3.8	83	-0.01
36	T 1,2-Dichloroethane	1.798	1.746	2.9	85	-0.02
37	IR 1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	87	0.00
38	T 1,1,1-Trichloroethane	0.451	0.419	7.1	84	0.00

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\21\  
 Data File : 03211201.D  
 Acq On : 21 Mar 2012 4:13  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 21 07:38:45 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
39 T	Isopropyl Acetate	0.204	0.195	4.4	84	-0.02
40 T	1-Butanol	0.323	0.351	-8.7	91	-0.05
41 T	Benzene	1.258	1.108	11.9	83	-0.02
42 T	Carbon Tetrachloride	0.352	0.330	6.2	81	-0.01
43 T	Cyclohexane	0.539	0.483	10.4	84	-0.02
44 T	tert-Amyl Methyl Ether	0.939	0.885	5.8	84	-0.02
45 T	1,2-Dichloropropane	0.334	0.306	8.4	84	-0.01
46 T	Bromodichloromethane	0.384	0.360	6.3	83	-0.01
47 T	Trichloroethene	0.355	0.306	13.8	81	-0.01
48 T	1,4-Dioxane	0.260	0.240	7.7	85	-0.02
49 T	2,2,4-Trimethylpentane (Iso	1.557	1.440	7.5	85	-0.01
50 T	Methyl Methacrylate	0.141	0.136	3.5	83	-0.02
51 T	n-Heptane	0.361	0.337	6.6	85	-0.01
52 T	cis-1,3-Dichloropropene	0.502	0.482	4.0	84	0.00
53 T	4-Methyl-2-pentanone	0.305	0.303	0.7	87	-0.02
54 T	trans-1,3-Dichloropropene	0.437	0.437	0.0	83	-0.01
55 T	1,1,2-Trichloroethane	0.292	0.269	7.9	82	-0.01
56 IR	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	87	0.00
57 S	Toluene-d8 (SS2)	2.306	2.280	1.1	86	0.00
58 T	Toluene	2.793	2.457	12.0	83	-0.01
59 T	2-Hexanone	1.490	1.438	3.5	88	-0.02
60 T	Dibromochloromethane	0.638	0.606	5.0	81	0.00
61 T	1,2-Dibromoethane	0.688	0.621	9.7	81	-0.01
62 T	n-Butyl Acetate	1.839	1.765	4.0	87	-0.01
63 T	n-Octane	0.645	0.625	3.1	89	-0.01
64 T	Tetrachloroethene	0.848	0.719	15.2	80	0.00
65 T	Chlorobenzene	1.780	1.542	13.4	82	0.00
66 T	Ethylbenzene	3.030	2.707	10.7	83	0.00
67 T	m- & p-Xylenes	2.419	2.215	8.4	85	-0.02
68 T	Bromoform	0.568	0.556	2.1	81	-0.01
69 T	Styrene	1.881	1.754	6.8	84	-0.01
70 T	o-Xylene	2.550	2.308	9.5	84	-0.01
71 T	n-Nonane	1.681	1.571	6.5	85	-0.01
72 T	1,1,2,2-Tetrachloroethane	1.160	1.066	8.1	83	-0.02
73 S	Bromofluorobenzene (SS3)	0.777	0.752	3.2	84	0.00
74 T	Cumene	3.299	2.988	9.4	86	-0.01
75 T	alpha-Pinene	1.611	1.480	8.1	81	-0.01
76 T	n-Propylbenzene	4.039	3.645	9.8	84	0.00

*Em 3/21/12*



Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\21\  
 Data File : 03211201.D  
 Acq On : 21 Mar 2012 4:13  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 21 07:38:45 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
77 T	3-Ethyltoluene	3.206	2.881	10.1	84	-0.01
78 T	4-Ethyltoluene	3.133	2.806	10.4	82	-0.02
79 T	1,3,5-Trimethylbenzene	2.600	2.325	10.6	83	-0.01
80 T	alpha-Methylstyrene	1.436	1.326	7.7	82	-0.01
81 T	2-Ethyltoluene	3.254	2.901	10.8	83	-0.01
82 T	1,2,4-Trimethylbenzene	2.664	2.417	9.3	83	-0.02
83 T	n-Decane	1.699	1.581	6.9	84	-0.01
84 T	Benzyl Chloride	1.948	2.028	-4.1	86	-0.02
85 T	1,3-Dichlorobenzene	1.607	1.306	18.7	80	-0.01
86 T	1,4-Dichlorobenzene	1.649	1.324	19.7	80	-0.02
87 T	sec-Butylbenzene	3.571	3.214	10.0	84	-0.01
88 T	4-Isopropyltoluene (p-Cymen)	3.554	3.220	9.4	83	-0.01
89 T	1,2,3-Trimethylbenzene	2.691	2.461	8.5	83	-0.01
90 T	1,2-Dichlorobenzene	1.531	1.292	15.6	81	-0.01
91 T	d-Limonene	1.067	1.038	2.7	84	-0.01
92 T	1,2-Dibromo-3-Chloropropane	0.517	0.488	5.6	81	-0.01
93 T	n-Undecane	1.678	1.592	5.1	85	0.00
94 T	1,2,4-Trichlorobenzene	1.122	0.982	12.5	80	-0.01
95 T	Naphthalene	3.881	3.509	9.6	82	-0.01
96 T	n-Dodecane	1.613	1.592	1.3	88	0.00
97 T	Hexachlorobutadiene	0.704	0.608	13.6	80	0.00
98 T	Cyclohexanone	0.986	1.003	-1.7	91	-0.02
99 T	tert-Butylbenzene	2.723	2.453	9.9	84	-0.02
100 T	n-Butylbenzene	2.817	2.521	10.5	83	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*EM 3/21/12*

Response Factor Report MS07

Method : J:\Ms07\METHODS\X7021712B.M (RTE Integrator)  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
 1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	Avg	%RSD
1) I Bromochloromethan	3.784	3.527	3.467	3.822	3.279	2.986	2.875	3.155	2.920	3.313	10.83
2) T Dichlorodifluorom		1.470	0.956	1.190	0.904	0.778	0.569	0.897	0.827	0.949	28.81
3) T Chloromethane	2.802	2.749	2.588	3.044	2.597	2.406	2.314	2.588	2.446	2.615	8.57
4) T Vinyl Chloride	1.894	1.496	1.211	1.613	1.335	1.208	1.105	1.433	1.367	1.407	17.11
5) T Bromomethane	1.511	1.416	1.357	1.517	1.295	1.237	1.195	1.312	1.224	1.340	8.92
6) T Chloroethane		1.548	1.214	1.122	1.080	1.193	1.090			1.208	14.52
7) T Acetone	3.051	2.887	2.831	3.145	2.692	2.565	2.447	2.691	2.536	2.761	8.59
8) T Trichlorofluorome	1.591	1.440	1.366	1.547	1.322	1.229	1.214	1.385	1.295	1.376	9.52
9) T 1,1-Dichloroethen	2.134	1.795	1.866	1.866	1.536	1.426	1.404	1.494	1.411	1.633	16.43
10) T Methylene Chlorid	1.408	1.325	1.295	1.376	1.169	1.081	1.062	1.170	1.097	1.220	10.87
11) T Trichlorotrifluor	1.972	1.710	1.589	1.762	1.505	1.406	1.393	1.512	1.439	1.588	12.18
12) T trans-1,2-Dichlor	3.265	3.020	3.016	3.383	2.925	2.756	2.681	2.937	2.841	2.980	7.60
13) T 1,1-Dichloroethan	5.168	4.740	4.762	5.045	4.335	4.137	4.166	4.690	4.471	4.613	7.90
14) T Methyl tert-Butyl	1.885	1.626	1.630	1.776	1.524	1.434	1.388	1.575	1.482	1.591	10.09
15) T cis-1,2-Dichloroe		3.041	3.257	2.586	2.390	2.280	2.524	2.374		2.636	14.03
16) T Chloroform	2.000	2.002	2.032	2.018	1.989	1.973	1.960	1.938	1.932	1.983	1.75
17) S 1,2-Dichloroethan	2.582	2.460	2.375	2.604	2.222	2.090	1.988	2.222	2.094	2.293	9.73
18) T 1,2-Dichloroethan	2.587	2.431	2.340	2.573	2.191	2.059	2.007	2.278	2.150	2.291	9.18
19) T 1,1,1-Trichloroet		7.536	7.534	6.055	5.665	5.395	5.903	5.468		6.222	14.88
20) T Benzene	2.280	1.842	1.769	1.950	1.676	1.579	1.549	1.771	1.692	1.790	12.41
21) T Carbon Tetrachlor											
22) I 1,4-Difluorobenze	0.417	0.397	0.373	0.409	0.354	0.331	0.319	0.358	0.336	0.366	9.68
23) T 1,2-Dichloropropa	0.491	0.464	0.436	0.489	0.420	0.398	0.394	0.447	0.416	0.439	8.26
24) T Bromodichlorometh	0.418	0.379	0.349	0.390	0.324	0.305	0.287	0.326	0.307	0.343	12.96
25) T Trichloroethene	0.324	0.290	0.270	0.278	0.242	0.231	0.227	0.258	0.243	0.263	11.99
26) T 1,4-Dioxane	0.567	0.523	0.508	0.540	0.499	0.484	0.488	0.559	0.522	0.521	5.67
27) T cis-1,3-Dichlorop	0.457	0.434	0.431	0.444	0.419	0.409	0.417	0.505	0.482	0.444	7.22
28) T trans-1,3-Dichlor											

(#) = Out of Range ### Number of calibration levels exceeded format ###  
 X7021712B.M Thu Mar 22 06:58:21 2012

Response Factor Report MS07

Method : J:\Ms07\METHODS\X7021712B.M (RTE Integrator)  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
 1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	Avg	%RSD
29) T 1,1,2-Trichloroet	0.315	0.296	0.286	0.305	0.263	0.248	0.242	0.274	0.257	0.276	9.32
30) S Toluene-d8 (SS2)	1.053	1.071	1.054	1.047	1.051	1.052	1.071	1.060	1.049	1.056	0.84
31) T Toluene	1.782	1.541	1.420	1.485	1.269	1.198	1.177	1.307	1.186	1.374	14.76
32) T 1,2-Dibromoethane	0.379	0.362	0.345	0.362	0.323	0.307	0.304	0.351	0.331	0.340	7.65
33) T Tetrachloroethene	0.423	0.394	0.371	0.400	0.341	0.320	0.311	0.347	0.326	0.359	11.03
-----ISTD-----											
34) I Chlorobenzene-d5	4.581	4.172	4.017	4.308	3.688	3.469	3.311	3.705	3.345	3.844	11.68
35) T Chlorobenzene	7.604	7.082	6.707	7.235	6.297	6.070	5.915	6.682	5.878	6.608	9.30
36) T Ethylbenzene	5.861	5.347	5.132	5.536	4.930	4.777	4.719	5.268	4.558	5.125	8.24
37) T m,p-Xylene	6.505	5.826	5.526	5.889	5.188	4.995	4.950	5.594	4.932	5.490	9.67
38) T o-Xylene	2.655	2.674	2.710	2.646	2.437	2.371	2.298	2.761	2.498	2.561	6.40
39) T 1,1,2,2-Tetrachlo	1.712	1.701	1.695	1.697	1.683	1.680	1.670	1.642	1.593	1.675	2.20
40) S Bromofluorobenzene	3.304	3.261	3.222	3.204	2.835	2.719	2.608	3.138	2.846	3.015	8.71
41) T 1,3-Dichlorobenze	3.470	3.303	3.208	3.208	2.803	2.671	2.579	3.174	2.891	3.034	10.13
42) T 1,4-Dichlorobenze	3.271	3.123	3.112	3.026	2.662	2.536	2.418	3.053	2.772	2.886	10.31
43) T 1,2-Dichlorobenze	2.184	2.118	2.169	1.753	1.641	1.608	1.575	2.139	1.978	1.907	13.66
44) T 1,2,4-Trichlorobe	6.857	6.355	6.704	5.143	5.107	5.172	5.297	8.126	7.328	6.232	17.82
45) T Naphthalene	1.486	1.423	1.403	1.217	1.055	1.013	0.975	1.227	1.145	1.216	15.40
46) T Hexachlorobutadie											

(#) = Out of Range ### Number of calibration levels exceeded format ###  
 X7021712B.M Thu Mar 22 06:58:21 2012

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_03\22\  
 Data File : 03221204.D  
 Acq On : 22 Mar 2012 5:21 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00069  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Mar 22 06:57:51 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane (IS1)	1.000	1.000	0.0	97	0.00
2 T	Dichlorodifluoromethane (CF	3.313	3.019	8.9	89	0.00
3 T	Chloromethane	0.949	0.890	6.2	95	0.00
4 T	Vinyl Chloride	2.615	2.487	4.9	92	0.00
5 T	Bromomethane	1.407	1.417	-0.7	102	0.00
6 T	Chloroethane	1.340	1.295	3.4	97	0.00
7 T	Acetone	1.208	1.294	-7.1	103	0.00
8 T	Trichlorofluoromethane	2.761	2.600	5.8	93	0.00
9 T	1,1-Dichloroethene	1.376	1.351	1.8	99	0.00
10 T	Methylene Chloride	1.633	1.511	7.5	95	0.00
11 T	Trichlorotrifluoroethane	1.220	1.203	1.4	99	0.00
12 T	trans-1,2-Dichloroethene	1.588	1.483	6.6	95	0.00
13 T	1,1-Dichloroethane	2.980	2.876	3.5	95	0.00
14 T	Methyl tert-Butyl Ether	4.613	4.232	8.3	94	0.00
15 T	cis-1,2-Dichloroethene	1.591	1.557	2.1	99	0.00
16 T	Chloroform	2.636	2.617	0.7	98	0.00
17 S	1,2-Dichloroethane-d4 (SS1)	1.983	1.977	0.3	96	0.00
18 T	1,2-Dichloroethane	2.293	2.073	9.6	90	0.00
19 T	1,1,1-Trichloroethane	2.291	2.199	4.0	97	0.00
20 T	Benzene	6.222	6.262	-0.6	100	0.00
21 T	Carbon Tetrachloride	1.790	1.688	5.7	97	0.00
22 I	1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	106	0.00
23 T	1,2-Dichloropropane	0.366	0.322	12.0	96	0.00
24 T	Bromodichloromethane	0.439	0.383	12.8	96	0.00
25 T	Trichloroethene	0.343	0.303	11.7	99	0.00
26 T	1,4-Dioxane	0.263	0.233	11.4	102	0.00
27 T	cis-1,3-Dichloropropene	0.521	0.474	9.0	100	0.00
28 T	trans-1,3-Dichloropropene	0.444	0.403	9.2	101	0.00
29 T	1,1,2-Trichloroethane	0.276	0.252	8.7	101	0.00
30 S	Toluene-d8 (SS2)	1.056	1.191	-12.8	120	0.00
31 T	Toluene	1.374	1.230	10.5	102	0.00
32 T	1,2-Dibromoethane	0.340	0.316	7.1	103	0.00
33 T	Tetrachloroethene	0.359	0.336	6.4	104	0.00
34 I	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	105	0.00
35 T	Chlorobenzene	3.844	3.647	5.1	104	0.00
36 T	Ethylbenzene	6.608	5.864	11.3	98	0.00
37 T	m,p-Xylene	5.125	4.630	9.7	99	0.00

*Handwritten signature/initials*

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_03\22\  
 Data File : 03221204.D  
 Acq On : 22 Mar 2012 5:21 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
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 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
38 T	o-Xylene	5.490	5.028	8.4	102	0.00
39 T	1,1,2,2-Tetrachloroethane	2.561	2.469	3.6	107	0.00
40 S	Bromofluorobenzene (SS3)	1.675	1.947	-16.2	122	0.00
41 T	1,3-Dichlorobenzene	3.015	3.092	-2.6	115	0.00
42 T	1,4-Dichlorobenzene	3.034	3.052	-0.6	115	0.00
43 T	1,2-Dichlorobenzene	2.886	2.957	-2.5	117	0.00
44 T	1,2,4-Trichlorobenzene	1.907	2.112	-10.7	135	0.00
45 T	Naphthalene	6.232	7.427	-19.2	153	0.00
46 T	Hexachlorobutadiene	1.216	1.227	-0.9	122	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*KR 3/22/12*

## LABORATORY REPORT

March 30, 2012

Christoher Cote  
Montana DEQ  
1100 N. Last Chance Gulch  
Helena, MT 59601

### **RE: Background Indoor Air Study**

Dear Christoher:

Enclosed are the results of the samples submitted to our laboratory on March 16, 2012. For your reference, these analyses have been assigned our service request number P1201043.

All analyses were performed according to our laboratory's NELAP and DoD-ELAP-approved quality assurance program. The test results meet requirements of the current NELAP and DoD-ELAP standards, where applicable, and except as noted in the laboratory case narrative provided. For a specific list of NELAP and DoD-ELAP-accredited analytes, refer to the certifications section at [www.caslab.com](http://www.caslab.com). Results are intended to be considered in their entirety and apply only to the samples analyzed and reported herein.

Columbia Analytical Services, Inc. is certified by the California Department of Health Services, NELAP Laboratory Certificate No. 02115CA; Arizona Department of Health Services, Certificate No. AZ0694; Florida Department of Health, NELAP Certification E871020; New Jersey Department of Environmental Protection, NELAP Laboratory Certification ID #CA009; New York State Department of Health, NELAP NY Lab ID No: 11221; Oregon Environmental Laboratory Accreditation Program, NELAP ID: CA20007; The American Industrial Hygiene Association, Laboratory #101661; United States Department of Defense Environmental Laboratory Accreditation Program (DoD-ELAP), Certificate No. L11-203; Pennsylvania Registration No. 68-03307; TX Commission of Environmental Quality, NELAP ID T104704413-11-2; Minnesota Department of Health, NELAP Certificate No. 362188; Washington State Department of Ecology, ELAP Lab ID: C946, State of Utah Department of Health, NELAP Certificate No. CA015272011-1; Los Angeles Department of Building and Safety, Approval No: TA00001. Each of the certifications listed above have an explicit Scope of Accreditation that applies to specific matrices/methods/analytes; therefore, please contact me for information corresponding to a particular certification.

If you have any questions, please call me at (805) 526-7161.

Respectfully submitted,

**Columbia Analytical Services, Inc.**

Kate Aguilera  
Project Manager

Client: Montana DEQ  
Project: Background Indoor Air Study

CAS Project No: P1201043

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## CASE NARRATIVE

The samples were received intact under chain of custody on March 16, 2012 and were stored in accordance with the analytical method requirements. Please refer to the sample acceptance check form for additional information. The results reported herein are applicable only to the condition of the samples at the time of sample receipt.

### Air-Phase Petroleum Hydrocarbons (APH) Analysis

The samples were analyzed for total aliphatic and aromatic gasoline range hydrocarbons by gas chromatography/mass spectrometry according to the Method for the Determination of Air-Phase Petroleum Hydrocarbons (APH), Massachusetts Department of Environmental Protection, Revision 1, December, 2009.

### Volatile Organic Compound Analysis

The samples were also analyzed in scan and SIM mode for selected volatile organic compounds in accordance with EPA Method TO-15 from the Compendium of Methods for the Determination of Toxic Organic Compounds in Ambient Air, Second Edition (EPA/625/R-96/010b), January, 1999. The analytical system was comprised of a gas chromatograph / mass spectrometer (GC/MS) interfaced to a whole-air preconcentrator.

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*The results of analyses are given in the attached laboratory report. All results are intended to be considered in their entirety, and Columbia Analytical Services, Inc. (CAS) is not responsible for utilization of less than the complete report.*

*Use of Columbia Analytical Services, Inc. (CAS) Name. Client shall not use CAS's name or trademark in any marketing or reporting materials, press releases or in any other manner ("Materials") whatsoever and shall not attribute to CAS any test result, tolerance or specification derived from CAS's data ("Attribution") without CAS's prior written consent, which may be withheld by CAS for any reason in its sole discretion. To request CAS's consent, Client shall provide copies of the proposed Materials or Attribution and describe in writing Client's proposed use of such Materials or Attribution. If CAS has not provided written approval of the Materials or Attribution within ten (10) days of receipt from Client, Client's request to use CAS's name or trademark in any Materials or Attribution shall be deemed denied. CAS may, in its discretion, reasonably charge Client for its time in reviewing Materials or Attribution requests. Client acknowledges and agrees that the unauthorized use of CAS's name or trademark may cause CAS to incur irreparable harm for which the recovery of money damages will be inadequate. Accordingly, Client acknowledges and agrees that a violation shall justify preliminary injunctive relief. For questions contact the laboratory.*

DETAIL SUMMARY REPORT

Client: Montana DEQ  
 Project ID: Background Indoor Air Study

Service Request: P1201043

Date Received: 3/16/2012  
 Time Received: 10:15

Client Sample ID	Lab Code	Matrix	Date Collected	Time Collected	Container ID	Pi1 (psig)	Pf1 (psig)	MA APH 1.0 - MA VOC PH Can	TO-15 - VOC Cans	TO-15 - VOC SIM
15-IF-031212	P1201043-001	Air	3/13/2012	14:10	AS00187	-2.79	3.54	X	X	X
52-IF-031212	P1201043-002	Air	3/13/2012	14:10	AC01845	-4.19	3.67	X	X	X
10-IF-031212	P1201043-003	Air	3/13/2012	15:55	AC01842	-1.14	3.52	X	X	X
11-IF-031212	P1201043-004	Air	3/13/2012	16:30	AS00188	-2.57	3.67	X	X	X
12-IF-031312	P1201043-005	Air	3/14/2012	17:10	AS00185	-3.74	3.69	X	X	X
13-IF-031312	P1201043-006	Air	3/14/2012	09:40	AS00190	-3.01	3.68	X	X	X
16-IF-031312	P1201043-007	Air	3/14/2012	12:00	AC01428	-3.45	3.74	X	X	X



**Air - Chain of Custody Record & Analytical Service Request**

**Requested Turnaround Time in Business Days (Surcharges) please circle**  
 1 Day (100%) 2 Day (75%) 3 Day (50%) 4 Day (35%) 5 Day (25%) 10 Day-Standard

CAS Project No. **R1201043**

**Company Name & Address (Reporting Information)**  
 Monitoring Department of Environmental Quality  
 PO Box 200901  
 Helena, MT 59620-0901

**Project Name**  
 Background Indoor Air Study

**Project Number**

**Analysis Method**  
 MA-APH

**Comments**  
 e.g. Actual Preservative or specific instructions

**Project Manager**  
 Chris Cote

**Phone**  
 406-841-5078

**Fax**  
 406-841-5050

**P.O. # / Billing Information**  
 Reference 45; Background Indoor Air Study  
 Bill to Christine at MDEQ

**CAS Contact:**

**Email Address for Result Reporting**  
 cote@2emtgvy.mt.gov; Markishy@edm.com

**Sampler (Print & Sign)**  
 Bob Roll + Pat Skibsted

Client Sample ID	Laboratory ID Number	Date Collected	Time Collected	Canister ID (Bar code # - AC, SC, etc.)	Flow Controller ID (Bar code # - FC #)	Canister Start Pressure "Hg	Canister End Pressure "Hg/psig	Sample Volume	Analysis Method	Comments
15-1E-031212	0262	3/13/12	14:10	EA00187	EA00419	-24	-6.5		TO-15	Analyze project
52-1E-031212	0403	3/13/12	14:10	EA01945	EA00317	-27	-7.0		TO-15 SIM	Specific list of compounds
10-1E-031212	0005	3/13/12	15:55	EA01942	EA00413	-28	-1.5		X	
11-1E-031212	0240	3/13/12	16:30	EA00188	EA00017	-28	-4.5		X	
12-1E-031312	0356	3/14/12	17:10	EA00185	EA00387	-27	-6.5		X	
13-1E-031312	0284	3/14/12	9:40	EA00190	EA00019	-28	-5.0		X	to project specific
16-1E-031312	0325	3/14/12	12:00	EA01428	EA00374	-27	-5.5		X	reparing / mtds

**Report Tier Levels - please select**

Tier I - Results (Default if not specified) \_\_\_\_\_  
 Tier II (Results + QC Summaries) \_\_\_\_\_  
 Tier III (Results + QC & Calibration Summaries)   
 Tier IV (Data Validation Package) 10% Surcharge \_\_\_\_\_

**Relinquished by: (Signature)** *Pat Roll* **Date:** 3/15/12 **Time:** 1100

**Received by: (Signature)** *[Signature]* **Date:** 3/15/12 **Time:** 1015

**Project Requirements (MRLs, QAPP)**  
 EDD required  / No   
 Type: *SV and EXHA Exhaust*

**Cooler / Blank Temperature** \_\_\_\_\_ °C

**Sample Acceptance Check Form**

Client: Montana DEQ Work order: P1201043  
 Project: Background Indoor Air Study  
 Sample(s) received on: 3/16/12 Date opened: 3/16/12 by: MZAMORA

**Note:** This form is used for all samples received by CAS. The use of this form for custody seals is strictly meant to indicate presence/absence and not as an indication of compliance or nonconformity. Thermal preservation and pH will only be evaluated either at the request of the client and/or as required by the method/SOP.

- |  | <u>Yes</u>                          | <u>No</u>                           | <u>N/A</u>                          |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| 1 Were <b>sample containers</b> properly marked with client sample ID?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 2 Container(s) <b>supplied by CAS</b> ?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 3 Did <b>sample containers</b> arrive in good condition?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 4 Were <b>chain-of-custody</b> papers used and filled out?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 5 Did <b>sample container labels</b> and/or tags agree with custody papers?                                      | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 6 Was <b>sample volume</b> received adequate for analysis?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 7 Are samples within specified holding times?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 8 Was proper <b>temperature</b> (thermal preservation) of cooler at receipt adhered to?                          | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 9 Was a <b>trip blank</b> received?  | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| 10 Were <b>custody seals</b> on outside of cooler/Box?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| Location of seal(s)? _____ Sealing Lid?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were signature and date included?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were seals intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were custody seals on outside of sample container?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| Location of seal(s)? _____ Sealing Lid?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were signature and date included?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were seals intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 11 Do containers have appropriate <b>preservation</b> , according to method/SOP or Client specified information? | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Is there a client indication that the submitted samples are <b>pH</b> preserved?                                 | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were <b>VOA vials</b> checked for presence/absence of air bubbles?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Does the client/method/SOP require that the analyst check the sample pH and <u>if necessary</u> alter it?        | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 12 <b>Tubes:</b> Are the tubes capped and intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Do they contain moisture?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 13 <b>Badges:</b> Are the badges properly capped and intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Are dual bed badges separated and individually capped and intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |

Lab Sample ID	Container Description	Required pH *	Received pH	Adjusted pH	VOA Headspace (Presence/Absence)	Receipt / Preservation Comments
P1201043-001.01	6.0 L Silonite Can					
P1201043-002.01	6.0 L Ambient Can					
P1201043-003.01	6.0 L Ambient Can					
P1201043-004.01	6.0 L Silonite Can					
P1201043-005.01	6.0 L Silonite Can					
P1201043-006.01	6.0 L Silonite Can					
P1201043-007.01	6.0 L Ambient Can					

Explain any discrepancies: (include lab sample ID numbers): \_\_\_\_\_

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 15-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-001

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00187

Date Collected: 3/13/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.79      Final Pressure (psig): 3.54

Canister Dilution Factor: 1.53

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	46	31	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	15	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	7.7	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 52-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-002

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01845

Date Collected: 3/13/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.19      Final Pressure (psig): 3.67

Canister Dilution Factor: 1.75

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	41	35	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	19	18	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	8.8	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 10-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-003

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01842

Date Collected: 3/13/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.14      Final Pressure (psig): 3.52

Canister Dilution Factor: 1.34

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	32	27	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	100	13	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	6.7	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 11-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-004

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Lusine Hakobyan/Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00188

Date Collected: 3/13/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.57      Final Pressure (psig): 3.67

Canister Dilution Factor: 1.51

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	43	30	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	15	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	8.6	7.6	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 12-IF-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-005

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00185

Date Collected: 3/14/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.74 Final Pressure (psig): 3.69

Canister Dilution Factor: 1.68

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	60	34	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	17	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	8.4	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 13-IF-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-006

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00190

Date Collected: 3/14/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.01      Final Pressure (psig): 3.68

Canister Dilution Factor: 1.57

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	33	31	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	16	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	7.9	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 16-IF-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-007

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Lusine Hakobyan/Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01428

Date Collected: 3/14/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.45      Final Pressure (psig): 3.74

Canister Dilution Factor: 1.64

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	210	33	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	16	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	11	8.2	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P120324-MB

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	20	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	10	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	5.0	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

LABORATORY CONTROL SAMPLE SUMMARY

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**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P120324-LCS

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: NA Liter(s)

Compound	Spike Amount ng	Result ng	% Recovery	CAS	Data Qualifier
				Acceptance Limits	
C5 - C8 Aliphatic Hydrocarbons	24.5	22.5	92	70-130	
C9 - C12 Aliphatic Hydrocarbons	25.0	23.3	93	70-130	
C9 - C10 Aromatic Hydrocarbons	48.5	43.0	89	70-130	

LABORATORY DUPLICATE SUMMARY RESULTS

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**Client:** Montana DEQ

**Client Sample ID:** 52-IF-031212

**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043

CAS Sample ID: P1201043-002DUP

Test Code: Massachusetts APH, Revision 1, December 2009

Date Collected: 3/13/12

Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8

Date Received: 3/16/12

Analyst: Lusine Hakobyan

Date Analyzed: 3/24/12

Sampling Media: 6.0 L Summa Canister

Volume(s) Analyzed: 1.00 Liter(s)

Test Notes:

Container ID: AC01845

Initial Pressure (psig): -4.19 Final Pressure (psig): 3.67

Canister Dilution Factor: 1.75

Compound	Sample Result	Duplicate Sample Result	Average	% RPD	RPD	Data
	µg/m <sup>3</sup>	µg/m <sup>3</sup>	µg/m <sup>3</sup>		Limit	Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	41.3	48.0	44.65	<b>15</b>	30	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	18.8	24.6	21.7	<b>27</b>	30	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	ND	-	-	30	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 15-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-001

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00187

Date Collected: 3/13/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.79      Final Pressure (psig): 3.54

Canister Dilution Factor: 1.53

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	3.7	0.77	2.2	0.44	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.0	0.77	0.41	0.15	
74-87-3	Chloromethane	ND	0.77	ND	0.37	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.77	ND	0.11	
75-01-4	Vinyl Chloride	ND	0.77	ND	0.30	
106-99-0	1,3-Butadiene	ND	0.77	ND	0.35	
74-83-9	Bromomethane	ND	0.77	ND	0.20	
75-00-3	Chloroethane	ND	0.77	ND	0.29	
64-17-5	Ethanol	480	7.7	260	4.1	
75-05-8	Acetonitrile	ND	0.77	ND	0.46	
107-02-8	Acrolein	ND	3.1	ND	1.3	
67-64-1	Acetone	32	7.7	14	3.2	
75-69-4	Trichlorofluoromethane	2.6	0.77	0.46	0.14	
67-63-0	2-Propanol (Isopropyl Alcohol)	17	1.5	6.8	0.62	
107-13-1	Acrylonitrile	ND	0.77	ND	0.35	
75-35-4	1,1-Dichloroethene	ND	0.77	ND	0.19	
75-09-2	Methylene Chloride	ND	0.77	ND	0.22	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.77	ND	0.24	
76-13-1	Trichlorotrifluoroethane	ND	0.77	ND	0.10	
75-15-0	Carbon Disulfide	ND	7.7	ND	2.5	
156-60-5	trans-1,2-Dichloroethene	ND	0.77	ND	0.19	
75-34-3	1,1-Dichloroethane	ND	0.77	ND	0.19	
1634-04-4	Methyl tert-Butyl Ether	ND	0.77	ND	0.21	
108-05-4	Vinyl Acetate	ND	7.7	ND	2.2	
78-93-3	2-Butanone (MEK)	ND	7.7	ND	2.6	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 15-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
CAS Sample ID: P1201043-001

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00187

Date Collected: 3/13/12  
Date Received: 3/16/12  
Date Analyzed: 3/24/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.79 Final Pressure (psig): 3.54

Canister Dilution Factor: 1.53

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
156-59-2	cis-1,2-Dichloroethene	ND	0.77	ND	0.19	
141-78-6	Ethyl Acetate	9.5	1.5	2.6	0.42	
110-54-3	n-Hexane	ND	0.77	ND	0.22	
67-66-3	Chloroform	ND	0.77	ND	0.16	
109-99-9	Tetrahydrofuran (THF)	ND	0.77	ND	0.26	
107-06-2	1,2-Dichloroethane	ND	0.77	ND	0.19	
71-55-6	1,1,1-Trichloroethane	1.4	0.77	0.25	0.14	
71-43-2	Benzene	0.78	0.77	0.25	0.24	
56-23-5	Carbon Tetrachloride	ND	0.77	ND	0.12	
110-82-7	Cyclohexane	ND	1.5	ND	0.44	
78-87-5	1,2-Dichloropropane	ND	0.77	ND	0.17	
75-27-4	Bromodichloromethane	ND	0.77	ND	0.11	
79-01-6	Trichloroethene	ND	0.77	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.77	ND	0.21	
80-62-6	Methyl Methacrylate	ND	1.5	ND	0.37	
142-82-5	n-Heptane	ND	0.77	ND	0.19	
10061-01-5	cis-1,3-Dichloropropene	ND	0.77	ND	0.17	
108-10-1	4-Methyl-2-pentanone	ND	0.77	ND	0.19	
10061-02-6	trans-1,3-Dichloropropene	ND	0.77	ND	0.17	
79-00-5	1,1,2-Trichloroethane	ND	0.77	ND	0.14	
108-88-3	Toluene	9.8	0.77	2.6	0.20	
591-78-6	2-Hexanone	ND	0.77	ND	0.19	
124-48-1	Dibromochloromethane	ND	0.77	ND	0.090	
106-93-4	1,2-Dibromoethane	ND	0.77	ND	0.10	
123-86-4	n-Butyl Acetate	2.0	0.77	0.42	0.16	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 15-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-001

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AS00187

**Date Collected:** 3/13/12  
**Date Received:** 3/16/12  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -2.79      Final Pressure (psig): 3.54

Canister Dilution Factor: 1.53

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
111-65-9	n-Octane	ND	0.77	ND	0.16	
127-18-4	Tetrachloroethene	ND	0.77	ND	0.11	
108-90-7	Chlorobenzene	ND	0.77	ND	0.17	
100-41-4	Ethylbenzene	ND	0.77	ND	0.18	
179601-23-1	m,p-Xylenes	<b>1.6</b>	1.5	<b>0.37</b>	0.35	
75-25-2	Bromoform	ND	0.77	ND	0.074	
100-42-5	Styrene	ND	0.77	ND	0.18	
95-47-6	o-Xylene	ND	0.77	ND	0.18	
111-84-2	n-Nonane	ND	0.77	ND	0.15	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.77	ND	0.11	
98-82-8	Cumene	ND	0.77	ND	0.16	
80-56-8	alpha-Pinene	<b>1.5</b>	0.77	<b>0.27</b>	0.14	
103-65-1	n-Propylbenzene	ND	0.77	ND	0.16	
622-96-8	4-Ethyltoluene	ND	0.77	ND	0.16	
108-67-8	1,3,5-Trimethylbenzene	ND	0.77	ND	0.16	
95-63-6	1,2,4-Trimethylbenzene	ND	0.77	ND	0.16	
100-44-7	Benzyl Chloride	ND	0.77	ND	0.15	
541-73-1	1,3-Dichlorobenzene	ND	0.77	ND	0.13	
106-46-7	1,4-Dichlorobenzene	ND	0.77	ND	0.13	
95-50-1	1,2-Dichlorobenzene	ND	0.77	ND	0.13	
5989-27-5	d-Limonene	<b>16</b>	0.77	<b>2.9</b>	0.14	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.77	ND	0.079	
120-82-1	1,2,4-Trichlorobenzene	ND	0.77	ND	0.10	
91-20-3	Naphthalene	ND	0.77	ND	0.15	
87-68-3	Hexachlorobutadiene	ND	0.77	ND	0.072	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 52-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
CAS Sample ID: P1201043-002

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01845

Date Collected: 3/13/12  
Date Received: 3/16/12  
Date Analyzed: 3/24/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.19 Final Pressure (psig): 3.67

Canister Dilution Factor: 1.75

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	3.4	0.88	2.0	0.51	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.0	0.88	0.40	0.18	
74-87-3	Chloromethane	ND	0.88	ND	0.42	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.88	ND	0.13	
75-01-4	Vinyl Chloride	ND	0.88	ND	0.34	
106-99-0	1,3-Butadiene	ND	0.88	ND	0.40	
74-83-9	Bromomethane	ND	0.88	ND	0.23	
75-00-3	Chloroethane	ND	0.88	ND	0.33	
64-17-5	Ethanol	480	8.8	250	4.6	
75-05-8	Acetonitrile	ND	0.88	ND	0.52	
107-02-8	Acrolein	ND	3.5	ND	1.5	
67-64-1	Acetone	31	8.8	13	3.7	
75-69-4	Trichlorofluoromethane	2.6	0.88	0.45	0.16	
67-63-0	2-Propanol (Isopropyl Alcohol)	15	1.8	6.2	0.71	
107-13-1	Acrylonitrile	ND	0.88	ND	0.40	
75-35-4	1,1-Dichloroethene	ND	0.88	ND	0.22	
75-09-2	Methylene Chloride	ND	0.88	ND	0.25	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.88	ND	0.28	
76-13-1	Trichlorotrifluoroethane	ND	0.88	ND	0.11	
75-15-0	Carbon Disulfide	ND	8.8	ND	2.8	
156-60-5	trans-1,2-Dichloroethene	ND	0.88	ND	0.22	
75-34-3	1,1-Dichloroethane	ND	0.88	ND	0.22	
1634-04-4	Methyl tert-Butyl Ether	ND	0.88	ND	0.24	
108-05-4	Vinyl Acetate	ND	8.8	ND	2.5	
78-93-3	2-Butanone (MEK)	ND	8.8	ND	3.0	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 52-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-002

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01845

Date Collected: 3/13/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.19      Final Pressure (psig): 3.67

Canister Dilution Factor: 1.75

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
156-59-2	cis-1,2-Dichloroethene	ND	0.88	ND	0.22	
141-78-6	Ethyl Acetate	9.7	1.8	2.7	0.49	
110-54-3	n-Hexane	ND	0.88	ND	0.25	
67-66-3	Chloroform	ND	0.88	ND	0.18	
109-99-9	Tetrahydrofuran (THF)	ND	0.88	ND	0.30	
107-06-2	1,2-Dichloroethane	ND	0.88	ND	0.22	
71-55-6	1,1,1-Trichloroethane	1.4	0.88	0.26	0.16	
71-43-2	Benzene	ND	0.88	ND	0.27	
56-23-5	Carbon Tetrachloride	ND	0.88	ND	0.14	
110-82-7	Cyclohexane	ND	1.8	ND	0.51	
78-87-5	1,2-Dichloropropane	ND	0.88	ND	0.19	
75-27-4	Bromodichloromethane	ND	0.88	ND	0.13	
79-01-6	Trichloroethene	ND	0.88	ND	0.16	
123-91-1	1,4-Dioxane	ND	0.88	ND	0.24	
80-62-6	Methyl Methacrylate	ND	1.8	ND	0.43	
142-82-5	n-Heptane	ND	0.88	ND	0.21	
10061-01-5	cis-1,3-Dichloropropene	ND	0.88	ND	0.19	
108-10-1	4-Methyl-2-pentanone	ND	0.88	ND	0.21	
10061-02-6	trans-1,3-Dichloropropene	ND	0.88	ND	0.19	
79-00-5	1,1,2-Trichloroethane	ND	0.88	ND	0.16	
108-88-3	Toluene	10	0.88	2.6	0.23	
591-78-6	2-Hexanone	ND	0.88	ND	0.21	
124-48-1	Dibromochloromethane	ND	0.88	ND	0.10	
106-93-4	1,2-Dibromoethane	ND	0.88	ND	0.11	
123-86-4	n-Butyl Acetate	2.1	0.88	0.44	0.18	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 52-IF-031212  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201043  
**CAS Sample ID:** P1201043-002

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01845

**Date Collected:** 3/13/12  
**Date Received:** 3/16/12  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -4.19      Final Pressure (psig): 3.67

Canister Dilution Factor: 1.75

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
111-65-9	n-Octane	ND	0.88	ND	0.19	
127-18-4	Tetrachloroethene	ND	0.88	ND	0.13	
108-90-7	Chlorobenzene	ND	0.88	ND	0.19	
100-41-4	Ethylbenzene	ND	0.88	ND	0.20	
179601-23-1	m,p-Xylenes	ND	1.8	ND	0.40	
75-25-2	Bromoform	ND	0.88	ND	0.085	
100-42-5	Styrene	ND	0.88	ND	0.21	
95-47-6	o-Xylene	ND	0.88	ND	0.20	
111-84-2	n-Nonane	ND	0.88	ND	0.17	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.88	ND	0.13	
98-82-8	Cumene	ND	0.88	ND	0.18	
80-56-8	alpha-Pinene	<b>1.6</b>	0.88	<b>0.28</b>	0.16	
103-65-1	n-Propylbenzene	ND	0.88	ND	0.18	
622-96-8	4-Ethyltoluene	ND	0.88	ND	0.18	
108-67-8	1,3,5-Trimethylbenzene	ND	0.88	ND	0.18	
95-63-6	1,2,4-Trimethylbenzene	ND	0.88	ND	0.18	
100-44-7	Benzyl Chloride	ND	0.88	ND	0.17	
541-73-1	1,3-Dichlorobenzene	ND	0.88	ND	0.15	
106-46-7	1,4-Dichlorobenzene	ND	0.88	ND	0.15	
95-50-1	1,2-Dichlorobenzene	ND	0.88	ND	0.15	
5989-27-5	d-Limonene	<b>18</b>	0.88	<b>3.2</b>	0.16	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.88	ND	0.091	
120-82-1	1,2,4-Trichlorobenzene	ND	0.88	ND	0.12	
91-20-3	Naphthalene	ND	0.88	ND	0.17	
87-68-3	Hexachlorobutadiene	ND	0.88	ND	0.082	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 10-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-003

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01842

**Date Collected:** 3/13/12  
**Date Received:** 3/16/12  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -1.14      Final Pressure (psig): 3.52

Canister Dilution Factor: 1.34

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	6.7	0.67	3.9	0.39	
75-71-8	Dichlorodifluoromethane (CFC 12)	1.9	0.67	0.39	0.14	
74-87-3	Chloromethane	ND	0.67	ND	0.32	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.67	ND	0.096	
75-01-4	Vinyl Chloride	ND	0.67	ND	0.26	
106-99-0	1,3-Butadiene	ND	0.67	ND	0.30	
74-83-9	Bromomethane	ND	0.67	ND	0.17	
75-00-3	Chloroethane	ND	0.67	ND	0.25	
64-17-5	Ethanol	190	6.7	99	3.6	
75-05-8	Acetonitrile	ND	0.67	ND	0.40	
107-02-8	Acrolein	ND	2.7	ND	1.2	
67-64-1	Acetone	25	6.7	11	2.8	
75-69-4	Trichlorofluoromethane	1.0	0.67	0.19	0.12	
67-63-0	2-Propanol (Isopropyl Alcohol)	36	1.3	15	0.55	
107-13-1	Acrylonitrile	ND	0.67	ND	0.31	
75-35-4	1,1-Dichloroethene	ND	0.67	ND	0.17	
75-09-2	Methylene Chloride	ND	0.67	ND	0.19	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.67	ND	0.21	
76-13-1	Trichlorotrifluoroethane	ND	0.67	ND	0.087	
75-15-0	Carbon Disulfide	ND	6.7	ND	2.2	
156-60-5	trans-1,2-Dichloroethene	ND	0.67	ND	0.17	
75-34-3	1,1-Dichloroethane	ND	0.67	ND	0.17	
1634-04-4	Methyl tert-Butyl Ether	ND	0.67	ND	0.19	
108-05-4	Vinyl Acetate	ND	6.7	ND	1.9	
78-93-3	2-Butanone (MEK)	ND	6.7	ND	2.3	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 10-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
CAS Sample ID: P1201043-003

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01842

Date Collected: 3/13/12  
Date Received: 3/16/12  
Date Analyzed: 3/24/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.14 Final Pressure (psig): 3.52

Canister Dilution Factor: 1.34

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
156-59-2	cis-1,2-Dichloroethene	ND	0.67	ND	0.17	
141-78-6	Ethyl Acetate	3.7	1.3	1.0	0.37	
110-54-3	n-Hexane	1.0	0.67	0.29	0.19	
67-66-3	Chloroform	ND	0.67	ND	0.14	
109-99-9	Tetrahydrofuran (THF)	ND	0.67	ND	0.23	
107-06-2	1,2-Dichloroethane	ND	0.67	ND	0.17	
71-55-6	1,1,1-Trichloroethane	ND	0.67	ND	0.12	
71-43-2	Benzene	0.81	0.67	0.25	0.21	
56-23-5	Carbon Tetrachloride	ND	0.67	ND	0.11	
110-82-7	Cyclohexane	ND	1.3	ND	0.39	
78-87-5	1,2-Dichloropropane	ND	0.67	ND	0.15	
75-27-4	Bromodichloromethane	ND	0.67	ND	0.10	
79-01-6	Trichloroethene	ND	0.67	ND	0.12	
123-91-1	1,4-Dioxane	ND	0.67	ND	0.19	
80-62-6	Methyl Methacrylate	ND	1.3	ND	0.33	
142-82-5	n-Heptane	0.79	0.67	0.19	0.16	
10061-01-5	cis-1,3-Dichloropropene	ND	0.67	ND	0.15	
108-10-1	4-Methyl-2-pentanone	ND	0.67	ND	0.16	
10061-02-6	trans-1,3-Dichloropropene	ND	0.67	ND	0.15	
79-00-5	1,1,2-Trichloroethane	ND	0.67	ND	0.12	
108-88-3	Toluene	6.6	0.67	1.8	0.18	
591-78-6	2-Hexanone	ND	0.67	ND	0.16	
124-48-1	Dibromochloromethane	ND	0.67	ND	0.079	
106-93-4	1,2-Dibromoethane	ND	0.67	ND	0.087	
123-86-4	n-Butyl Acetate	0.84	0.67	0.18	0.14	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 10-IF-031212  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201043  
**CAS Sample ID:** P1201043-003

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01842

**Date Collected:** 3/13/12  
**Date Received:** 3/16/12  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -1.14      Final Pressure (psig): 3.52

Canister Dilution Factor: 1.34

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
111-65-9	n-Octane	ND	0.67	ND	0.14	
127-18-4	Tetrachloroethene	ND	0.67	ND	0.099	
108-90-7	Chlorobenzene	ND	0.67	ND	0.15	
100-41-4	Ethylbenzene	ND	0.67	ND	0.15	
179601-23-1	m,p-Xylenes	<b>1.5</b>	1.3	<b>0.34</b>	0.31	
75-25-2	Bromoform	ND	0.67	ND	0.065	
100-42-5	Styrene	ND	0.67	ND	0.16	
95-47-6	o-Xylene	ND	0.67	ND	0.15	
111-84-2	n-Nonane	ND	0.67	ND	0.13	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.67	ND	0.098	
98-82-8	Cumene	ND	0.67	ND	0.14	
80-56-8	alpha-Pinene	<b>1.5</b>	0.67	<b>0.26</b>	0.12	
103-65-1	n-Propylbenzene	ND	0.67	ND	0.14	
622-96-8	4-Ethyltoluene	ND	0.67	ND	0.14	
108-67-8	1,3,5-Trimethylbenzene	ND	0.67	ND	0.14	
95-63-6	1,2,4-Trimethylbenzene	ND	0.67	ND	0.14	
100-44-7	Benzyl Chloride	ND	0.67	ND	0.13	
541-73-1	1,3-Dichlorobenzene	ND	0.67	ND	0.11	
106-46-7	1,4-Dichlorobenzene	ND	0.67	ND	0.11	
95-50-1	1,2-Dichlorobenzene	ND	0.67	ND	0.11	
5989-27-5	d-Limonene	<b>10</b>	0.67	<b>1.8</b>	0.12	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.67	ND	0.069	
120-82-1	1,2,4-Trichlorobenzene	ND	0.67	ND	0.090	
91-20-3	Naphthalene	ND	0.67	ND	0.13	
87-68-3	Hexachlorobutadiene	ND	0.67	ND	0.063	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 11-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-004

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Lusine Hakobyan/Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AS00188

**Date Collected:** 3/13/12  
**Date Received:** 3/16/12  
**Date Analyzed:** 3/24/12 & 3/26/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.050 Liter(s)

Initial Pressure (psig): -2.57 Final Pressure (psig): 3.67

Canister Dilution Factor: 1.51

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	5.7	0.76	3.3	0.44	
75-71-8	Dichlorodifluoromethane (CFC 12)	1.9	0.76	0.39	0.15	
74-87-3	Chloromethane	ND	0.76	ND	0.37	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.76	ND	0.11	
75-01-4	Vinyl Chloride	ND	0.76	ND	0.30	
106-99-0	1,3-Butadiene	ND	0.76	ND	0.34	
74-83-9	Bromomethane	ND	0.76	ND	0.19	
75-00-3	Chloroethane	ND	0.76	ND	0.29	
64-17-5	Ethanol	5,900	150	3,100	80	D
75-05-8	Acetonitrile	ND	0.76	ND	0.45	
107-02-8	Acrolein	ND	3.0	ND	1.3	
67-64-1	Acetone	62	7.6	26	3.2	
75-69-4	Trichlorofluoromethane	1.1	0.76	0.19	0.13	
67-63-0	2-Propanol (Isopropyl Alcohol)	58	1.5	24	0.61	
107-13-1	Acrylonitrile	ND	0.76	ND	0.35	
75-35-4	1,1-Dichloroethene	ND	0.76	ND	0.19	
75-09-2	Methylene Chloride	ND	0.76	ND	0.22	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.76	ND	0.24	
76-13-1	Trichlorotrifluoroethane	ND	0.76	ND	0.099	
75-15-0	Carbon Disulfide	ND	7.6	ND	2.4	
156-60-5	trans-1,2-Dichloroethene	ND	0.76	ND	0.19	
75-34-3	1,1-Dichloroethane	ND	0.76	ND	0.19	
1634-04-4	Methyl tert-Butyl Ether	ND	0.76	ND	0.21	
108-05-4	Vinyl Acetate	ND	7.6	ND	2.1	
78-93-3	2-Butanone (MEK)	ND	7.6	ND	2.6	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 11-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
CAS Sample ID: P1201043-004

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Lusine Hakobyan/Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00188

Date Collected: 3/13/12  
Date Received: 3/16/12  
Date Analyzed: 3/24/12 & 3/26/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.050 Liter(s)

Initial Pressure (psig): -2.57 Final Pressure (psig): 3.67

Canister Dilution Factor: 1.51

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
156-59-2	cis-1,2-Dichloroethene	ND	0.76	ND	0.19	
141-78-6	Ethyl Acetate	16	1.5	4.4	0.42	
110-54-3	n-Hexane	0.80	0.76	0.23	0.21	
67-66-3	Chloroform	1.3	0.76	0.27	0.15	
109-99-9	Tetrahydrofuran (THF)	ND	0.76	ND	0.26	
107-06-2	1,2-Dichloroethane	ND	0.76	ND	0.19	
71-55-6	1,1,1-Trichloroethane	ND	0.76	ND	0.14	
71-43-2	Benzene	ND	0.76	ND	0.24	
56-23-5	Carbon Tetrachloride	ND	0.76	ND	0.12	
110-82-7	Cyclohexane	ND	1.5	ND	0.44	
78-87-5	1,2-Dichloropropane	ND	0.76	ND	0.16	
75-27-4	Bromodichloromethane	ND	0.76	ND	0.11	
79-01-6	Trichloroethene	ND	0.76	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.76	ND	0.21	
80-62-6	Methyl Methacrylate	ND	1.5	ND	0.37	
142-82-5	n-Heptane	ND	0.76	ND	0.18	
10061-01-5	cis-1,3-Dichloropropene	ND	0.76	ND	0.17	
108-10-1	4-Methyl-2-pentanone	ND	0.76	ND	0.18	
10061-02-6	trans-1,3-Dichloropropene	ND	0.76	ND	0.17	
79-00-5	1,1,2-Trichloroethane	ND	0.76	ND	0.14	
108-88-3	Toluene	1.5	0.76	0.41	0.20	
591-78-6	2-Hexanone	ND	0.76	ND	0.18	
124-48-1	Dibromochloromethane	ND	0.76	ND	0.089	
106-93-4	1,2-Dibromoethane	ND	0.76	ND	0.098	
123-86-4	n-Butyl Acetate	ND	0.76	ND	0.16	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 11-IF-031212  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201043  
**CAS Sample ID:** P1201043-004

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Lusine Hakobyan/Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AS00188

**Date Collected:** 3/13/12  
**Date Received:** 3/16/12  
**Date Analyzed:** 3/24/12 & 3/26/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.050 Liter(s)

Initial Pressure (psig): -2.57      Final Pressure (psig): 3.67

Canister Dilution Factor: 1.51

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
111-65-9	n-Octane	ND	0.76	ND	0.16	
127-18-4	Tetrachloroethene	ND	0.76	ND	0.11	
108-90-7	Chlorobenzene	ND	0.76	ND	0.16	
100-41-4	Ethylbenzene	ND	0.76	ND	0.17	
179601-23-1	m,p-Xylenes	ND	1.5	ND	0.35	
75-25-2	Bromoform	ND	0.76	ND	0.073	
100-42-5	Styrene	<b>0.87</b>	0.76	<b>0.21</b>	0.18	
95-47-6	o-Xylene	ND	0.76	ND	0.17	
111-84-2	n-Nonane	ND	0.76	ND	0.14	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.76	ND	0.11	
98-82-8	Cumene	ND	0.76	ND	0.15	
80-56-8	alpha-Pinene	<b>4.0</b>	0.76	<b>0.71</b>	0.14	
103-65-1	n-Propylbenzene	ND	0.76	ND	0.15	
622-96-8	4-Ethyltoluene	ND	0.76	ND	0.15	
108-67-8	1,3,5-Trimethylbenzene	<b>0.85</b>	0.76	<b>0.17</b>	0.15	
95-63-6	1,2,4-Trimethylbenzene	<b>1.5</b>	0.76	<b>0.31</b>	0.15	
100-44-7	Benzyl Chloride	ND	0.76	ND	0.15	
541-73-1	1,3-Dichlorobenzene	ND	0.76	ND	0.13	
106-46-7	1,4-Dichlorobenzene	ND	0.76	ND	0.13	
95-50-1	1,2-Dichlorobenzene	ND	0.76	ND	0.13	
5989-27-5	d-Limonene	<b>9.0</b>	0.76	<b>1.6</b>	0.14	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.76	ND	0.078	
120-82-1	1,2,4-Trichlorobenzene	ND	0.76	ND	0.10	
91-20-3	Naphthalene	ND	0.76	ND	0.14	
87-68-3	Hexachlorobutadiene	ND	0.76	ND	0.071	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 12-IF-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-005

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AS00185

**Date Collected:** 3/14/12  
**Date Received:** 3/16/12  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -3.74      Final Pressure (psig): 3.69

Canister Dilution Factor: 1.68

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	2.8	0.84	1.6	0.49	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.0	0.84	0.41	0.17	
74-87-3	Chloromethane	ND	0.84	ND	0.41	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.84	ND	0.12	
75-01-4	Vinyl Chloride	ND	0.84	ND	0.33	
106-99-0	1,3-Butadiene	ND	0.84	ND	0.38	
74-83-9	Bromomethane	ND	0.84	ND	0.22	
75-00-3	Chloroethane	ND	0.84	ND	0.32	
64-17-5	Ethanol	86	8.4	46	4.5	
75-05-8	Acetonitrile	ND	0.84	ND	0.50	
107-02-8	Acrolein	ND	3.4	ND	1.5	
67-64-1	Acetone	27	8.4	11	3.5	
75-69-4	Trichlorofluoromethane	3.9	0.84	0.69	0.15	
67-63-0	2-Propanol (Isopropyl Alcohol)	8.6	1.7	3.5	0.68	
107-13-1	Acrylonitrile	ND	0.84	ND	0.39	
75-35-4	1,1-Dichloroethene	ND	0.84	ND	0.21	
75-09-2	Methylene Chloride	ND	0.84	ND	0.24	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.84	ND	0.27	
76-13-1	Trichlorotrifluoroethane	ND	0.84	ND	0.11	
75-15-0	Carbon Disulfide	ND	8.4	ND	2.7	
156-60-5	trans-1,2-Dichloroethene	ND	0.84	ND	0.21	
75-34-3	1,1-Dichloroethane	ND	0.84	ND	0.21	
1634-04-4	Methyl tert-Butyl Ether	ND	0.84	ND	0.23	
108-05-4	Vinyl Acetate	ND	8.4	ND	2.4	
78-93-3	2-Butanone (MEK)	ND	8.4	ND	2.8	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 12-IF-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-005

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00185

Date Collected: 3/14/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.74 Final Pressure (psig): 3.69

Canister Dilution Factor: 1.68

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
156-59-2	cis-1,2-Dichloroethene	ND	0.84	ND	0.21	
141-78-6	Ethyl Acetate	4.1	1.7	1.1	0.47	
110-54-3	n-Hexane	1.9	0.84	0.53	0.24	
67-66-3	Chloroform	ND	0.84	ND	0.17	
109-99-9	Tetrahydrofuran (THF)	ND	0.84	ND	0.28	
107-06-2	1,2-Dichloroethane	2.4	0.84	0.59	0.21	
71-55-6	1,1,1-Trichloroethane	ND	0.84	ND	0.15	
71-43-2	Benzene	1.8	0.84	0.55	0.26	
56-23-5	Carbon Tetrachloride	ND	0.84	ND	0.13	
110-82-7	Cyclohexane	ND	1.7	ND	0.49	
78-87-5	1,2-Dichloropropane	ND	0.84	ND	0.18	
75-27-4	Bromodichloromethane	ND	0.84	ND	0.13	
79-01-6	Trichloroethene	ND	0.84	ND	0.16	
123-91-1	1,4-Dioxane	ND	0.84	ND	0.23	
80-62-6	Methyl Methacrylate	ND	1.7	ND	0.41	
142-82-5	n-Heptane	0.87	0.84	0.21	0.21	
10061-01-5	cis-1,3-Dichloropropene	ND	0.84	ND	0.19	
108-10-1	4-Methyl-2-pentanone	ND	0.84	ND	0.21	
10061-02-6	trans-1,3-Dichloropropene	ND	0.84	ND	0.19	
79-00-5	1,1,2-Trichloroethane	ND	0.84	ND	0.15	
108-88-3	Toluene	5.4	0.84	1.4	0.22	
591-78-6	2-Hexanone	ND	0.84	ND	0.21	
124-48-1	Dibromochloromethane	ND	0.84	ND	0.099	
106-93-4	1,2-Dibromoethane	ND	0.84	ND	0.11	
123-86-4	n-Butyl Acetate	ND	0.84	ND	0.18	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 12-IF-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
CAS Sample ID: P1201043-005

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00185

Date Collected: 3/14/12  
Date Received: 3/16/12  
Date Analyzed: 3/24/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.74 Final Pressure (psig): 3.69

Canister Dilution Factor: 1.68

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
111-65-9	n-Octane	ND	0.84	ND	0.18	
127-18-4	Tetrachloroethene	ND	0.84	ND	0.12	
108-90-7	Chlorobenzene	ND	0.84	ND	0.18	
100-41-4	Ethylbenzene	<b>0.85</b>	0.84	<b>0.20</b>	0.19	
179601-23-1	m,p-Xylenes	<b>2.9</b>	1.7	<b>0.66</b>	0.39	
75-25-2	Bromoform	ND	0.84	ND	0.081	
100-42-5	Styrene	ND	0.84	ND	0.20	
95-47-6	o-Xylene	<b>0.90</b>	0.84	<b>0.21</b>	0.19	
111-84-2	n-Nonane	ND	0.84	ND	0.16	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.84	ND	0.12	
98-82-8	Cumene	ND	0.84	ND	0.17	
80-56-8	alpha-Pinene	<b>1.1</b>	0.84	<b>0.19</b>	0.15	
103-65-1	n-Propylbenzene	ND	0.84	ND	0.17	
622-96-8	4-Ethyltoluene	ND	0.84	ND	0.17	
108-67-8	1,3,5-Trimethylbenzene	ND	0.84	ND	0.17	
95-63-6	1,2,4-Trimethylbenzene	<b>1.0</b>	0.84	<b>0.21</b>	0.17	
100-44-7	Benzyl Chloride	ND	0.84	ND	0.16	
541-73-1	1,3-Dichlorobenzene	ND	0.84	ND	0.14	
106-46-7	1,4-Dichlorobenzene	ND	0.84	ND	0.14	
95-50-1	1,2-Dichlorobenzene	ND	0.84	ND	0.14	
5989-27-5	d-Limonene	<b>5.8</b>	0.84	<b>1.0</b>	0.15	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.84	ND	0.087	
120-82-1	1,2,4-Trichlorobenzene	ND	0.84	ND	0.11	
91-20-3	Naphthalene	ND	0.84	ND	0.16	
87-68-3	Hexachlorobutadiene	ND	0.84	ND	0.079	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 13-IF-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-006

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AS00190

**Date Collected:** 3/14/12  
**Date Received:** 3/16/12  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -3.01      Final Pressure (psig): 3.68

Canister Dilution Factor: 1.57

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	6.9	0.79	4.0	0.46	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.0	0.79	0.40	0.16	
74-87-3	Chloromethane	ND	0.79	ND	0.38	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.79	ND	0.11	
75-01-4	Vinyl Chloride	ND	0.79	ND	0.31	
106-99-0	1,3-Butadiene	ND	0.79	ND	0.35	
74-83-9	Bromomethane	ND	0.79	ND	0.20	
75-00-3	Chloroethane	ND	0.79	ND	0.30	
64-17-5	Ethanol	570	7.9	310	4.2	
75-05-8	Acetonitrile	ND	0.79	ND	0.47	
107-02-8	Acrolein	4.8	3.1	2.1	1.4	
67-64-1	Acetone	35	7.9	15	3.3	
75-69-4	Trichlorofluoromethane	1.1	0.79	0.19	0.14	
67-63-0	2-Propanol (Isopropyl Alcohol)	40	1.6	16	0.64	
107-13-1	Acrylonitrile	ND	0.79	ND	0.36	
75-35-4	1,1-Dichloroethene	ND	0.79	ND	0.20	
75-09-2	Methylene Chloride	ND	0.79	ND	0.23	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.79	ND	0.25	
76-13-1	Trichlorotrifluoroethane	ND	0.79	ND	0.10	
75-15-0	Carbon Disulfide	ND	7.9	ND	2.5	
156-60-5	trans-1,2-Dichloroethene	ND	0.79	ND	0.20	
75-34-3	1,1-Dichloroethane	ND	0.79	ND	0.19	
1634-04-4	Methyl tert-Butyl Ether	ND	0.79	ND	0.22	
108-05-4	Vinyl Acetate	ND	7.9	ND	2.2	
78-93-3	2-Butanone (MEK)	ND	7.9	ND	2.7	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 13-IF-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-006

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00190

Date Collected: 3/14/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.01      Final Pressure (psig): 3.68

Canister Dilution Factor: 1.57

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
156-59-2	cis-1,2-Dichloroethene	ND	0.79	ND	0.20	
141-78-6	Ethyl Acetate	4.5	1.6	1.2	0.44	
110-54-3	n-Hexane	ND	0.79	ND	0.22	
67-66-3	Chloroform	ND	0.79	ND	0.16	
109-99-9	Tetrahydrofuran (THF)	ND	0.79	ND	0.27	
107-06-2	1,2-Dichloroethane	ND	0.79	ND	0.19	
71-55-6	1,1,1-Trichloroethane	ND	0.79	ND	0.14	
71-43-2	Benzene	ND	0.79	ND	0.25	
56-23-5	Carbon Tetrachloride	ND	0.79	ND	0.12	
110-82-7	Cyclohexane	ND	1.6	ND	0.46	
78-87-5	1,2-Dichloropropane	ND	0.79	ND	0.17	
75-27-4	Bromodichloromethane	ND	0.79	ND	0.12	
79-01-6	Trichloroethene	ND	0.79	ND	0.15	
123-91-1	1,4-Dioxane	ND	0.79	ND	0.22	
80-62-6	Methyl Methacrylate	ND	1.6	ND	0.38	
142-82-5	n-Heptane	1.1	0.79	0.28	0.19	
10061-01-5	cis-1,3-Dichloropropene	ND	0.79	ND	0.17	
108-10-1	4-Methyl-2-pentanone	ND	0.79	ND	0.19	
10061-02-6	trans-1,3-Dichloropropene	ND	0.79	ND	0.17	
79-00-5	1,1,2-Trichloroethane	ND	0.79	ND	0.14	
108-88-3	Toluene	8.4	0.79	2.2	0.21	
591-78-6	2-Hexanone	ND	0.79	ND	0.19	
124-48-1	Dibromochloromethane	ND	0.79	ND	0.092	
106-93-4	1,2-Dibromoethane	ND	0.79	ND	0.10	
123-86-4	n-Butyl Acetate	1.4	0.79	0.30	0.17	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 13-IF-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-006

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AS00190

**Date Collected:** 3/14/12  
**Date Received:** 3/16/12  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -3.01      Final Pressure (psig): 3.68

Canister Dilution Factor: 1.57

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
111-65-9	n-Octane	1.3	0.79	0.28	0.17	
127-18-4	Tetrachloroethene	ND	0.79	ND	0.12	
108-90-7	Chlorobenzene	ND	0.79	ND	0.17	
100-41-4	Ethylbenzene	ND	0.79	ND	0.18	
179601-23-1	m,p-Xylenes	ND	1.6	ND	0.36	
75-25-2	Bromoform	ND	0.79	ND	0.076	
100-42-5	Styrene	ND	0.79	ND	0.18	
95-47-6	o-Xylene	ND	0.79	ND	0.18	
111-84-2	n-Nonane	ND	0.79	ND	0.15	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.79	ND	0.11	
98-82-8	Cumene	ND	0.79	ND	0.16	
80-56-8	alpha-Pinene	4.0	0.79	0.72	0.14	
103-65-1	n-Propylbenzene	ND	0.79	ND	0.16	
622-96-8	4-Ethyltoluene	ND	0.79	ND	0.16	
108-67-8	1,3,5-Trimethylbenzene	ND	0.79	ND	0.16	
95-63-6	1,2,4-Trimethylbenzene	ND	0.79	ND	0.16	
100-44-7	Benzyl Chloride	ND	0.79	ND	0.15	
541-73-1	1,3-Dichlorobenzene	ND	0.79	ND	0.13	
106-46-7	1,4-Dichlorobenzene	ND	0.79	ND	0.13	
95-50-1	1,2-Dichlorobenzene	ND	0.79	ND	0.13	
5989-27-5	d-Limonene	50	0.79	8.9	0.14	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.79	ND	0.081	
120-82-1	1,2,4-Trichlorobenzene	ND	0.79	ND	0.11	
91-20-3	Naphthalene	ND	0.79	ND	0.15	
87-68-3	Hexachlorobutadiene	ND	0.79	ND	0.074	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 16-IF-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-007

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Lusine Hakobyan/Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01428

**Date Collected:** 3/14/12  
**Date Received:** 3/16/12  
**Date Analyzed:** 3/24/12 & 3/26/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.050 Liter(s)

Initial Pressure (psig): -3.45      Final Pressure (psig): 3.74

Canister Dilution Factor: 1.64

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	11	0.82	6.1	0.48	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.0	0.82	0.40	0.17	
74-87-3	Chloromethane	ND	0.82	ND	0.40	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.82	ND	0.12	
75-01-4	Vinyl Chloride	ND	0.82	ND	0.32	
106-99-0	1,3-Butadiene	ND	0.82	ND	0.37	
74-83-9	Bromomethane	ND	0.82	ND	0.21	
75-00-3	Chloroethane	ND	0.82	ND	0.31	
64-17-5	Ethanol	1,800	160	980	87	D
75-05-8	Acetonitrile	ND	0.82	ND	0.49	
107-02-8	Acrolein	3.5	3.3	1.5	1.4	
67-64-1	Acetone	80	8.2	34	3.5	
75-69-4	Trichlorofluoromethane	1.1	0.82	0.20	0.15	
67-63-0	2-Propanol (Isopropyl Alcohol)	30	1.6	12	0.67	
107-13-1	Acrylonitrile	ND	0.82	ND	0.38	
75-35-4	1,1-Dichloroethene	ND	0.82	ND	0.21	
75-09-2	Methylene Chloride	ND	0.82	ND	0.24	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.82	ND	0.26	
76-13-1	Trichlorotrifluoroethane	ND	0.82	ND	0.11	
75-15-0	Carbon Disulfide	ND	8.2	ND	2.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.82	ND	0.21	
75-34-3	1,1-Dichloroethane	ND	0.82	ND	0.20	
1634-04-4	Methyl tert-Butyl Ether	ND	0.82	ND	0.23	
108-05-4	Vinyl Acetate	ND	8.2	ND	2.3	
78-93-3	2-Butanone (MEK)	ND	8.2	ND	2.8	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 16-IF-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-007

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Lusine Hakobyan/Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01428

**Date Collected:** 3/14/12  
**Date Received:** 3/16/12  
**Date Analyzed:** 3/24/12 & 3/26/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.050 Liter(s)

Initial Pressure (psig): -3.45      Final Pressure (psig): 3.74

Canister Dilution Factor: 1.64

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
156-59-2	cis-1,2-Dichloroethene	ND	0.82	ND	0.21	
141-78-6	Ethyl Acetate	5.1	1.6	1.4	0.46	
110-54-3	n-Hexane	7.7	0.82	2.2	0.23	
67-66-3	Chloroform	2.0	0.82	0.41	0.17	
109-99-9	Tetrahydrofuran (THF)	ND	0.82	ND	0.28	
107-06-2	1,2-Dichloroethane	ND	0.82	ND	0.20	
71-55-6	1,1,1-Trichloroethane	ND	0.82	ND	0.15	
71-43-2	Benzene	7.6	0.82	2.4	0.26	
56-23-5	Carbon Tetrachloride	ND	0.82	ND	0.13	
110-82-7	Cyclohexane	ND	1.6	ND	0.48	
78-87-5	1,2-Dichloropropane	ND	0.82	ND	0.18	
75-27-4	Bromodichloromethane	ND	0.82	ND	0.12	
79-01-6	Trichloroethene	ND	0.82	ND	0.15	
123-91-1	1,4-Dioxane	ND	0.82	ND	0.23	
80-62-6	Methyl Methacrylate	ND	1.6	ND	0.40	
142-82-5	n-Heptane	3.2	0.82	0.79	0.20	
10061-01-5	cis-1,3-Dichloropropene	ND	0.82	ND	0.18	
108-10-1	4-Methyl-2-pentanone	ND	0.82	ND	0.20	
10061-02-6	trans-1,3-Dichloropropene	ND	0.82	ND	0.18	
79-00-5	1,1,2-Trichloroethane	ND	0.82	ND	0.15	
108-88-3	Toluene	30	0.82	7.9	0.22	
591-78-6	2-Hexanone	ND	0.82	ND	0.20	
124-48-1	Dibromochloromethane	ND	0.82	ND	0.096	
106-93-4	1,2-Dibromoethane	ND	0.82	ND	0.11	
123-86-4	n-Butyl Acetate	0.83	0.82	0.17	0.17	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 16-IF-031312  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201043  
**CAS Sample ID:** P1201043-007

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Lusine Hakobyan/Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01428

**Date Collected:** 3/14/12  
**Date Received:** 3/16/12  
**Date Analyzed:** 3/24/12 & 3/26/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.050 Liter(s)

Initial Pressure (psig): -3.45      Final Pressure (psig): 3.74

Canister Dilution Factor: 1.64

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
111-65-9	n-Octane	1.0	0.82	0.22	0.18	
127-18-4	Tetrachloroethene	ND	0.82	ND	0.12	
108-90-7	Chlorobenzene	ND	0.82	ND	0.18	
100-41-4	Ethylbenzene	3.0	0.82	0.70	0.19	
179601-23-1	m,p-Xylenes	12	1.6	2.7	0.38	
75-25-2	Bromoform	ND	0.82	ND	0.079	
100-42-5	Styrene	ND	0.82	ND	0.19	
95-47-6	o-Xylene	3.2	0.82	0.73	0.19	
111-84-2	n-Nonane	1.1	0.82	0.20	0.16	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.82	ND	0.12	
98-82-8	Cumene	ND	0.82	ND	0.17	
80-56-8	alpha-Pinene	5.0	0.82	0.90	0.15	
103-65-1	n-Propylbenzene	ND	0.82	ND	0.17	
622-96-8	4-Ethyltoluene	ND	0.82	ND	0.17	
108-67-8	1,3,5-Trimethylbenzene	ND	0.82	ND	0.17	
95-63-6	1,2,4-Trimethylbenzene	2.1	0.82	0.44	0.17	
100-44-7	Benzyl Chloride	ND	0.82	ND	0.16	
541-73-1	1,3-Dichlorobenzene	ND	0.82	ND	0.14	
106-46-7	1,4-Dichlorobenzene	ND	0.82	ND	0.14	
95-50-1	1,2-Dichlorobenzene	ND	0.82	ND	0.14	
5989-27-5	d-Limonene	16	0.82	2.9	0.15	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.82	ND	0.085	
120-82-1	1,2,4-Trichlorobenzene	ND	0.82	ND	0.11	
91-20-3	Naphthalene	ND	0.82	ND	0.16	
87-68-3	Hexachlorobutadiene	ND	0.82	ND	0.077	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
CAS Sample ID: P120324-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/24/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	ND	0.50	ND	0.29	
75-71-8	Dichlorodifluoromethane (CFC 12)	ND	0.50	ND	0.10	
74-87-3	Chloromethane	ND	0.50	ND	0.24	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.50	ND	0.072	
75-01-4	Vinyl Chloride	ND	0.50	ND	0.20	
106-99-0	1,3-Butadiene	ND	0.50	ND	0.23	
74-83-9	Bromomethane	ND	0.50	ND	0.13	
75-00-3	Chloroethane	ND	0.50	ND	0.19	
64-17-5	Ethanol	ND	5.0	ND	2.7	
75-05-8	Acetonitrile	ND	0.50	ND	0.30	
107-02-8	Acrolein	ND	2.0	ND	0.87	
67-64-1	Acetone	ND	5.0	ND	2.1	
75-69-4	Trichlorofluoromethane	ND	0.50	ND	0.089	
67-63-0	2-Propanol (Isopropyl Alcohol)	ND	1.0	ND	0.41	
107-13-1	Acrylonitrile	ND	0.50	ND	0.23	
75-35-4	1,1-Dichloroethene	ND	0.50	ND	0.13	
75-09-2	Methylene Chloride	ND	0.50	ND	0.14	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.50	ND	0.16	
76-13-1	Trichlorotrifluoroethane	ND	0.50	ND	0.065	
75-15-0	Carbon Disulfide	ND	5.0	ND	1.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.50	ND	0.13	
75-34-3	1,1-Dichloroethane	ND	0.50	ND	0.12	
1634-04-4	Methyl tert-Butyl Ether	ND	0.50	ND	0.14	
108-05-4	Vinyl Acetate	ND	5.0	ND	1.4	
78-93-3	2-Butanone (MEK)	ND	5.0	ND	1.7	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
CAS Sample ID: P120324-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/24/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
156-59-2	cis-1,2-Dichloroethene	ND	0.50	ND	0.13	
141-78-6	Ethyl Acetate	ND	1.0	ND	0.28	
110-54-3	n-Hexane	ND	0.50	ND	0.14	
67-66-3	Chloroform	ND	0.50	ND	0.10	
109-99-9	Tetrahydrofuran (THF)	ND	0.50	ND	0.17	
107-06-2	1,2-Dichloroethane	ND	0.50	ND	0.12	
71-55-6	1,1,1-Trichloroethane	ND	0.50	ND	0.092	
71-43-2	Benzene	ND	0.50	ND	0.16	
56-23-5	Carbon Tetrachloride	ND	0.50	ND	0.080	
110-82-7	Cyclohexane	ND	1.0	ND	0.29	
78-87-5	1,2-Dichloropropane	ND	0.50	ND	0.11	
75-27-4	Bromodichloromethane	ND	0.50	ND	0.075	
79-01-6	Trichloroethene	ND	0.50	ND	0.093	
123-91-1	1,4-Dioxane	ND	0.50	ND	0.14	
80-62-6	Methyl Methacrylate	ND	1.0	ND	0.24	
142-82-5	n-Heptane	ND	0.50	ND	0.12	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-10-1	4-Methyl-2-pentanone	ND	0.50	ND	0.12	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	ND	0.11	
79-00-5	1,1,2-Trichloroethane	ND	0.50	ND	0.092	
108-88-3	Toluene	ND	0.50	ND	0.13	
591-78-6	2-Hexanone	ND	0.50	ND	0.12	
124-48-1	Dibromochloromethane	ND	0.50	ND	0.059	
106-93-4	1,2-Dibromoethane	ND	0.50	ND	0.065	
123-86-4	n-Butyl Acetate	ND	0.50	ND	0.11	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P120324-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
111-65-9	n-Octane	ND	0.50	ND	0.11	
127-18-4	Tetrachloroethene	ND	0.50	ND	0.074	
108-90-7	Chlorobenzene	ND	0.50	ND	0.11	
100-41-4	Ethylbenzene	ND	0.50	ND	0.12	
179601-23-1	m,p-Xylenes	ND	1.0	ND	0.23	
75-25-2	Bromoform	ND	0.50	ND	0.048	
100-42-5	Styrene	ND	0.50	ND	0.12	
95-47-6	o-Xylene	ND	0.50	ND	0.12	
111-84-2	n-Nonane	ND	0.50	ND	0.095	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	ND	0.073	
98-82-8	Cumene	ND	0.50	ND	0.10	
80-56-8	alpha-Pinene	ND	0.50	ND	0.090	
103-65-1	n-Propylbenzene	ND	0.50	ND	0.10	
622-96-8	4-Ethyltoluene	ND	0.50	ND	0.10	
108-67-8	1,3,5-Trimethylbenzene	ND	0.50	ND	0.10	
95-63-6	1,2,4-Trimethylbenzene	ND	0.50	ND	0.10	
100-44-7	Benzyl Chloride	ND	0.50	ND	0.097	
541-73-1	1,3-Dichlorobenzene	ND	0.50	ND	0.083	
106-46-7	1,4-Dichlorobenzene	ND	0.50	ND	0.083	
95-50-1	1,2-Dichlorobenzene	ND	0.50	ND	0.083	
5989-27-5	d-Limonene	ND	0.50	ND	0.090	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.50	ND	0.052	
120-82-1	1,2,4-Trichlorobenzene	ND	0.50	ND	0.067	
91-20-3	Naphthalene	ND	0.50	ND	0.095	
87-68-3	Hexachlorobutadiene	ND	0.50	ND	0.047	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
CAS Sample ID: P120326-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/26/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	ND	0.50	ND	0.29	
75-71-8	Dichlorodifluoromethane (CFC 12)	ND	0.50	ND	0.10	
74-87-3	Chloromethane	ND	0.50	ND	0.24	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.50	ND	0.072	
75-01-4	Vinyl Chloride	ND	0.50	ND	0.20	
106-99-0	1,3-Butadiene	ND	0.50	ND	0.23	
74-83-9	Bromomethane	ND	0.50	ND	0.13	
75-00-3	Chloroethane	ND	0.50	ND	0.19	
64-17-5	Ethanol	ND	5.0	ND	2.7	
75-05-8	Acetonitrile	ND	0.50	ND	0.30	
107-02-8	Acrolein	ND	2.0	ND	0.87	
67-64-1	Acetone	ND	5.0	ND	2.1	
75-69-4	Trichlorofluoromethane	ND	0.50	ND	0.089	
67-63-0	2-Propanol (Isopropyl Alcohol)	ND	1.0	ND	0.41	
107-13-1	Acrylonitrile	ND	0.50	ND	0.23	
75-35-4	1,1-Dichloroethene	ND	0.50	ND	0.13	
75-09-2	Methylene Chloride	ND	0.50	ND	0.14	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.50	ND	0.16	
76-13-1	Trichlorotrifluoroethane	ND	0.50	ND	0.065	
75-15-0	Carbon Disulfide	ND	5.0	ND	1.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.50	ND	0.13	
75-34-3	1,1-Dichloroethane	ND	0.50	ND	0.12	
1634-04-4	Methyl tert-Butyl Ether	ND	0.50	ND	0.14	
108-05-4	Vinyl Acetate	ND	5.0	ND	1.4	
78-93-3	2-Butanone (MEK)	ND	5.0	ND	1.7	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
CAS Sample ID: P120326-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/26/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
156-59-2	cis-1,2-Dichloroethene	ND	0.50	ND	0.13	
141-78-6	Ethyl Acetate	ND	1.0	ND	0.28	
110-54-3	n-Hexane	ND	0.50	ND	0.14	
67-66-3	Chloroform	ND	0.50	ND	0.10	
109-99-9	Tetrahydrofuran (THF)	ND	0.50	ND	0.17	
107-06-2	1,2-Dichloroethane	ND	0.50	ND	0.12	
71-55-6	1,1,1-Trichloroethane	ND	0.50	ND	0.092	
71-43-2	Benzene	ND	0.50	ND	0.16	
56-23-5	Carbon Tetrachloride	ND	0.50	ND	0.080	
110-82-7	Cyclohexane	ND	1.0	ND	0.29	
78-87-5	1,2-Dichloropropane	ND	0.50	ND	0.11	
75-27-4	Bromodichloromethane	ND	0.50	ND	0.075	
79-01-6	Trichloroethene	ND	0.50	ND	0.093	
123-91-1	1,4-Dioxane	ND	0.50	ND	0.14	
80-62-6	Methyl Methacrylate	ND	1.0	ND	0.24	
142-82-5	n-Heptane	ND	0.50	ND	0.12	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-10-1	4-Methyl-2-pentanone	ND	0.50	ND	0.12	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	ND	0.11	
79-00-5	1,1,2-Trichloroethane	ND	0.50	ND	0.092	
108-88-3	Toluene	ND	0.50	ND	0.13	
591-78-6	2-Hexanone	ND	0.50	ND	0.12	
124-48-1	Dibromochloromethane	ND	0.50	ND	0.059	
106-93-4	1,2-Dibromoethane	ND	0.50	ND	0.065	
123-86-4	n-Butyl Acetate	ND	0.50	ND	0.11	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P120326-MB

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/26/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
111-65-9	n-Octane	ND	0.50	ND	0.11	
127-18-4	Tetrachloroethene	ND	0.50	ND	0.074	
108-90-7	Chlorobenzene	ND	0.50	ND	0.11	
100-41-4	Ethylbenzene	ND	0.50	ND	0.12	
179601-23-1	m,p-Xylenes	ND	1.0	ND	0.23	
75-25-2	Bromoform	ND	0.50	ND	0.048	
100-42-5	Styrene	ND	0.50	ND	0.12	
95-47-6	o-Xylene	ND	0.50	ND	0.12	
111-84-2	n-Nonane	ND	0.50	ND	0.095	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	ND	0.073	
98-82-8	Cumene	ND	0.50	ND	0.10	
80-56-8	alpha-Pinene	ND	0.50	ND	0.090	
103-65-1	n-Propylbenzene	ND	0.50	ND	0.10	
622-96-8	4-Ethyltoluene	ND	0.50	ND	0.10	
108-67-8	1,3,5-Trimethylbenzene	ND	0.50	ND	0.10	
95-63-6	1,2,4-Trimethylbenzene	ND	0.50	ND	0.10	
100-44-7	Benzyl Chloride	ND	0.50	ND	0.097	
541-73-1	1,3-Dichlorobenzene	ND	0.50	ND	0.083	
106-46-7	1,4-Dichlorobenzene	ND	0.50	ND	0.083	
95-50-1	1,2-Dichlorobenzene	ND	0.50	ND	0.083	
5989-27-5	d-Limonene	ND	0.50	ND	0.090	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.50	ND	0.052	
120-82-1	1,2,4-Trichlorobenzene	ND	0.50	ND	0.067	
91-20-3	Naphthalene	ND	0.50	ND	0.095	
87-68-3	Hexachlorobutadiene	ND	0.50	ND	0.047	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister(s)  
 Test Notes:

Date(s) Collected: 3/13 - 3/14/12  
 Date(s) Received: 3/16/12  
 Date(s) Analyzed: 3/24 - 3/26/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		Percent Recovered	Percent Recovered	Percent Recovered		
Method Blank	P120324-MB	102	100	99	70-130	
Method Blank	P120326-MB	104	99	96	70-130	
Lab Control Sample	P120324-LCS	102	100	100	70-130	
Lab Control Sample	P120326-LCS	105	98	100	70-130	
15-IF-031212	P1201043-001	102	100	97	70-130	
52-IF-031212	P1201043-002	101	100	98	70-130	
52-IF-031212	P1201043-002DUP	102	99	95	70-130	
10-IF-031212	P1201043-003	102	99	96	70-130	
11-IF-031212	P1201043-004	102	99	97	70-130	
12-IF-031312	P1201043-005	102	99	99	70-130	
13-IF-031312	P1201043-006	103	97	98	70-130	
16-IF-031312	P1201043-007	102	99	99	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.



## LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P120324-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
115-07-1	Propene	198	183	92	54-137	
75-71-8	Dichlorodifluoromethane (CFC 12)	196	156	80	55-112	
74-87-3	Chloromethane	190	155	82	66-122	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	200	154	77	64-113	
75-01-4	Vinyl Chloride	194	154	79	68-115	
106-99-0	1,3-Butadiene	204	182	89	74-142	
74-83-9	Bromomethane	194	160	82	72-124	
75-00-3	Chloroethane	196	157	80	69-115	
64-17-5	Ethanol	928	800	86	67-127	
75-05-8	Acetonitrile	194	178	92	63-126	
107-02-8	Acrolein	198	164	83	62-127	
67-64-1	Acetone	1,010	833	82	67-106	
75-69-4	Trichlorofluoromethane	202	161	80	66-121	
67-63-0	2-Propanol (Isopropyl Alcohol)	382	304	80	56-112	
107-13-1	Acrylonitrile	198	183	92	78-128	
75-35-4	1,1-Dichloroethene	212	181	85	74-116	
75-09-2	Methylene Chloride	206	168	82	69-103	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	208	199	96	76-142	
76-13-1	Trichlorotrifluoroethane	206	175	85	69-118	
75-15-0	Carbon Disulfide	208	169	81	71-112	
156-60-5	trans-1,2-Dichloroethene	196	172	88	73-121	
75-34-3	1,1-Dichloroethane	200	176	88	71-118	
1634-04-4	Methyl tert-Butyl Ether	198	177	89	72-115	
108-05-4	Vinyl Acetate	952	902	95	51-150	
78-93-3	2-Butanone (MEK)	206	187	91	80-130	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P120324-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
156-59-2	cis-1,2-Dichloroethene	206	179	87	73-117	
141-78-6	Ethyl Acetate	398	357	90	79-126	
110-54-3	n-Hexane	198	163	82	68-109	
67-66-3	Chloroform	214	176	82	67-118	
109-99-9	Tetrahydrofuran (THF)	202	175	87	57-130	
107-06-2	1,2-Dichloroethane	200	175	88	62-121	
71-55-6	1,1,1-Trichloroethane	198	169	85	70-116	
71-43-2	Benzene	198	170	86	66-103	
56-23-5	Carbon Tetrachloride	202	178	88	68-123	
110-82-7	Cyclohexane	390	323	83	73-111	
78-87-5	1,2-Dichloropropane	198	166	84	74-114	
75-27-4	Bromodichloromethane	198	172	87	75-120	
79-01-6	Trichloroethene	194	157	81	65-109	
123-91-1	1,4-Dioxane	200	173	87	74-120	
80-62-6	Methyl Methacrylate	400	361	90	80-124	
142-82-5	n-Heptane	196	171	87	75-114	
10061-01-5	cis-1,3-Dichloropropene	188	167	89	79-120	
108-10-1	4-Methyl-2-pentanone	204	183	90	79-128	
10061-02-6	trans-1,3-Dichloropropene	210	193	92	83-131	
79-00-5	1,1,2-Trichloroethane	196	168	86	76-116	
108-88-3	Toluene	202	167	83	64-115	
591-78-6	2-Hexanone	222	196	88	73-120	
124-48-1	Dibromochloromethane	206	186	90	72-137	
106-93-4	1,2-Dibromoethane	200	169	85	70-126	
123-86-4	n-Butyl Acetate	220	191	87	65-127	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P120324-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
111-65-9	n-Octane	200	182	91	72-118	
127-18-4	Tetrachloroethene	184	148	80	63-123	
108-90-7	Chlorobenzene	202	165	82	65-117	
100-41-4	Ethylbenzene	198	164	83	69-118	
179601-23-1	m,p-Xylenes	392	336	86	68-119	
75-25-2	Bromoform	208	177	85	79-150	
100-42-5	Styrene	200	177	89	74-127	
95-47-6	o-Xylene	194	164	85	68-118	
111-84-2	n-Nonane	196	168	86	72-116	
79-34-5	1,1,2,2-Tetrachloroethane	192	159	83	72-135	
98-82-8	Cumene	190	161	85	68-119	
80-56-8	alpha-Pinene	186	163	88	70-123	
103-65-1	n-Propylbenzene	192	160	83	69-119	
622-96-8	4-Ethyltoluene	198	165	83	68-121	
108-67-8	1,3,5-Trimethylbenzene	200	169	85	67-118	
95-63-6	1,2,4-Trimethylbenzene	194	168	87	66-122	
100-44-7	Benzyl Chloride	200	192	96	73-144	
541-73-1	1,3-Dichlorobenzene	200	155	78	64-122	
106-46-7	1,4-Dichlorobenzene	206	153	74	65-125	
95-50-1	1,2-Dichlorobenzene	198	154	78	63-128	
5989-27-5	d-Limonene	200	182	91	72-126	
96-12-8	1,2-Dibromo-3-chloropropane	196	171	87	72-139	
120-82-1	1,2,4-Trichlorobenzene	196	157	80	65-139	
91-20-3	Naphthalene	172	141	82	60-142	
87-68-3	Hexachlorobutadiene	202	161	80	58-137	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P120326-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/26/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
115-07-1	Propene	198	191	96	54-137	
75-71-8	Dichlorodifluoromethane (CFC 12)	196	157	80	55-112	
74-87-3	Chloromethane	190	153	81	66-122	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	200	155	78	64-113	
75-01-4	Vinyl Chloride	194	154	79	68-115	
106-99-0	1,3-Butadiene	204	183	90	74-142	
74-83-9	Bromomethane	194	160	82	72-124	
75-00-3	Chloroethane	196	159	81	69-115	
64-17-5	Ethanol	928	796	86	67-127	
75-05-8	Acetonitrile	194	182	94	63-126	
107-02-8	Acrolein	198	166	84	62-127	
67-64-1	Acetone	1,010	839	83	67-106	
75-69-4	Trichlorofluoromethane	202	163	81	66-121	
67-63-0	2-Propanol (Isopropyl Alcohol)	382	293	77	56-112	
107-13-1	Acrylonitrile	198	185	93	78-128	
75-35-4	1,1-Dichloroethene	212	181	85	74-116	
75-09-2	Methylene Chloride	206	166	81	69-103	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	208	203	98	76-142	
76-13-1	Trichlorotrifluoroethane	206	173	84	69-118	
75-15-0	Carbon Disulfide	208	170	82	71-112	
156-60-5	trans-1,2-Dichloroethene	196	172	88	73-121	
75-34-3	1,1-Dichloroethane	200	178	89	71-118	
1634-04-4	Methyl tert-Butyl Ether	198	178	90	72-115	
108-05-4	Vinyl Acetate	952	885	93	51-150	
78-93-3	2-Butanone (MEK)	206	186	90	80-130	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P120326-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/26/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
156-59-2	cis-1,2-Dichloroethene	206	180	87	73-117	
141-78-6	Ethyl Acetate	398	356	89	79-126	
110-54-3	n-Hexane	198	166	84	68-109	
67-66-3	Chloroform	214	177	83	67-118	
109-99-9	Tetrahydrofuran (THF)	202	173	86	57-130	
107-06-2	1,2-Dichloroethane	200	179	90	62-121	
71-55-6	1,1,1-Trichloroethane	198	171	86	70-116	
71-43-2	Benzene	198	170	86	66-103	
56-23-5	Carbon Tetrachloride	202	178	88	68-123	
110-82-7	Cyclohexane	390	323	83	73-111	
78-87-5	1,2-Dichloropropane	198	167	84	74-114	
75-27-4	Bromodichloromethane	198	173	87	75-120	
79-01-6	Trichloroethene	194	156	80	65-109	
123-91-1	1,4-Dioxane	200	172	86	74-120	
80-62-6	Methyl Methacrylate	400	360	90	80-124	
142-82-5	n-Heptane	196	171	87	75-114	
10061-01-5	cis-1,3-Dichloropropene	188	169	90	79-120	
108-10-1	4-Methyl-2-pentanone	204	189	93	79-128	
10061-02-6	trans-1,3-Dichloropropene	210	195	93	83-131	
79-00-5	1,1,2-Trichloroethane	196	169	86	76-116	
108-88-3	Toluene	202	161	80	64-115	
591-78-6	2-Hexanone	222	195	88	73-120	
124-48-1	Dibromochloromethane	206	178	86	72-137	
106-93-4	1,2-Dibromoethane	200	162	81	70-126	
123-86-4	n-Butyl Acetate	220	189	86	65-127	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
CAS Sample ID: P120326-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/26/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
111-65-9	n-Octane	200	176	88	72-118	
127-18-4	Tetrachloroethene	184	141	77	63-123	
108-90-7	Chlorobenzene	202	158	78	65-117	
100-41-4	Ethylbenzene	198	161	81	69-118	
179601-23-1	m,p-Xylenes	392	325	83	68-119	
75-25-2	Bromoform	208	169	81	79-150	
100-42-5	Styrene	200	169	85	74-127	
95-47-6	o-Xylene	194	158	81	68-118	
111-84-2	n-Nonane	196	166	85	72-116	
79-34-5	1,1,2,2-Tetrachloroethane	192	155	81	72-135	
98-82-8	Cumene	190	156	82	68-119	
80-56-8	alpha-Pinene	186	157	84	70-123	
103-65-1	n-Propylbenzene	192	156	81	69-119	
622-96-8	4-Ethyltoluene	198	159	80	68-121	
108-67-8	1,3,5-Trimethylbenzene	200	163	82	67-118	
95-63-6	1,2,4-Trimethylbenzene	194	163	84	66-122	
100-44-7	Benzyl Chloride	200	185	93	73-144	
541-73-1	1,3-Dichlorobenzene	200	148	74	64-122	
106-46-7	1,4-Dichlorobenzene	206	146	71	65-125	
95-50-1	1,2-Dichlorobenzene	198	148	75	63-128	
5989-27-5	d-Limonene	200	177	89	72-126	
96-12-8	1,2-Dibromo-3-chloropropane	196	164	84	72-139	
120-82-1	1,2,4-Trichlorobenzene	196	150	77	65-139	
91-20-3	Naphthalene	172	134	78	60-142	
87-68-3	Hexachlorobutadiene	202	152	75	58-137	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result. Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

## LABORATORY DUPLICATE SUMMARY RESULTS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 52-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-002DUP

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01845

**Date Collected:** 3/13/12  
**Date Received:** 3/16/12  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -4.19

Final Pressure (psig): 3.67

Canister Dilution Factor: 1.75

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
Propene	3.36	1.95	3.17	1.84	3.265	6	25	
Dichlorodifluoromethane (CFC 12)	1.97	0.398	1.95	0.394	1.96	1	25	
Chloromethane	ND	ND	ND	ND	-	-	25	
1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	ND	ND	ND	-	-	25	
Vinyl Chloride	ND	ND	ND	ND	-	-	25	
1,3-Butadiene	ND	ND	ND	ND	-	-	25	
Bromomethane	ND	ND	ND	ND	-	-	25	
Chloroethane	ND	ND	ND	ND	-	-	25	
Ethanol	478	254	477	253	477.5	0.2	25	
Acetonitrile	ND	ND	ND	ND	-	-	25	
Acrolein	ND	ND	ND	ND	-	-	25	
Acetone	30.5	12.9	30.5	12.9	30.5	0	25	
Trichlorofluoromethane	2.55	0.455	2.56	0.456	2.555	0.4	25	
2-Propanol (Isopropyl Alcohol)	15.2	6.20	15.6	6.36	15.4	3	25	
Acrylonitrile	ND	ND	ND	ND	-	-	25	
1,1-Dichloroethene	ND	ND	ND	ND	-	-	25	
Methylene Chloride	ND	ND	ND	ND	-	-	25	
3-Chloro-1-propene (Allyl Chloride)	ND	ND	ND	ND	-	-	25	
Trichlorotrifluoroethane	ND	ND	ND	ND	-	-	25	
Carbon Disulfide	ND	ND	ND	ND	-	-	25	
trans-1,2-Dichloroethene	ND	ND	ND	ND	-	-	25	
1,1-Dichloroethane	ND	ND	ND	ND	-	-	25	
Methyl tert-Butyl Ether	ND	ND	ND	ND	-	-	25	
Vinyl Acetate	ND	ND	ND	ND	-	-	25	
2-Butanone (MEK)	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

## LABORATORY DUPLICATE SUMMARY RESULTS

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**Client:** Montana DEQ  
**Client Sample ID:** 52-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-002DUP

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01845

**Date Collected:** 3/13/12  
**Date Received:** 3/16/12  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -4.19

Final Pressure (psig): 3.67

Canister Dilution Factor: 1.75

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
cis-1,2-Dichloroethene	ND	ND	ND	ND	-	-	25	
Ethyl Acetate	9.65	2.68	9.51	2.64	9.58	1	25	
n-Hexane	ND	ND	ND	ND	-	-	25	
Chloroform	ND	ND	ND	ND	-	-	25	
Tetrahydrofuran (THF)	ND	ND	ND	ND	-	-	25	
1,2-Dichloroethane	ND	ND	ND	ND	-	-	25	
1,1,1-Trichloroethane	1.42	0.260	1.40	0.256	1.41	1	25	
Benzene	ND	ND	ND	ND	-	-	25	
Carbon Tetrachloride	ND	ND	ND	ND	-	-	25	
Cyclohexane	ND	ND	ND	ND	-	-	25	
1,2-Dichloropropane	ND	ND	ND	ND	-	-	25	
Bromodichloromethane	ND	ND	ND	ND	-	-	25	
Trichloroethene	ND	ND	ND	ND	-	-	25	
1,4-Dioxane	ND	ND	ND	ND	-	-	25	
Methyl Methacrylate	ND	ND	ND	ND	-	-	25	
n-Heptane	ND	ND	ND	ND	-	-	25	
cis-1,3-Dichloropropene	ND	ND	ND	ND	-	-	25	
4-Methyl-2-pentanone	ND	ND	ND	ND	-	-	25	
trans-1,3-Dichloropropene	ND	ND	ND	ND	-	-	25	
1,1,2-Trichloroethane	ND	ND	ND	ND	-	-	25	
Toluene	9.98	2.65	9.86	2.62	9.92	1	25	
2-Hexanone	ND	ND	ND	ND	-	-	25	
Dibromochloromethane	ND	ND	ND	ND	-	-	25	
1,2-Dibromoethane	ND	ND	ND	ND	-	-	25	
n-Butyl Acetate	2.10	0.443	2.08	0.439	2.09	1	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.



LABORATORY DUPLICATE SUMMARY RESULTS

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**Client:** Montana DEQ  
**Client Sample ID:** 52-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-002DUP

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01845

Date Collected: 3/13/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.19

Final Pressure (psig): 3.67

Canister Dilution Factor: 1.75

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
n-Octane	ND	ND	ND	ND	-	-	25	
Tetrachloroethene	ND	ND	ND	ND	-	-	25	
Chlorobenzene	ND	ND	ND	ND	-	-	25	
Ethylbenzene	ND	ND	ND	ND	-	-	25	
m,p-Xylenes	ND	ND	ND	ND	-	-	25	
Bromoform	ND	ND	ND	ND	-	-	25	
Styrene	ND	ND	ND	ND	-	-	25	
o-Xylene	ND	ND	ND	ND	-	-	25	
n-Nonane	ND	ND	ND	ND	-	-	25	
1,1,2,2-Tetrachloroethane	ND	ND	ND	ND	-	-	25	
Cumene	ND	ND	ND	ND	-	-	25	
alpha-Pinene	1.56	0.280	1.58	0.284	1.57	<b>1</b>	25	
n-Propylbenzene	ND	ND	ND	ND	-	-	25	
4-Ethyltoluene	ND	ND	ND	ND	-	-	25	
1,3,5-Trimethylbenzene	ND	ND	ND	ND	-	-	25	
1,2,4-Trimethylbenzene	ND	ND	ND	ND	-	-	25	
Benzyl Chloride	ND	ND	ND	ND	-	-	25	
1,3-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
1,4-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
1,2-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
d-Limonene	17.6	3.16	17.5	3.14	17.55	<b>0.6</b>	25	
1,2-Dibromo-3-chloropropane	ND	ND	ND	ND	-	-	25	
1,2,4-Trichlorobenzene	ND	ND	ND	ND	-	-	25	
Naphthalene	ND	ND	ND	ND	-	-	25	
Hexachlorobutadiene	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 15-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
CAS Sample ID: P1201043-001

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00187

Date Collected: 3/13/12  
Date Received: 3/16/12  
Date Analyzed: 3/27/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.79      Final Pressure (psig): 3.54

Canister Dilution Factor: 1.53

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.038	ND	0.015	
107-06-2	1,2-Dichloroethane	<b>0.19</b>	0.038	<b>0.048</b>	0.0095	
71-43-2	Benzene	<b>0.76</b>	0.11	<b>0.24</b>	0.036	
79-01-6	Trichloroethene	ND	0.038	ND	0.0071	
79-00-5	1,1,2-Trichloroethane	ND	0.15	ND	0.028	
127-18-4	Tetrachloroethene	<b>0.098</b>	0.038	<b>0.015</b>	0.0056	
100-41-4	Ethylbenzene	<b>0.53</b>	0.15	<b>0.12</b>	0.035	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.038	ND	0.0056	
91-20-3	Naphthalene	<b>0.24</b>	0.15	<b>0.046</b>	0.029	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 52-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-002

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01845

Date Collected: 3/13/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.19      Final Pressure (psig): 3.67

Canister Dilution Factor: 1.75

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.044	ND	0.017	
107-06-2	1,2-Dichloroethane	<b>0.19</b>	0.044	<b>0.048</b>	0.011	
71-43-2	Benzene	<b>0.75</b>	0.13	<b>0.24</b>	0.041	
79-01-6	Trichloroethene	ND	0.044	ND	0.0081	
79-00-5	1,1,2-Trichloroethane	ND	0.18	ND	0.032	
127-18-4	Tetrachloroethene	<b>0.14</b>	0.044	<b>0.020</b>	0.0065	
100-41-4	Ethylbenzene	<b>0.54</b>	0.18	<b>0.12</b>	0.040	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.044	ND	0.0064	
91-20-3	Naphthalene	<b>0.28</b>	0.18	<b>0.053</b>	0.033	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 10-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-003

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01842

Date Collected: 3/13/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.14      Final Pressure (psig): 3.52

Canister Dilution Factor: 1.34

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.034	ND	0.013	
107-06-2	1,2-Dichloroethane	<b>0.088</b>	0.034	<b>0.022</b>	0.0083	
71-43-2	Benzene	<b>0.81</b>	0.10	<b>0.25</b>	0.031	
79-01-6	Trichloroethene	ND	0.034	ND	0.0062	
79-00-5	1,1,2-Trichloroethane	ND	0.13	ND	0.025	
127-18-4	Tetrachloroethene	<b>0.069</b>	0.034	<b>0.010</b>	0.0049	
100-41-4	Ethylbenzene	<b>0.48</b>	0.13	<b>0.11</b>	0.031	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.034	ND	0.0049	
91-20-3	Naphthalene	<b>0.21</b>	0.13	<b>0.039</b>	0.026	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 11-IF-031212  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-004

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00188

Date Collected: 3/13/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.57      Final Pressure (psig): 3.67

Canister Dilution Factor: 1.51

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.038	ND	0.015	
107-06-2	1,2-Dichloroethane	<b>0.085</b>	0.038	<b>0.021</b>	0.0093	
71-43-2	Benzene	<b>0.52</b>	0.11	<b>0.16</b>	0.035	
79-01-6	Trichloroethene	ND	0.038	ND	0.0070	
79-00-5	1,1,2-Trichloroethane	ND	0.15	ND	0.028	
127-18-4	Tetrachloroethene	<b>0.074</b>	0.038	<b>0.011</b>	0.0056	
100-41-4	Ethylbenzene	<b>0.20</b>	0.15	<b>0.045</b>	0.035	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.038	ND	0.0055	
91-20-3	Naphthalene	ND	0.15	ND	0.029	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 12-IF-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-005

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00185

Date Collected: 3/14/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.74      Final Pressure (psig): 3.69

Canister Dilution Factor: 1.68

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.042	ND	0.016	
107-06-2	1,2-Dichloroethane	<b>2.1</b>	0.042	<b>0.51</b>	0.010	
71-43-2	Benzene	<b>1.7</b>	0.13	<b>0.53</b>	0.039	
79-01-6	Trichloroethene	<b>0.094</b>	0.042	<b>0.018</b>	0.0078	
79-00-5	1,1,2-Trichloroethane	ND	0.17	ND	0.031	
127-18-4	Tetrachloroethene	<b>0.13</b>	0.042	<b>0.020</b>	0.0062	
100-41-4	Ethylbenzene	<b>0.79</b>	0.17	<b>0.18</b>	0.039	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.042	ND	0.0061	
91-20-3	Naphthalene	<b>0.23</b>	0.17	<b>0.043</b>	0.032	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 13-IF-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-006

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00190

Date Collected: 3/14/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.01      Final Pressure (psig): 3.68

Canister Dilution Factor: 1.57

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.039	ND	0.015	
107-06-2	1,2-Dichloroethane	<b>0.17</b>	0.039	<b>0.043</b>	0.0097	
71-43-2	Benzene	<b>0.52</b>	0.12	<b>0.16</b>	0.037	
79-01-6	Trichloroethene	<b>0.042</b>	0.039	<b>0.0079</b>	0.0073	
79-00-5	1,1,2-Trichloroethane	ND	0.16	ND	0.029	
127-18-4	Tetrachloroethene	<b>0.083</b>	0.039	<b>0.012</b>	0.0058	
100-41-4	Ethylbenzene	<b>0.21</b>	0.16	<b>0.048</b>	0.036	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.039	ND	0.0057	
91-20-3	Naphthalene	<b>0.46</b>	0.16	<b>0.087</b>	0.030	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 16-IF-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P1201043-007

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01428

Date Collected: 3/14/12  
 Date Received: 3/16/12  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.45 Final Pressure (psig): 3.74

Canister Dilution Factor: 1.64

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.041	ND	0.016	
107-06-2	1,2-Dichloroethane	<b>0.12</b>	0.041	<b>0.029</b>	0.010	
71-43-2	Benzene	<b>7.3</b>	0.12	<b>2.3</b>	0.039	
79-01-6	Trichloroethene	ND	0.041	ND	0.0076	
79-00-5	1,1,2-Trichloroethane	ND	0.16	ND	0.030	
127-18-4	Tetrachloroethene	<b>0.070</b>	0.041	<b>0.010</b>	0.0060	
100-41-4	Ethylbenzene	<b>2.9</b>	0.16	<b>0.67</b>	0.038	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.041	ND	0.0060	
91-20-3	Naphthalene	<b>0.29</b>	0.16	<b>0.056</b>	0.031	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P120327-MB

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.025	ND	0.0098	
107-06-2	1,2-Dichloroethane	ND	0.025	ND	0.0062	
71-43-2	Benzene	ND	0.075	ND	0.023	
79-01-6	Trichloroethene	ND	0.025	ND	0.0047	
79-00-5	1,1,2-Trichloroethane	ND	0.10	ND	0.018	
127-18-4	Tetrachloroethene	ND	0.025	ND	0.0037	
100-41-4	Ethylbenzene	ND	0.10	ND	0.023	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.025	ND	0.0036	
91-20-3	Naphthalene	ND	0.10	ND	0.019	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

SURROGATE SPIKE RECOVERY RESULTS

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**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister(s)  
 Test Notes:

Date(s) Collected: 3/13 - 3/14/12  
 Date(s) Received: 3/16/12  
 Date(s) Analyzed: 3/27/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		% Recovered	% Recovered	% Recovered		
Method Blank	P120327-MB	101	113	114	70-130	
Lab Control Sample	P120327-LCS	102	112	118	70-130	
15-IF-031212	P1201043-001	104	114	111	70-130	
52-IF-031212	P1201043-002	102	113	108	70-130	
10-IF-031212	P1201043-003	101	112	113	70-130	
11-IF-031212	P1201043-004	100	114	115	70-130	
12-IF-031312	P1201043-005	99	112	115	70-130	
13-IF-031312	P1201043-006	99	112	117	70-130	
16-IF-031312	P1201043-007	99	112	119	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.

Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201043  
 CAS Sample ID: P120327-LCS

**Test Code:** EPA TO-15 SIM  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
**Analyst:** Karen Ryan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/27/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
75-01-4	Vinyl Chloride	3.88	3.08	79	56-127	
107-06-2	1,2-Dichloroethane	4.00	3.10	78	51-140	
71-43-2	Benzene	3.96	3.49	88	56-125	
79-01-6	Trichloroethene	3.88	2.95	76	51-127	
79-00-5	1,1,2-Trichloroethane	3.92	2.95	75	49-137	
127-18-4	Tetrachloroethene	3.68	2.87	78	58-134	
100-41-4	Ethylbenzene	3.96	3.00	76	56-126	
79-34-5	1,1,2,2-Tetrachloroethane	3.84	2.91	76	53-148	
91-20-3	Naphthalene	3.44	3.05	89	25-149	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

Massachusetts APH  
Hydrocarbon Ranges

ICAL Method: M8011412A.M

ICAL Date: 1/14/12

Instrument ID: MS08

areas

masses

	0.5	1	5	25	50	100	0.5	1	5	25	50	100
<b>C5-C8 Aliphatics</b>												
Isopentane	60478	112606	488243	2480105	4710733	9222808	0.520	1.04	5.20	26.0	52.0	104
n-Hexane	59953	110631	507126	2577865	4931102	9678199	0.490	0.98	4.90	24.5	49.0	98
Cyclohexane	63203	115230	548910	2874726	5524402	10890911	0.500	1.00	5.00	25.0	50.0	100
2,3-Dimethylpentane	68648	127222	585183	2998017	5742104	11319597	0.520	1.04	5.20	26.0	52.0	104
n-Heptane	66143	120286	558074	2884817	5509801	10873046	0.490	0.98	4.90	24.5	49.0	98
n-Octane	76169	137829	632891	3286017	6205035	12092514	0.505	1.01	5.05	25.3	50.5	101
<b>area sum:</b>	<b>394594</b>	<b>723804</b>	<b>3320427</b>	<b>17101547</b>	<b>32623177</b>	<b>64077075</b>	<b>3.025</b>	<b>6.050</b>	<b>30.25</b>	<b>151.30</b>	<b>302.50</b>	<b>605.00</b>

**C9-C12 Aliphatics**

	0.5	1	5	25	50	100	0.5	1	5	25	50	100
2,3-Dimethylheptane	78780	147724	682518	3501487	6691428	13148840	0.485	0.97	4.85	24.3	48.5	97
n-Nonane	77464	140680	662221	3417128	6555165	12940625	0.485	0.97	4.85	24.3	48.5	97
n-Decane	80593	147724	705259	3651028	7021424	13870530	0.490	0.98	4.90	24.5	49.0	98
2-Butylcyclohexane	88915	166782	776903	4026931	7749936	15415340	0.495	0.99	4.95	24.8	49.5	99
n-Undecane	80867	152557	719618	3772331	7263077	14385881	0.480	0.96	4.80	24.0	48.0	96
n-Dodecane	70045	143746	687255	3673524	7115012	14041583	0.500	1.00	5.00	25.0	50.0	100
<b>area sum:</b>	<b>476664</b>	<b>899213</b>	<b>4233774</b>	<b>22042429</b>	<b>42396042</b>	<b>83802799</b>	<b>2.935</b>	<b>5.870</b>	<b>29.35</b>	<b>146.90</b>	<b>293.50</b>	<b>587.00</b>

**C9-C10 Aromatics**

	0.5	1	5	25	50	100	0.5	1	5	25	50	100
Isopropylbenzene	10658	19169	87439	448287	849124	1695737	0.485	0.97	4.85	24.3	48.5	97
3-Ethyltoluene	11805	20564	97823	504510	977020	1973178	0.470	0.94	4.70	23.5	47.0	94
1,3,5-Trimethylbenzene	15759	29260	134073	700133	1353305	2726630	0.490	0.98	4.90	24.5	49.0	98
1,2,3-Trimethylbenzene	17195	31328	143473	757946	1493373	3063059	0.495	0.99	4.95	24.8	49.5	99
p-Isopropyltoluene	9490	18116	83401	440962	872799	1814314	0.460	0.92	4.60	23.0	46.0	92
<b>area sum:</b>	<b>64907</b>	<b>118437</b>	<b>546209</b>	<b>2851838</b>	<b>5545621</b>	<b>11272918</b>	<b>2.400</b>	<b>4.800</b>	<b>24.00</b>	<b>120.10</b>	<b>240.00</b>	<b>480.00</b>

Massachusetts APH  
Hydrocarbon Ranges

ICAL: M8011412A.M

ICAL Date: 1/14/12

Instrument ID: MS08

areas

<u>Internal Standards (TIC)</u>	0.5	1	5	25	50	100
Bromochloromethane (IS1)	1312793	1307309	1293586	1289283	1284194	1283817
1,4-Difluorobenzene (IS2)	2547079	2541159	2512498	2475255	2422119	2352159
Chlorobenzene-d5 (IS3)	2929189	2913374	2880258	2895080	2878314	2908396

Internal Standards (EIC)

Bromochloromethane (IS1)	249696	247899	245985	245585	244327	243561
1,4-Difluorobenzene (IS2)	1153657	1149947	1133233	1136223	1127203	1138730
Chlorobenzene-d5 (IS3)	1007702	1004638	991212	992925	982812	986950

Surrogates (TIC)

	0.5	1	5	25	50	100
1,2-Dichloroethane-d4	1162497	1154617	1105630	1126201	1103958	1095945
♀ Toluene-d8	3508018	3490493	3455382	3466688	3437986	3457323
♂ p-Bromofluorobenzene	2298220	2284797	2197239	2227396	2192159	2197999

C5-C8 Aliphatics

	0.5	1	5	25	50	100	RRF <sub>avg</sub>	%RSD
	2.8268	2.6009	2.4215	2.4870	2.3919	2.3252	2.509	7.24

C9-C12 Aliphatics

	0.5	1	5	25	50	100	RRF <sub>avg</sub>	%RSD
	4.0291	3.8120	3.6383	3.7780	3.6744	3.6163	3.758	4.09

C9-C10 Aromatics

	0.5	1	5	25	50	100	RRF <sub>avg</sub>	%RSD
	0.6709	0.6140	0.5740	0.5979	0.5878	0.5949	0.607	5.63

**Massachusetts APH**  
**Continuing Calibration Verification Check Sheet**

Data File Name: 03241202.D  
 Data File Path: J:\MS08\Data\2012\_03\24\  
 Operator: EM  
 Date Acquired: 3/24/12 1:17  
 Acq. Method File: TO15.M  
 Sample Name: 25ng TO-15/MAPH CCV STD  
 Misc Info: S25-03151201/S25-03161202  
 Instrument Name: MS08

Enter RRFs from current ICAL!

Internal Standards	RT	Area
7) 1,4-Difluorobenzene (IS2)	15.89	1097014
16) Chlorobenzene-d5 (IS3)	21.69	957299

C5-C8 Aliphatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
3) Isopentane	8.19	2331141	2.445	147.5	-2.53	-30	30	Pass
4) n-Hexane	13.09	2450350						
9) Cyclohexane	15.82	2698497						
10) 2,3-Dimethylpentane	16.17	2863333	Spike	ICAL				
11) n-Heptane	17.35	2745668	Amt (ng)	RRF				
14) n-Octane	20.69	3146714	151.30	2.509				
		<b>16235703</b>						

C9-C12 Aliphatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
18) 2,3-Dimethylheptane	22.31	3331254	3.712	145.1	-1.23	-30	30	Pass
19) n-Nonane	23.30	3243877						
25) n-Decane	25.27	3468097						
28) Butylcyclohexane	25.99	3790882	Spike	ICAL				
29) n-Undecane	26.77	3564701	Amt (ng)	RRF				
30) n-Dodecane	28.00	3481141	146.90	3.758				
		<b>20879952</b>						

C9-C10 Aromatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
22) Isopropylbenzene	23.78	435563	0.589	116.6	-2.92	-30	30	Pass
23) 1-Methyl-3-ethylbenzene	24.52	480909						
24) 1,3,5-Trimethylbenzene	24.66	662082						
26) p-Isopropyltoluene	25.68	417535	Spike	ICAL				
27) 1,2,3-Trimethylbenzene	25.68	713901	Amt (ng)	RRF				
		<b>2709990</b>	120.1	0.607				

EM 3/24/12

Response Factor Report MS08

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Mon Dec 12 09:45:45 2011  
 Response Via : Initial Calibration

Calibration Files

0.1 =12091118.D 0.2 =12091119.D 0.5 =12091120.D 1.0 =12091121.D 5.0 =12091122.D 25 =12091123.D  
 50 =12091124.D 100 =12091125.D

Compound	0.1	0.2	0.5	1.0	5.0	25	50	100	Avg	%RSD
1) IR Bromochloromethane...										
2) T Propene	2.487	1.965	1.952	1.692	2.194	2.303	2.185	2.036	2.102	11.62
3) T Dichlorodifluo...	3.545	2.805	2.940	2.748	2.432	2.361	2.278	2.199	2.664	16.73
4) T Chloromethane	3.290	2.811	2.875	2.512	2.176	2.286	1.977	1.664	2.449	21.62
5) T 1,2-Dichloro-1...	1.901	1.661	1.575	1.458	1.359	1.333	1.200	1.190	1.460	16.68
6) T Vinyl Chloride	2.906	2.590	2.631	2.374	2.145	2.129	1.950	1.939	2.333	15.07
7) T 1,3-Butadiene	2.145	1.788	1.790	1.718	1.581	1.689	1.621	1.616	1.743	10.33
8) T Bromomethane	1.687	1.328	1.349	1.212	1.094	1.142	1.111	1.092	1.252	16.21
9) T Chloroethane	1.671	1.259	1.333	1.193	1.101	1.091	1.046	1.044	1.217	17.29
10) T Ethanol	2.040	1.345	1.186	1.085	1.022	1.096	1.076	1.064	1.239	27.34
11) T Acetonitrile	4.365	2.973	2.723	2.546	2.480	2.562	2.482	2.468	2.825	22.85
12) T Acrolein	1.328	0.931	0.907	0.861	0.772	0.814	0.789	0.779	0.897	20.48
13) T Acetone	1.645	1.316	1.249	1.129	0.994	1.015	0.981	0.964	1.161	20.26
14) T Trichlorofluor...	2.776	2.385	2.496	2.272	2.097	2.135	2.042	2.011	2.277	11.59
15) T 2-Propanol (Is...		4.463	4.536	4.168	2.920	2.763	2.577	2.672	3.443	26.09
16) T Acrylonitrile	2.701	1.907	1.878	1.783	1.745	1.813	1.765	1.738	1.916	16.85
17) T 1,1-Dichloroet...	1.458	1.418	1.452	1.263	1.206	1.223	1.172	1.152	1.293	9.95
18) T 2-Methyl-2-Pro...	4.464	3.710	3.939	3.616	3.558	3.584			3.812	9.14
19) T Methylene Chlo...			1.958	1.578	1.294	1.273	1.220	1.201	1.421	20.88
20) T 3-Chloro-1-pro...	2.347	2.260	2.345	2.220	2.150	2.296	2.230	2.191	2.255	3.15
21) T Trichlorotrifl...	1.348	1.244	1.299	1.193	1.089	1.098	1.062	1.035	1.171	10.01
22) T Carbon Disulfide		6.210	5.678	5.057	4.654	4.699	4.488	4.424	5.030	13.40
23) T trans-1,2-Dich...	2.436	2.237	2.206	2.142	1.951	1.953	1.883	1.835	2.080	10.00
24) T 1,1-Dichloroet...	2.896	2.757	2.747	2.500	2.307	2.379	2.288	2.236	2.514	10.07
25) T Methyl tert-Bu...	5.128	4.611	4.650	4.307	4.119	4.185	4.042	3.984	4.378	8.93
26) T Vinyl Acetate	0.430	0.368	0.423	0.399	0.393	0.410	0.402	0.398	0.403	4.70
27) T 2-Butanone (MEK)	1.265	1.003	1.038	0.972	0.927	0.964	0.928	0.794	0.986	13.56
28) T cis-1,2-Dichlo...	2.442	2.006	2.101	1.938	1.826	1.836	1.764	1.720	1.960	12.05
29) T Diisopropyl Ether	1.446	1.223	1.323	1.262	1.171	1.188	1.153	1.148	1.239	8.26
30) T Ethyl Acetate	0.611	0.554	0.573	0.545	0.517	0.535	0.520	0.516	0.546	6.02
31) T n-Hexane	3.541	3.006	3.181	2.805	2.582	2.633	2.545	2.543	2.855	12.70

Response Factor Report MS08

Method Path	J:\MS08\Methods\										
Method File	R8120911A.M										
Title	EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)										
32) T	Chloroform	2.679	2.408	2.409	2.204	1.983	2.054	1.974	1.937	2.206	12.18
33) S	1,2-Dichloroet...	1.586	1.577	1.582	1.590	1.535	1.533	1.520	1.504	1.553	2.19
34) T	Tetrahydrofura...	1.274	0.932	1.100	0.938	0.910	0.879	0.830	0.821	0.961	15.96
35) T	Ethyl tert-But...	1.999	1.832	2.033	1.818	1.732	1.757	1.708	1.687	1.821	7.18
36) T	1,2-Dichloroet...	2.099	1.992	1.976	1.806	1.683	1.700	1.599	1.531	1.798	11.39
37) IR	1,4-Difluorobenzen...	-----ISTD-----									
38) T	1,1,1-Trichloror...	0.531	0.465	0.486	0.453	0.419	0.434	0.413	0.403	0.451	9.46
39) T	Isopropyl Acetate	0.227	0.205	0.220	0.206	0.194	0.201	0.192	0.188	0.204	6.71
40) T	1-Butanol	0.348	0.319	0.300	0.320	0.335	0.324	0.318	0.323	0.323	4.66
41) T	Benzene	1.633	1.385	1.359	1.227	1.136	1.158	1.101	1.070	1.258	15.11
42) T	Carbon Tetrach...	0.394	0.356	0.369	0.340	0.334	0.352	0.337	0.333	0.352	6.03
43) T	Cyclohexane	0.666	0.576	0.588	0.539	0.492	0.500	0.479	0.469	0.539	12.59
44) T	tert-Amyl Meth...	1.092	0.926	0.998	0.937	0.900	0.918	0.881	0.858	0.939	7.95
45) T	1,2-Dichloropr...	0.397	0.361	0.368	0.324	0.312	0.316	0.300	0.293	0.334	11.12
46) T	Bromodichlorom...	0.458	0.387	0.392	0.378	0.362	0.378	0.362	0.353	0.384	8.57
47) T	Trichloroethene	0.466	0.375	0.379	0.344	0.323	0.328	0.316	0.309	0.355	14.62
48) T	1,4-Dioxane	0.305	0.282	0.288	0.248	0.245	0.246	0.237	0.230	0.260	10.55
49) T	2,2,4-Trimethy...	1.964	1.629	1.686	1.546	1.447	1.461	1.387	1.337	1.557	12.98
50) T	Methyl Methacr...	0.156	0.139	0.150	0.135	0.136	0.142	0.136	0.135	0.141	5.53
51) T	n-Heptane	0.427	0.373	0.385	0.364	0.341	0.344	0.332	0.324	0.361	9.35
52) T	cis-1,3-Dichlo...	0.571	0.502	0.521	0.491	0.482	0.500	0.481	0.471	0.502	6.33
53) T	4-Methyl-2-pen...	0.359	0.295	0.308	0.303	0.297	0.303	0.289	0.284	0.305	7.68
54) T	trans-1,3-Dich...	0.476	0.413	0.443	0.421	0.429	0.453	0.437	0.427	0.437	4.59
55) T	1,1,2-Trichlor...	0.340	0.301	0.310	0.284	0.276	0.285	0.271	0.265	0.292	8.44
56) IR	Chlorobenzene-d5 (...)	-----ISTD-----									
57) S	Toluene-d8 (SS2)	2.305	2.306	2.305	2.309	2.306	2.300	2.310	2.304	2.306	0.14
58) T	Toluene	3.623	2.920	3.025	2.738	2.571	2.574	2.473	2.421	2.793	14.19
59) T	2-Hexanone	2.049	1.429	1.515	1.403	1.395	1.422	1.367	1.337	1.490	15.57
60) T	Dibromochlorom...	0.696	0.618	0.636	0.636	0.612	0.648	0.629	0.625	0.638	4.14
61) T	1,2-Dibromoethane	0.831	0.730	0.718	0.658	0.641	0.662	0.638	0.628	0.688	9.97
62) T	n-Butyl Acetate	2.338	1.850	1.857	1.721	1.704	1.765	1.737	1.737	1.839	11.40
63) T	n-Octane	0.785	0.652	0.714	0.643	0.602	0.611	0.580	0.572	0.645	11.28
64) T	Tetrachloroethene	1.090	0.904	0.904	0.826	0.767	0.779	0.756	0.755	0.848	13.63
65) T	Chlorobenzene	2.375	1.868	1.886	1.744	1.610	1.637	1.574	1.548	1.780	15.30
66) T	Ethylbenzene	3.846	3.171	3.188	2.956	2.803	2.848	2.743	2.680	3.030	12.51
67) T	m- & p-Xylenes	3.081	2.499	2.611	2.392	2.216	2.273	2.170	2.112	2.419	13.08
68) T	Bromoform	0.622	0.494	0.556	0.541	0.551	0.598	0.589	0.594	0.568	7.19
69) T	Styrene	2.478	1.895	1.937	1.800	1.735	1.803	1.726	1.675	1.881	13.64
70) T	o-Xylene	3.279	2.708	2.726	2.479	2.331	2.373	2.283	2.224	2.550	13.65



Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M

Title	: EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)										
71) T	n-Nonane	2.005	1.753	1.847	1.675	1.581	1.597	1.518	1.470	1.681	10.69
72) T	1,1,2,2-Tetrac...	1.348	1.164	1.262	1.150	1.083	1.121	1.085	1.066	1.160	8.48
73) S	Bromofluoroben...	0.774	0.771	0.779	0.781	0.775	0.778	0.782	0.780	0.777	0.50
74) T	Cumene	4.522	3.495	3.524	3.169	2.962	3.002	2.884	2.831	3.299	16.98
75) T	alpha-Pinene	1.782	1.636	1.734	1.601	1.537	1.578	1.518	1.498	1.611	6.35
76) T	n-Propylbenzene	5.441	4.229	4.306	3.924	3.710	3.754	3.553	3.394	4.039	16.01
77) T	3-Ethyltoluene	4.091	3.373	3.461	3.038	2.964	2.981	2.893	2.845	3.206	13.14
78) T	4-Ethyltoluene	4.003	3.202	3.301	3.151	2.837	2.983	2.810	2.776	3.133	12.84
79) T	1,3,5-Trimethy...	3.324	2.662	2.723	2.537	2.378	2.426	2.344	2.403	2.600	12.43
80) T	alpha-Methylst...	1.742	1.386	1.449	1.371	1.390	1.402	1.354	1.399	1.436	8.80
81) T	2-Ethyltoluene	4.143	3.362	3.430	3.169	2.975	3.048	2.919	2.985	3.254	12.43
82) T	1,2,4-Trimethy...	3.356	2.663	2.791	2.551	2.433	2.513	2.454	2.548	2.664	11.37
83) T	n-Decane	1.969	1.702	1.844	1.705	1.582	1.626	1.554	1.609	1.699	8.39
84) T	Benzyl Chloride	2.373	1.648	1.772	1.688	1.799	2.058	2.054	2.190	1.948	13.35
85) T	1,3-Dichlorobe...	2.498	1.658	1.649	1.455	1.377	1.409	1.374	1.436	1.607	23.49
86) T	1,4-Dichlorobe...	2.655	1.785	1.618	1.492	1.380	1.428	1.388	1.449	1.649	25.99
87) T	sec-Butylbenzene	4.527	3.773	3.816	3.471	3.264	3.328	3.214	3.176	3.571	12.80
88) T	4-Isopropyltol...	4.426	3.674	3.683	3.416	3.254	3.356	3.253	3.366	3.554	10.98
89) T	1,2,3-Trimethy...	3.266	2.650	2.852	2.588	2.471	2.573	2.506	2.619	2.691	9.63
90) T	1,2-Dichlorobe...	2.173	1.589	1.545	1.435	1.332	1.387	1.370	1.415	1.531	17.89
91) T	d-Limonene	1.124	0.998	1.106	1.069	1.064	1.072	1.040	1.065	1.067	3.58
92) T	1,2-Dibromo-3-...	0.563	0.488	0.501	0.486	0.494	0.525	0.532	0.542	0.517	5.50
93) T	n-Undecane	1.922	1.637	1.755	1.671	1.608	1.630	1.608	1.592	1.678	6.63
94) T	1,2,4-Trichlor...		1.335	1.165	1.062	1.017	1.071	1.086	1.118	1.122	9.32
95) T	Naphthalene		4.848	3.923	3.638	3.497	3.721	3.755	3.783	3.881	11.50
96) T	n-Dodecane	1.866	1.599	1.669	1.586	1.545	1.575	1.553	1.509	1.613	6.99
97) T	Hexachlorobuta...	0.841	0.749	0.730	0.673	0.635	0.664	0.663	0.676	0.704	9.50
98) T	Cyclohexanone	1.172	0.979	1.021	0.969	0.952	0.962	0.931	0.903	0.986	8.37
99) T	tert-Butylbenzene	3.486	2.799	2.880	2.658	2.449	2.533	2.453	2.528	2.723	12.71
100) T	n-Butylbenzene	3.596	2.904	2.958	2.670	2.572	2.625	2.613	2.593	2.817	12.33

(#) = Out of Range

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\24\  
 Data File : 03241201.D  
 Acq On : 24 Mar 2012 00:34  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 24 08:11:33 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	IR Bromochloromethane (IS1)	1.000	1.000	0.0	83	-0.01
2	T Propene	2.102	1.911	9.1	69	0.00
3	T Dichlorodifluoromethane (CF	2.664	2.198	17.5	77	0.01
4	T Chloromethane	2.449	2.091	14.6	76	0.00
5	T 1,2-Dichloro-1,1,2,2-tetra	1.460	1.164	20.3	73	0.00
6	T Vinyl Chloride	2.333	1.919	17.7	75	0.00
7	T 1,3-Butadiene	1.743	1.594	8.5	78	0.00
8	T Bromomethane	1.252	1.056	15.7	77	0.00
9	T Chloroethane	1.217	1.019	16.3	78	0.00
10	T Ethanol	1.239	1.162	6.2	88	-0.07
11	T Acetonitrile	2.825	2.548	9.8	83	-0.05
12	T Acrolein	0.897	0.744	17.1	76	-0.02
13	T Acetone	1.161	0.998	14.0	82	-0.04
14	T Trichlorofluoromethane	2.277	2.018	11.4	79	0.00
15	T 2-Propanol (Isopropanol)	3.443	3.218	6.5	97	-0.05
16	T Acrylonitrile	1.916	1.760	8.1	81	-0.03
17	T 1,1-Dichloroethene	1.293	1.152	10.9	78	0.00
18	T 2-Methyl-2-Propanol (tert-B	3.812	3.168	16.9	74	-0.03
19	T Methylene Chloride	1.421	1.188	16.4	78	-0.01
20	T 3-Chloro-1-propene (Allyl C	2.255	2.246	0.4	81	-0.01
21	T Trichlorotrifluoroethane	1.171	1.027	12.3	78	0.00
22	T Carbon Disulfide	5.030	4.402	12.5	78	0.00
23	T trans-1,2-Dichloroethene	2.080	1.877	9.8	80	0.00
24	T 1,1-Dichloroethane	2.514	2.309	8.2	81	-0.01
25	T Methyl tert-Butyl Ether	4.378	4.054	7.4	81	0.00
26	T Vinyl Acetate	0.403	0.395	2.0	80	-0.03
27	T 2-Butanone (MEK)	0.986	0.934	5.3	81	-0.02
28	T cis-1,2-Dichloroethene	1.960	1.765	9.9	80	-0.01
29	T Diisopropyl Ether	1.239	1.143	7.7	80	0.00
30	T Ethyl Acetate	0.546	0.519	4.9	81	-0.02
31	T n-Hexane	2.855	2.469	13.5	78	0.00
32	T Chloroform	2.206	1.955	11.4	79	-0.03
33	S 1,2-Dichloroethane-d4 (SS1)	1.553	1.606	-3.4	87	-0.01
34	T Tetrahydrofuran (THF)	0.961	0.863	10.2	82	0.00
35	T Ethyl tert-Butyl Ether	1.821	1.683	7.6	80	0.00
36	T 1,2-Dichloroethane	1.798	1.636	9.0	80	-0.01
37	IR 1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	86	0.00
38	T 1,1,1-Trichloroethane	0.451	0.399	11.5	79	0.00

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\24\  
 Data File : 03241201.D  
 Acq On : 24 Mar 2012 00:34  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 24 08:11:33 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
39 T	Isopropyl Acetate	0.204	0.187	8.3	80	-0.02
40 T	1-Butanol	0.323	0.343	-6.2	88	-0.05
41 T	Benzene	1.258	1.072	14.8	80	-0.01
42 T	Carbon Tetrachloride	0.352	0.318	9.7	78	-0.01
43 T	Cyclohexane	0.539	0.465	13.7	80	-0.01
44 T	tert-Amyl Methyl Ether	0.939	0.851	9.4	80	0.00
45 T	1,2-Dichloropropane	0.334	0.293	12.3	80	-0.01
46 T	Bromodichloromethane	0.384	0.350	8.9	80	0.00
47 T	Trichloroethene	0.355	0.296	16.6	78	-0.01
48 T	1,4-Dioxane	0.260	0.235	9.6	82	-0.02
49 T	2,2,4-Trimethylpentane (Iso	1.557	1.379	11.4	81	0.00
50 T	Methyl Methacrylate	0.141	0.132	6.4	81	-0.02
51 T	n-Heptane	0.361	0.325	10.0	81	-0.01
52 T	cis-1,3-Dichloropropene	0.502	0.468	6.8	81	0.00
53 T	4-Methyl-2-pentanone	0.305	0.289	5.2	82	-0.02
54 T	trans-1,3-Dichloropropene	0.437	0.422	3.4	80	0.00
55 T	1,1,2-Trichloroethane	0.292	0.259	11.3	78	-0.01
56 IR	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	86	0.00
57 S	Toluene-d8 (SS2)	2.306	2.302	0.2	86	0.00
58 T	Toluene	2.793	2.399	14.1	80	0.00
59 T	2-Hexanone	1.490	1.370	8.1	83	-0.01
60 T	Dibromochloromethane	0.638	0.588	7.8	78	0.00
61 T	1,2-Dibromoethane	0.688	0.601	12.6	78	0.00
62 T	n-Butyl Acetate	1.839	1.689	8.2	82	-0.01
63 T	n-Octane	0.645	0.606	6.0	85	0.00
64 T	Tetrachloroethene	0.848	0.706	16.7	78	0.00
65 T	Chlorobenzene	1.780	1.511	15.1	79	0.00
66 T	Ethylbenzene	3.030	2.603	14.1	78	0.00
67 T	m- & p-Xylenes	2.419	2.148	11.2	81	-0.01
68 T	Bromoform	0.568	0.547	3.7	78	0.00
69 T	Styrene	1.881	1.716	8.8	82	-0.01
70 T	o-Xylene	2.550	2.246	11.9	81	-0.01
71 T	n-Nonane	1.681	1.511	10.1	81	0.00
72 T	1,1,2,2-Tetrachloroethane	1.160	1.028	11.4	79	-0.01
73 S	Bromofluorobenzene (SS3)	0.777	0.795	-2.3	88	0.00
74 T	Cumene	3.299	2.937	11.0	84	0.00
75 T	alpha-Pinene	1.611	1.440	10.6	78	0.00
76 T	n-Propylbenzene	4.039	3.517	12.9	80	0.00

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\24\  
 Data File : 03241201.D  
 Acq On : 24 Mar 2012 00:34  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 24 08:11:33 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
77 T	3-Ethyltoluene	3.206	2.756	14.0	79	0.00
78 T	4-Ethyltoluene	3.133	2.746	12.4	79	-0.01
79 T	1,3,5-Trimethylbenzene	2.600	2.260	13.1	80	-0.01
80 T	alpha-Methylstyrene	1.436	1.296	9.7	79	-0.01
81 T	2-Ethyltoluene	3.254	2.808	13.7	79	-0.01
82 T	1,2,4-Trimethylbenzene	2.664	2.335	12.3	80	-0.01
83 T	n-Decane	1.699	1.527	10.1	81	-0.01
84 T	Benzyl Chloride	1.948	2.052	-5.3	86	-0.02
85 T	1,3-Dichlorobenzene	1.607	1.288	19.9	79	-0.01
86 T	1,4-Dichlorobenzene	1.649	1.298	21.3	78	-0.01
87 T	sec-Butylbenzene	3.571	3.106	13.0	80	-0.01
88 T	4-Isopropyltoluene (p-Cymen	3.554	3.122	12.2	80	-0.01
89 T	1,2,3-Trimethylbenzene	2.691	2.373	11.8	79	0.00
90 T	1,2-Dichlorobenzene	1.531	1.255	18.0	78	-0.01
91 T	d-Limonene	1.067	0.988	7.4	79	-0.01
92 T	1,2-Dibromo-3-Chloropropane	0.517	0.478	7.5	78	-0.01
93 T	n-Undecane	1.678	1.536	8.5	81	0.00
94 T	1,2,4-Trichlorobenzene	1.122	0.967	13.8	77	-0.01
95 T	Naphthalene	3.881	3.454	11.0	80	0.00
96 T	n-Dodecane	1.613	1.520	5.8	83	0.00
97 T	Hexachlorobutadiene	0.704	0.595	15.5	77	0.00
98 T	Cyclohexanone	0.986	0.960	2.6	86	-0.02
99 T	tert-Butylbenzene	2.723	2.380	12.6	81	-0.01
100 T	n-Butylbenzene	2.817	2.441	13.3	80	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\26\  
 Data File : 03261201.D  
 Acq On : 26 Mar 2012 8:12  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 26 10:52:52 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
1	IR Bromochloromethane (IS1)	1.000	1.000	0.0	84	-0.01
2	T Propene	2.102	1.914	8.9	70	0.01
3	T Dichlorodifluoromethane (CF	2.664	2.234	16.1	80	0.01
4	T Chloromethane	2.449	2.138	12.7	79	0.01
5	T 1,2-Dichloro-1,1,2,2-tetra	1.460	1.163	20.3	74	0.00
6	T Vinyl Chloride	2.333	1.948	16.5	77	0.00
7	T 1,3-Butadiene	1.743	1.643	5.7	82	0.00
8	T Bromomethane	1.252	1.073	14.3	79	0.00
9	T Chloroethane	1.217	1.032	15.2	80	0.00
10	T Ethanol	1.239	1.177	5.0	91	-0.06
11	T Acetonitrile	2.825	2.611	7.6	86	-0.04
12	T Acrolein	0.897	0.758	15.5	79	-0.02
13	T Acetone	1.161	1.022	12.0	85	-0.04
14	T Trichlorofluoromethane	2.277	2.061	9.5	81	0.00
15	T 2-Propanol (Isopropanol)	3.443	3.491	-1.4	107	-0.04
16	T Acrylonitrile	1.916	1.792	6.5	83	-0.03
17	T 1,1-Dichloroethene	1.293	1.150	11.1	79	0.00
18	T 2-Methyl-2-Propanol (tert-B	3.812	3.044	20.1	72	-0.02
19	T Methylene Chloride	1.421	1.194	16.0	79	0.00
20	T 3-Chloro-1-propene (Allyl C	2.255	2.342	-3.9	86	-0.01
21	T Trichlorotrifluoroethane	1.171	1.029	12.1	79	0.00
22	T Carbon Disulfide	5.030	4.487	10.8	81	0.00
23	T trans-1,2-Dichloroethene	2.080	1.929	7.3	83	0.00
24	T 1,1-Dichloroethane	2.514	2.351	6.5	83	-0.01
25	T Methyl tert-Butyl Ether	4.378	4.109	6.1	83	0.00
26	T Vinyl Acetate	0.403	0.398	1.2	82	-0.03
27	T 2-Butanone (MEK)	0.986	0.944	4.3	83	-0.02
28	T cis-1,2-Dichloroethene	1.960	1.798	8.3	83	-0.01
29	T Diisopropyl Ether	1.239	1.141	7.9	81	0.00
30	T Ethyl Acetate	0.546	0.526	3.7	83	-0.02
31	T n-Hexane	2.855	2.516	11.9	81	0.00
32	T Chloroform	2.206	1.991	9.7	82	-0.03
33	S 1,2-Dichloroethane-d4 (SS1)	1.553	1.635	-5.3	90	-0.01
34	T Tetrahydrofuran (THF)	0.961	0.875	8.9	84	0.00
35	T Ethyl tert-Butyl Ether	1.821	1.692	7.1	81	-0.01
36	T 1,2-Dichloroethane	1.798	1.684	6.3	84	-0.01
37	IR 1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	88	0.00
38	T 1,1,1-Trichloroethane	0.451	0.401	11.1	81	0.00

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\26\  
 Data File : 03261201.D  
 Acq On : 26 Mar 2012 8:12  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 26 10:52:52 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
39 T	Isopropyl Acetate	0.204	0.190	6.9	83	-0.02
40 T	1-Butanol	0.323	0.349	-8.0	92	-0.05
41 T	Benzene	1.258	1.071	14.9	81	-0.01
42 T	Carbon Tetrachloride	0.352	0.319	9.4	80	-0.01
43 T	Cyclohexane	0.539	0.465	13.7	82	-0.01
44 T	tert-Amyl Methyl Ether	0.939	0.856	8.8	82	-0.01
45 T	1,2-Dichloropropane	0.334	0.294	12.0	82	-0.01
46 T	Bromodichloromethane	0.384	0.349	9.1	81	0.00
47 T	Trichloroethene	0.355	0.294	17.2	79	-0.01
48 T	1,4-Dioxane	0.260	0.234	10.0	83	-0.02
49 T	2,2,4-Trimethylpentane (Iso	1.557	1.386	11.0	83	-0.01
50 T	Methyl Methacrylate	0.141	0.133	5.7	82	-0.02
51 T	n-Heptane	0.361	0.327	9.4	84	-0.01
52 T	cis-1,3-Dichloropropene	0.502	0.470	6.4	83	0.00
53 T	4-Methyl-2-pentanone	0.305	0.294	3.6	85	-0.02
54 T	trans-1,3-Dichloropropene	0.437	0.425	2.7	82	-0.01
55 T	1,1,2-Trichloroethane	0.292	0.259	11.3	80	-0.01
56 IR	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	89	0.00
57 S	Toluene-d8 (SS2)	2.306	2.288	0.8	88	0.00
58 T	Toluene	2.793	2.371	15.1	82	0.00
59 T	2-Hexanone	1.490	1.390	6.7	87	-0.02
60 T	Dibromochloromethane	0.638	0.586	8.2	80	0.00
61 T	1,2-Dibromoethane	0.688	0.597	13.2	80	-0.01
62 T	n-Butyl Acetate	1.839	1.708	7.1	86	-0.01
63 T	n-Octane	0.645	0.610	5.4	88	-0.01
64 T	Tetrachloroethene	0.848	0.690	18.6	78	0.00
65 T	Chlorobenzene	1.780	1.493	16.1	81	0.00
66 T	Ethylbenzene	3.030	2.671	11.8	83	0.00
67 T	m- & p-Xylenes	2.419	2.135	11.7	83	-0.02
68 T	Bromoform	0.568	0.534	6.0	79	-0.01
69 T	Styrene	1.881	1.689	10.2	83	-0.01
70 T	o-Xylene	2.550	2.229	12.6	83	-0.01
71 T	n-Nonane	1.681	1.513	10.0	84	-0.01
72 T	1,1,2,2-Tetrachloroethane	1.160	1.030	11.2	81	-0.02
73 S	Bromofluorobenzene (SS3)	0.777	0.764	1.7	87	0.00
74 T	Cumene	3.299	2.910	11.8	86	0.00
75 T	alpha-Pinene	1.611	1.440	10.6	81	0.00
76 T	n-Propylbenzene	4.039	3.488	13.6	82	0.00

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\26\  
 Data File : 03261201.D  
 Acq On : 26 Mar 2012 8:12  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 26 10:52:52 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
77 T	3-Ethyltoluene	3.206	2.755	14.1	82	-0.01
78 T	4-Ethyltoluene	3.133	2.693	14.0	80	-0.01
79 T	1,3,5-Trimethylbenzene	2.600	2.250	13.5	82	-0.01
80 T	alpha-Methylstyrene	1.436	1.270	11.6	80	-0.01
81 T	2-Ethyltoluene	3.254	2.799	14.0	81	-0.01
82 T	1,2,4-Trimethylbenzene	2.664	2.317	13.0	82	-0.02
83 T	n-Decane	1.699	1.524	10.3	83	-0.01
84 T	Benzyl Chloride	1.948	2.075	-6.5	89	-0.02
85 T	1,3-Dichlorobenzene	1.607	1.267	21.2	80	-0.01
86 T	1,4-Dichlorobenzene	1.649	1.269	23.0	79	-0.02
87 T	sec-Butylbenzene	3.571	3.090	13.5	82	-0.01
88 T	4-Isopropyltoluene (p-Cymen)	3.554	3.089	13.1	81	-0.01
89 T	1,2,3-Trimethylbenzene	2.691	2.373	11.8	82	-0.01
90 T	1,2-Dichlorobenzene	1.531	1.237	19.2	79	-0.01
91 T	d-Limonene	1.067	0.991	7.1	82	-0.01
92 T	1,2-Dibromo-3-Chloropropane	0.517	0.479	7.4	81	-0.01
93 T	n-Undecane	1.678	1.541	8.2	84	0.00
94 T	1,2,4-Trichlorobenzene	1.122	0.944	15.9	78	-0.01
95 T	Naphthalene	3.881	3.423	11.8	81	-0.01
96 T	n-Dodecane	1.613	1.536	4.8	86	0.00
97 T	Hexachlorobutadiene	0.704	0.581	17.5	77	0.00
98 T	Cyclohexanone	0.986	0.965	2.1	89	-0.02
99 T	tert-Butylbenzene	2.723	2.359	13.4	82	-0.01
100 T	n-Butylbenzene	2.817	2.442	13.3	82	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*EM 3/26/12*

Response Factor Report MS07

Method : J:\Ms07\METHODS\X7021712B.M (RTE Integrator)  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
 1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	Avg	%RSD
1) I Bromochloromethan					ISTD						
2) T Dichlorodifluorom	3.784	3.527	3.467	3.822	3.279	2.986	2.875	3.155	2.920	3.313	10.83
3) T Chloromethane		1.470	0.956	1.190	0.904	0.778	0.569	0.897	0.827	0.949	28.81
4) T Vinyl Chloride	2.802	2.749	2.588	3.044	2.597	2.406	2.314	2.588	2.446	2.615	8.57
5) T Bromomethane	1.894	1.496	1.211	1.613	1.335	1.208	1.105	1.433	1.367	1.407	17.11
6) T Chloroethane	1.511	1.416	1.357	1.517	1.295	1.237	1.195	1.312	1.224	1.340	8.92
7) T Acetone				1.548	1.214	1.122	1.080	1.193	1.090	1.208	14.52
8) T Trichlorofluorome	3.051	2.887	2.831	3.145	2.692	2.565	2.447	2.691	2.536	2.761	8.59
9) T 1,1-Dichloroethen	1.591	1.440	1.366	1.547	1.322	1.229	1.214	1.385	1.295	1.376	9.52
10) T Methylene Chlorid		2.134	1.795	1.866	1.536	1.426	1.404	1.494	1.411	1.633	16.43
11) T Trichlorotrifluor	1.408	1.325	1.295	1.376	1.169	1.081	1.062	1.170	1.097	1.220	10.87
12) T trans-1,2-Dichlor	1.972	1.710	1.589	1.762	1.505	1.406	1.393	1.512	1.439	1.588	12.18
13) T 1,1-Dichloroethan	3.265	3.020	3.016	3.383	2.925	2.756	2.681	2.937	2.841	2.980	7.60
14) T Methyl tert-Butyl	5.168	4.740	4.762	5.045	4.335	4.137	4.166	4.690	4.471	4.613	7.90
15) T cis-1,2-Dichloroe	1.885	1.626	1.630	1.776	1.524	1.434	1.388	1.575	1.482	1.591	10.09
16) T Chloroform			3.041	3.257	2.586	2.390	2.280	2.524	2.374	2.636	14.03
17) S 1,2-Dichloroethan	2.000	2.002	2.032	2.018	1.989	1.973	1.960	1.938	1.932	1.983	1.75
18) T 1,2-Dichloroethan	2.582	2.460	2.375	2.604	2.222	2.090	1.988	2.222	2.094	2.293	9.73
19) T 1,1,1-Trichloroet	2.587	2.431	2.340	2.573	2.191	2.059	2.007	2.278	2.150	2.291	9.18
20) T Benzene			7.536	7.534	6.055	5.665	5.395	5.903	5.468	6.222	14.88
21) T Carbon Tetrachlor	2.280	1.842	1.769	1.950	1.676	1.579	1.549	1.771	1.692	1.790	12.41
22) I 1,4-Difluorobenze					ISTD						
23) T 1,2-Dichloropropa	0.417	0.397	0.373	0.409	0.354	0.331	0.319	0.358	0.336	0.366	9.68
24) T Bromodichlorometh	0.491	0.464	0.436	0.489	0.420	0.398	0.394	0.447	0.416	0.439	8.26
25) T Trichloroethene	0.418	0.379	0.349	0.390	0.324	0.305	0.287	0.326	0.307	0.343	12.96
26) T 1,4-Dioxane	0.324	0.290	0.270	0.278	0.242	0.231	0.227	0.258	0.243	0.263	11.99
27) T cis-1,3-Dichlorop	0.567	0.523	0.508	0.540	0.499	0.484	0.488	0.559	0.522	0.521	5.67
28) T trans-1,3-Dichlor	0.457	0.434	0.431	0.444	0.419	0.409	0.417	0.505	0.482	0.444	7.22

75 of 88

(#) = Out of Range ### Number of calibration levels exceeded format ###  
 X7021712B.M Tue Mar 27 13:12:09 2012



Response Factor Report MS07

Method : J:\Ms07\METHODS\X7021712B.M (RTE Integrator)  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
 1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	Avg	%RSD
29) T 1,1,2-Trichloroet	0.315	0.296	0.286	0.305	0.263	0.248	0.242	0.274	0.257	0.276	9.32
30) S Toluene-d8 (SS2)	1.053	1.071	1.054	1.047	1.051	1.052	1.071	1.060	1.049	1.056	0.84
31) T Toluene	1.782	1.541	1.420	1.485	1.269	1.198	1.177	1.307	1.186	1.374	14.76
32) T 1,2-Dibromoethane	0.379	0.362	0.345	0.362	0.323	0.307	0.304	0.351	0.331	0.340	7.65
33) T Tetrachloroethene	0.423	0.394	0.371	0.400	0.341	0.320	0.311	0.347	0.326	0.359	11.03
-----ISTD-----											
34) I Chlorobenzene-d5											
35) T Chlorobenzene	4.581	4.172	4.017	4.308	3.688	3.469	3.311	3.705	3.345	3.844	11.68
36) T Ethylbenzene	7.604	7.082	6.707	7.235	6.297	6.070	5.915	6.682	5.878	6.608	9.30
37) T m,p-Xylene	5.861	5.347	5.132	5.536	4.930	4.777	4.719	5.268	4.558	5.125	8.24
38) T o-Xylene	6.505	5.826	5.526	5.889	5.188	4.995	4.950	5.594	4.932	5.490	9.67
39) T 1,1,2,2-Tetrachlo	2.655	2.674	2.710	2.646	2.437	2.371	2.298	2.761	2.498	2.561	6.40
40) S Bromofluorobenzene	1.712	1.701	1.695	1.697	1.683	1.680	1.670	1.642	1.593	1.675	2.20
41) T 1,3-Dichlorobenze	3.304	3.261	3.222	3.204	2.835	2.719	2.608	3.138	2.846	3.015	8.71
42) T 1,4-Dichlorobenze	3.470	3.303	3.208	3.208	2.803	2.671	2.579	3.174	2.891	3.034	10.13
43) T 1,2-Dichlorobenze	3.271	3.123	3.112	3.026	2.662	2.536	2.418	3.053	2.772	2.886	10.31
44) T 1,2,4-Trichlorobe	2.184	2.118	2.169	1.753	1.641	1.608	1.575	2.139	1.978	1.907	13.66
45) T Naphthalene	6.857	6.355	6.704	5.143	5.107	5.172	5.297	8.126	7.328	6.232	17.82
46) T Hexachlorobutadie	1.486	1.423	1.403	1.217	1.055	1.013	0.975	1.227	1.145	1.216	15.40

(#) = Out of Range ### Number of calibration levels exceeded format ###  
 X7021712B.M Tue Mar 27 13:12:09 2012

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_03\27\  
 Data File : 03271203.D  
 Acq On : 27 Mar 2012 12:37 pm  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00069  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Mar 27 13:11:37 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane (IS1)	1.000	1.000	0.0	94	0.00
2 T	Dichlorodifluoromethane (CF)	3.313	2.627	20.7	76	0.00
3 T	Chloromethane	0.949	0.752	20.8	78	0.00
4 T	Vinyl Chloride	2.615	2.114	19.2	77	0.00
5 T	Bromomethane	1.407	1.229	12.7	87	0.00
6 T	Chloroethane	1.340	1.087	18.9	79	0.00
7 T	Acetone	1.208	1.101	8.9	86	0.00
8 T	Trichlorofluoromethane	2.761	2.220	19.6	78	0.00
9 T	1,1-Dichloroethene	1.376	1.159	15.8	83	0.00
10 T	Methylene Chloride	1.633	1.304	20.1	80	0.00
11 T	Trichlorotrifluoroethane	1.220	1.044	14.4	84	0.00
12 T	trans-1,2-Dichloroethene	1.588	1.286	19.0	81	0.00
13 T	1,1-Dichloroethane	2.980	2.440	18.1	79	0.00
14 T	Methyl tert-Butyl Ether	4.613	3.724	19.3	81	0.00
15 T	cis-1,2-Dichloroethene	1.591	1.352	15.0	84	0.00
16 T	Chloroform	2.636	2.203	16.4	80	0.00
17 S	1,2-Dichloroethane-d4 (SS1)	1.983	1.936	2.4	92	0.00
18 T	1,2-Dichloroethane	2.293	1.762	23.2	75	0.00
19 T	1,1,1-Trichloroethane	2.291	1.897	17.2	82	0.00
20 T	Benzene	6.222	5.363	13.8	84	0.00
21 T	Carbon Tetrachloride	1.790	1.447	19.2	82	0.00
22 I	1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	103	0.00
23 T	1,2-Dichloropropane	0.366	0.276	24.6	80	0.00
24 T	Bromodichloromethane	0.439	0.326	25.7	80	0.00
25 T	Trichloroethene	0.343	0.263	23.3	84	0.00
26 T	1,4-Dioxane	0.263	0.203	22.8	87	0.00
27 T	cis-1,3-Dichloropropene	0.521	0.407	21.9	84	0.00
28 T	trans-1,3-Dichloropropene	0.444	0.346	22.1	85	0.00
29 T	1,1,2-Trichloroethane	0.276	0.213	22.8	83	0.00
30 S	Toluene-d8 (SS2)	1.056	1.183	-12.0	116	0.00
31 T	Toluene	1.374	1.058	23.0	86	0.00
32 T	1,2-Dibromoethane	0.340	0.270	20.6	86	0.00
33 T	Tetrachloroethene	0.359	0.293	18.4	89	0.00
34 I	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	103	0.00
35 T	Chlorobenzene	3.844	3.159	17.8	88	0.00
36 T	Ethylbenzene	6.608	5.258	20.4	86	0.00
37 T	m,p-Xylene	5.125	4.158	18.9	87	0.00

*KR3/2/11*

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_03\27\  
 Data File : 03271203.D  
 Acq On : 27 Mar 2012 12:37 pm  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00069  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Mar 27 13:11:37 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
38 T	o-Xylene	5.490	4.453	18.9	89	0.00
39 T	1,1,2,2-Tetrachloroethane	2.561	2.083	18.7	88	0.00
40 S	Bromofluorobenzene (SS3)	1.675	1.976	-18.0	121	0.00
41 T	1,3-Dichlorobenzene	3.015	2.675	11.3	97	0.00
42 T	1,4-Dichlorobenzene	3.034	2.637	13.1	97	0.00
43 T	1,2-Dichlorobenzene	2.886	2.578	10.7	100	0.00
44 T	1,2,4-Trichlorobenzene	1.907	1.894	0.7	119	0.00
45 T	Naphthalene	6.232	6.662	-6.9	135	0.00
46 T	Hexachlorobutadiene	1.216	1.119	8.0	109	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*KR 3/27/12*

## LABORATORY REPORT

April 3, 2012

Christopher Cote  
Montana DEQ  
1100 N. Last Chance Gulch  
Helena, MT 59601

### RE: Background Indoor Air Study

Dear Christopher:

Enclosed are the results of the samples submitted to our laboratory on March 20, 2012. For your reference, these analyses have been assigned our service request number P1201074.

All analyses were performed according to our laboratory's NELAP and DoD-ELAP-approved quality assurance program. The test results meet requirements of the current NELAP and DoD-ELAP standards, where applicable, and except as noted in the laboratory case narrative provided. For a specific list of NELAP and DoD-ELAP-accredited analytes, refer to the certifications section at [www.caslab.com](http://www.caslab.com). Results are intended to be considered in their entirety and apply only to the samples analyzed and reported herein.

Columbia Analytical Services, Inc. is certified by the California Department of Health Services, NELAP Laboratory Certificate No. 02115CA; Arizona Department of Health Services, Certificate No. AZ0694; Florida Department of Health, NELAP Certification E871020; New Jersey Department of Environmental Protection, NELAP Laboratory Certification ID #CA009; New York State Department of Health, NELAP NY Lab ID No: 11221; Oregon Environmental Laboratory Accreditation Program, NELAP ID: CA20007; The American Industrial Hygiene Association, Laboratory #101661; United States Department of Defense Environmental Laboratory Accreditation Program (DoD-ELAP), Certificate No. L11-203; Pennsylvania Registration No. 68-03307; TX Commission of Environmental Quality, NELAP ID T104704413-11-2; Minnesota Department of Health, NELAP Certificate No. 362188; Washington State Department of Ecology, ELAP Lab ID: C946, State of Utah Department of Health, NELAP Certificate No. CA015272011-1; Los Angeles Department of Building and Safety, Approval No: TA00001. Each of the certifications listed above have an explicit Scope of Accreditation that applies to specific matrices/methods/analytes; therefore, please contact me for information corresponding to a particular certification.

If you have any questions, please call me at (805) 526-7161.

Respectfully submitted,

**Columbia Analytical Services, Inc.**

For Kate Aguilera  
Project Manager

Client: Montana DEQ  
Project: Background Indoor Air Study

CAS Project No: P1201074

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## CASE NARRATIVE

The samples were received intact under chain of custody on March 20, 2012 and were stored in accordance with the analytical method requirements. Please refer to the sample acceptance check form for additional information. The results reported herein are applicable only to the condition of the samples at the time of sample receipt.

### Air-Phase Petroleum Hydrocarbons (APH) Analysis

The samples were analyzed for total aliphatic and aromatic gasoline range hydrocarbons by gas chromatography/mass spectrometry according to the Method for the Determination of Air-Phase Petroleum Hydrocarbons (APH), Massachusetts Department of Environmental Protection, Revision 1, December, 2009.

### Volatile Organic Compound Analysis

The samples were also analyzed in Scan and SIM mode for selected volatile organic compounds in accordance with EPA Method TO-15 from the Compendium of Methods for the Determination of Toxic Organic Compounds in Ambient Air, Second Edition (EPA/625/R-96/010b), January, 1999. The analytical system was comprised of a gas chromatograph / mass spectrometer (GC/MS) interfaced to a whole-air preconcentrator.

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*The results of analyses are given in the attached laboratory report. All results are intended to be considered in their entirety, and Columbia Analytical Services, Inc. (CAS) is not responsible for utilization of less than the complete report.*

*Use of Columbia Analytical Services, Inc. (CAS) Name. Client shall not use CAS's name or trademark in any marketing or reporting materials, press releases or in any other manner ("Materials") whatsoever and shall not attribute to CAS any test result, tolerance or specification derived from CAS's data ("Attribution") without CAS's prior written consent, which may be withheld by CAS for any reason in its sole discretion. To request CAS's consent, Client shall provide copies of the proposed Materials or Attribution and describe in writing Client's proposed use of such Materials or Attribution. If CAS has not provided written approval of the Materials or Attribution within ten (10) days of receipt from Client, Client's request to use CAS's name or trademark in any Materials or Attribution shall be deemed denied. CAS may, in its discretion, reasonably charge Client for its time in reviewing Materials or Attribution requests. Client acknowledges and agrees that the unauthorized use of CAS's name or trademark may cause CAS to incur irreparable harm for which the recovery of money damages will be inadequate. Accordingly, Client acknowledges and agrees that a violation shall justify preliminary injunctive relief. For questions contact the laboratory.*

DETAIL SUMMARY REPORT

Client: Montana DEQ  
 Project ID: Background Indoor Air Study

Service Request: P1201074

Date Received: 3/20/2012  
 Time Received: 10:05

Client Sample ID	Lab Code	Matrix	Date Collected	Time Collected	Container ID	Pi1 (psig)	Pf1 (psig)	MA APH 1.0 - MA VOC PH Can	TO-15 - VOC Cans	TO-15 - VOC SIM
51-1F-030512	P1201074-001	Air	3/5/2012	20:30	AC01472	-4.07	3.59	X	X	X
04-1F-031312	P1201074-002	Air	3/13/2012	16:32	AC01828	-4.19	3.58	X	X	X

**Requested Turnaround Time in Business Days (Surcharges) please circle**  
 1 Day (100%) 2 Day (75%) 3 Day (50%) 4 Day (35%) 5 Day (25%) 10 Day-Standard

CAS Project No. **P1201074**

Company Name & Address (Reporting Information)							Project Name		CAS Contact:		Comments e.g. Actual Preservative or specific instructions	
Project Manager							Project Number	Whole Range Analysis Method				
Company Name & Address (Reporting Information) MT DEQ Helera AT 59620 P.O. # / Billing Information reference as: Background Indoor Air Study Bill to Chris's Cote, MDEQ							Project Number B0 Box 200901		CAS Contact: Whole Range		Comments e.g. Actual Preservative or specific instructions	
Phone 406 B44 5038 406-841-5050 Fax 406 B44 5038 406-841-5050							Project Name Helera AT 59620		CAS Contact: Whole Range			Comments e.g. Actual Preservative or specific instructions
Email Address for Result Reporting cote22@mt.deq.net							Project Name Background Indoor Air Study		CAS Contact: Whole Range		Comments e.g. Actual Preservative or specific instructions	
Client Sample ID	Laboratory ID Number	Date Collected	Time Collected	Canister ID (Bar code # - AC, SC, etc.)	Flow Controller ID (Bar code # - FC #)	Canister Start Pressure "Hg	Canister End Pressure "Hg/psig	Sample Volume				
51-1F-030512	①-417	3/5/12	830P						X	TO 15	PH	Analyze
04-1F-031312	②-417	3/13/12	432P	AL01020	FLA0003	255	9.0		X			Project specific list of compounds to project specific reporting limits

**Report Tier Levels - please select**

Tier I - Results (Default if not specified) \_\_\_\_\_  
 Tier II (Results + QC Summaries) \_\_\_\_\_  
 Tier III (Results + QC & Calibration Summaries) X  
 Tier IV (Data Validation Package) 10% Surcharge \_\_\_\_\_

EDD required Yes/No  
 Type: SVI Excel

Relinquished by: (Signature) [Signature] Date: 3/13/12 Time: 3:15P  
 Received by: (Signature) [Signature] Date: 3/13/12 Time: 3:15P

Cooler / Blank Temperature \_\_\_\_\_ °C

**Sample Acceptance Check Form**

Client: Montana DEQ Work order: P1201074

Project: Background Indoor Air Study

Sample(s) received on: 3/20/12 Date opened: 3/20/12 by: MZAMORA

*Note:* This form is used for all samples received by CAS. The use of this form for custody seals is strictly meant to indicate presence/absence and not as an indication of compliance or nonconformity. Thermal preservation and pH will only be evaluated either at the request of the client and/or as required by the method/SOP.

- |  | <u>Yes</u>                          | <u>No</u>                           | <u>N/A</u>                          |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| 1 Were <b>sample containers</b> properly marked with client sample ID?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 2 Container(s) <b>supplied by CAS</b> ?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 3 Did <b>sample containers</b> arrive in good condition?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 4 Were <b>chain-of-custody</b> papers used and filled out?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 5 Did <b>sample container labels</b> and/or tags agree with custody papers?                                      | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 6 Was <b>sample volume</b> received adequate for analysis?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 7 Are samples within specified holding times?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 8 Was proper <b>temperature</b> (thermal preservation) of cooler at receipt adhered to?                          | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 9 Was a <b>trip blank</b> received?  | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| 10 Were <b>custody seals</b> on outside of cooler/Box?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| Location of seal(s)? _____ Sealing Lid?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were signature and date included?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were seals intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were custody seals on outside of sample container?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| Location of seal(s)? _____ Sealing Lid?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were signature and date included?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were seals intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 11 Do containers have appropriate <b>preservation</b> , according to method/SOP or Client specified information? | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Is there a client indication that the submitted samples are <b>pH</b> preserved?                                 | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were <b>VOA vials</b> checked for presence/absence of air bubbles?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Does the client/method/SOP require that the analyst check the sample pH and <u>if necessary</u> alter it?        | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 12 <b>Tubes:</b> Are the tubes capped and intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Do they contain moisture?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 13 <b>Badges:</b> Are the badges properly capped and intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Are dual bed badges separated and individually capped and intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |

Lab Sample ID	Container Description	Required pH *	Received pH	Adjusted pH	VOA Headspace (Presence/Absence)	Receipt / Preservation Comments
P1201074-001.01	6.0 L Ambient Can					
P1201074-002.01	6.0 L Ambient Can					

Explain any discrepancies: (include lab sample ID numbers): \_\_\_\_\_



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 51-1F-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
 CAS Sample ID: P1201074-001

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Lusine Hakobyan/Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01472

Date Collected: 3/5/12  
 Date Received: 3/20/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.07      Final Pressure (psig): 3.59

Canister Dilution Factor: 1.72

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	180	34	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	40	17	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	8.6	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 04-1F-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
 CAS Sample ID: P1201074-002

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01828

Date Collected: 3/13/12  
 Date Received: 3/20/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.19      Final Pressure (psig): 3.58

Canister Dilution Factor: 1.74

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	35	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	<b>21</b>	17	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	8.7	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
 CAS Sample ID: P120324-MB

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	20	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	10	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	5.0	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
 CAS Sample ID: P120324-LCS

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: NA Liter(s)

Compound	Spike Amount ng	Result ng	% Recovery	CAS	Data Qualifier
				Acceptance Limits	
C5 - C8 Aliphatic Hydrocarbons	24.5	22.5	92	70-130	
C9 - C12 Aliphatic Hydrocarbons	25.0	23.3	93	70-130	
C9 - C10 Aromatic Hydrocarbons	48.5	43.0	89	70-130	

## RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 51-1F-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
 CAS Sample ID: P1201074-001

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Lusine Hakobyan/Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01472

**Date Collected:** 3/5/12  
**Date Received:** 3/20/12  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -4.07 Final Pressure (psig): 3.59

Canister Dilution Factor: 1.72

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	5.9	0.86	3.4	0.50	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.1	0.86	0.42	0.17	
74-87-3	Chloromethane	ND	0.86	ND	0.42	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.86	ND	0.12	
106-99-0	1,3-Butadiene	ND	0.86	ND	0.39	
74-83-9	Bromomethane	ND	0.86	ND	0.22	
75-00-3	Chloroethane	ND	0.86	ND	0.33	
64-17-5	Ethanol	1,800	86	940	46	D
75-05-8	Acetonitrile	ND	0.86	ND	0.51	
107-02-8	Acrolein	3.8	3.4	1.7	1.5	
67-64-1	Acetone	76	8.6	32	3.6	
75-69-4	Trichlorofluoromethane	1.1	0.86	0.20	0.15	
67-63-0	2-Propanol (Isopropyl Alcohol)	34	1.7	14	0.70	
107-13-1	Acrylonitrile	ND	0.86	ND	0.40	
75-35-4	1,1-Dichloroethene	ND	0.86	ND	0.22	
75-09-2	Methylene Chloride	ND	0.86	ND	0.25	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.86	ND	0.27	
76-13-1	Trichlorotrifluoroethane	ND	0.86	ND	0.11	
75-15-0	Carbon Disulfide	ND	8.6	ND	2.8	
156-60-5	trans-1,2-Dichloroethene	ND	0.86	ND	0.22	
75-34-3	1,1-Dichloroethane	ND	0.86	ND	0.21	
1634-04-4	Methyl tert-Butyl Ether	ND	0.86	ND	0.24	
108-05-4	Vinyl Acetate	18	8.6	5.1	2.4	
78-93-3	2-Butanone (MEK)	ND	8.6	ND	2.9	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 51-1F-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
 CAS Sample ID: P1201074-001

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Lusine Hakobyan/Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01472

Date Collected: 3/5/12  
 Date Received: 3/20/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -4.07      Final Pressure (psig): 3.59

Canister Dilution Factor: 1.72

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
156-59-2	cis-1,2-Dichloroethene	ND	0.86	ND	0.22	
141-78-6	Ethyl Acetate	<b>24</b>	1.7	<b>6.6</b>	0.48	
110-54-3	n-Hexane	<b>6.8</b>	0.86	<b>1.9</b>	0.24	
67-66-3	Chloroform	ND	0.86	ND	0.18	
109-99-9	Tetrahydrofuran (THF)	ND	0.86	ND	0.29	
71-55-6	1,1,1-Trichloroethane	ND	0.86	ND	0.16	
56-23-5	Carbon Tetrachloride	ND	0.86	ND	0.14	
110-82-7	Cyclohexane	ND	1.7	ND	0.50	
78-87-5	1,2-Dichloropropane	ND	0.86	ND	0.19	
75-27-4	Bromodichloromethane	ND	0.86	ND	0.13	
123-91-1	1,4-Dioxane	ND	0.86	ND	0.24	
80-62-6	Methyl Methacrylate	<b>3.2</b>	1.7	<b>0.79</b>	0.42	
142-82-5	n-Heptane	<b>2.7</b>	0.86	<b>0.66</b>	0.21	
10061-01-5	cis-1,3-Dichloropropene	ND	0.86	ND	0.19	
108-10-1	4-Methyl-2-pentanone	ND	0.86	ND	0.21	
10061-02-6	trans-1,3-Dichloropropene	ND	0.86	ND	0.19	
108-88-3	Toluene	<b>29</b>	0.86	<b>7.7</b>	0.23	
591-78-6	2-Hexanone	<b>1.0</b>	0.86	<b>0.25</b>	0.21	
124-48-1	Dibromochloromethane	ND	0.86	ND	0.10	
106-93-4	1,2-Dibromoethane	ND	0.86	ND	0.11	
123-86-4	n-Butyl Acetate	<b>3.0</b>	0.86	<b>0.64</b>	0.18	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 51-1F-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
 CAS Sample ID: P1201074-001

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Lusine Hakobyan/Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01472

Date Collected: 3/5/12  
 Date Received: 3/20/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -4.07 Final Pressure (psig): 3.59

Canister Dilution Factor: 1.72

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
111-65-9	n-Octane	1.2	0.86	0.25	0.18	
108-90-7	Chlorobenzene	ND	0.86	ND	0.19	
179601-23-1	m,p-Xylenes	6.3	1.7	1.4	0.40	
75-25-2	Bromoform	ND	0.86	ND	0.083	
100-42-5	Styrene	2.1	0.86	0.49	0.20	
95-47-6	o-Xylene	2.0	0.86	0.45	0.20	
111-84-2	n-Nonane	ND	0.86	ND	0.16	
98-82-8	Cumene	ND	0.86	ND	0.18	
80-56-8	alpha-Pinene	10	0.86	1.8	0.15	
103-65-1	n-Propylbenzene	ND	0.86	ND	0.18	
622-96-8	4-Ethyltoluene	ND	0.86	ND	0.18	
108-67-8	1,3,5-Trimethylbenzene	ND	0.86	ND	0.18	
95-63-6	1,2,4-Trimethylbenzene	1.4	0.86	0.28	0.18	
100-44-7	Benzyl Chloride	ND	0.86	ND	0.17	
541-73-1	1,3-Dichlorobenzene	ND	0.86	ND	0.14	
106-46-7	1,4-Dichlorobenzene	ND	0.86	ND	0.14	
95-50-1	1,2-Dichlorobenzene	ND	0.86	ND	0.14	
5989-27-5	d-Limonene	160	0.86	29	0.15	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.86	ND	0.089	
120-82-1	1,2,4-Trichlorobenzene	ND	0.86	ND	0.12	
87-68-3	Hexachlorobutadiene	ND	0.86	ND	0.081	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 04-1F-031312  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201074  
**CAS Sample ID:** P1201074-002

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01828

**Date Collected:** 3/13/12  
**Date Received:** 3/20/12  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -4.19      Final Pressure (psig): 3.58

Canister Dilution Factor: 1.74

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	7.4	0.87	4.3	0.51	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.1	0.87	0.42	0.18	
74-87-3	Chloromethane	ND	0.87	ND	0.42	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.87	ND	0.12	
106-99-0	1,3-Butadiene	ND	0.87	ND	0.39	
74-83-9	Bromomethane	ND	0.87	ND	0.22	
75-00-3	Chloroethane	ND	0.87	ND	0.33	
64-17-5	Ethanol	200	8.7	110	4.6	
75-05-8	Acetonitrile	ND	0.87	ND	0.52	
107-02-8	Acrolein	ND	3.5	ND	1.5	
67-64-1	Acetone	24	8.7	10	3.7	
75-69-4	Trichlorofluoromethane	1.1	0.87	0.19	0.15	
67-63-0	2-Propanol (Isopropyl Alcohol)	45	1.7	18	0.71	
107-13-1	Acrylonitrile	ND	0.87	ND	0.40	
75-35-4	1,1-Dichloroethene	ND	0.87	ND	0.22	
75-09-2	Methylene Chloride	ND	0.87	ND	0.25	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.87	ND	0.28	
76-13-1	Trichlorotrifluoroethane	ND	0.87	ND	0.11	
75-15-0	Carbon Disulfide	ND	8.7	ND	2.8	
156-60-5	trans-1,2-Dichloroethene	ND	0.87	ND	0.22	
75-34-3	1,1-Dichloroethane	ND	0.87	ND	0.22	
1634-04-4	Methyl tert-Butyl Ether	ND	0.87	ND	0.24	
108-05-4	Vinyl Acetate	ND	8.7	ND	2.5	
78-93-3	2-Butanone (MEK)	ND	8.7	ND	3.0	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



## RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 04-1F-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
 CAS Sample ID: P1201074-002

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01828

**Date Collected:** 3/13/12  
**Date Received:** 3/20/12  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -4.19      Final Pressure (psig): 3.58

Canister Dilution Factor: 1.74

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
156-59-2	cis-1,2-Dichloroethene	ND	0.87	ND	0.22	
141-78-6	Ethyl Acetate	3.5	1.7	0.98	0.48	
110-54-3	n-Hexane	ND	0.87	ND	0.25	
67-66-3	Chloroform	ND	0.87	ND	0.18	
109-99-9	Tetrahydrofuran (THF)	ND	0.87	ND	0.30	
71-55-6	1,1,1-Trichloroethane	ND	0.87	ND	0.16	
56-23-5	Carbon Tetrachloride	ND	0.87	ND	0.14	
110-82-7	Cyclohexane	ND	1.7	ND	0.51	
78-87-5	1,2-Dichloropropane	ND	0.87	ND	0.19	
75-27-4	Bromodichloromethane	ND	0.87	ND	0.13	
123-91-1	1,4-Dioxane	ND	0.87	ND	0.24	
80-62-6	Methyl Methacrylate	ND	1.7	ND	0.43	
142-82-5	n-Heptane	ND	0.87	ND	0.21	
10061-01-5	cis-1,3-Dichloropropene	ND	0.87	ND	0.19	
108-10-1	4-Methyl-2-pentanone	ND	0.87	ND	0.21	
10061-02-6	trans-1,3-Dichloropropene	ND	0.87	ND	0.19	
108-88-3	Toluene	ND	0.87	ND	0.23	
591-78-6	2-Hexanone	ND	0.87	ND	0.21	
124-48-1	Dibromochloromethane	ND	0.87	ND	0.10	
106-93-4	1,2-Dibromoethane	ND	0.87	ND	0.11	
123-86-4	n-Butyl Acetate	ND	0.87	ND	0.18	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 04-1F-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
 CAS Sample ID: P1201074-002

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01828

Date Collected: 3/13/12  
 Date Received: 3/20/12  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.19 Final Pressure (psig): 3.58

Canister Dilution Factor: 1.74

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
111-65-9	n-Octane	ND	0.87	ND	0.19	
108-90-7	Chlorobenzene	ND	0.87	ND	0.19	
179601-23-1	m,p-Xylenes	ND	1.7	ND	0.40	
75-25-2	Bromoform	ND	0.87	ND	0.084	
100-42-5	Styrene	ND	0.87	ND	0.20	
95-47-6	o-Xylene	ND	0.87	ND	0.20	
111-84-2	n-Nonane	ND	0.87	ND	0.17	
98-82-8	Cumene	ND	0.87	ND	0.18	
80-56-8	alpha-Pinene	<b>1.1</b>	0.87	<b>0.20</b>	0.16	
103-65-1	n-Propylbenzene	ND	0.87	ND	0.18	
622-96-8	4-Ethyltoluene	ND	0.87	ND	0.18	
108-67-8	1,3,5-Trimethylbenzene	ND	0.87	ND	0.18	
95-63-6	1,2,4-Trimethylbenzene	ND	0.87	ND	0.18	
100-44-7	Benzyl Chloride	ND	0.87	ND	0.17	
541-73-1	1,3-Dichlorobenzene	ND	0.87	ND	0.14	
106-46-7	1,4-Dichlorobenzene	<b>36</b>	0.87	<b>6.0</b>	0.14	
95-50-1	1,2-Dichlorobenzene	ND	0.87	ND	0.14	
5989-27-5	d-Limonene	<b>33</b>	0.87	<b>6.0</b>	0.16	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.87	ND	0.090	
120-82-1	1,2,4-Trichlorobenzene	ND	0.87	ND	0.12	
87-68-3	Hexachlorobutadiene	ND	0.87	ND	0.082	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
 CAS Sample ID: P120324-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/24/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	ND	0.50	ND	0.29	
75-71-8	Dichlorodifluoromethane (CFC 12)	ND	0.50	ND	0.10	
74-87-3	Chloromethane	ND	0.50	ND	0.24	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.50	ND	0.072	
106-99-0	1,3-Butadiene	ND	0.50	ND	0.23	
74-83-9	Bromomethane	ND	0.50	ND	0.13	
75-00-3	Chloroethane	ND	0.50	ND	0.19	
64-17-5	Ethanol	ND	5.0	ND	2.7	
75-05-8	Acetonitrile	ND	0.50	ND	0.30	
107-02-8	Acrolein	ND	2.0	ND	0.87	
67-64-1	Acetone	ND	5.0	ND	2.1	
75-69-4	Trichlorofluoromethane	ND	0.50	ND	0.089	
67-63-0	2-Propanol (Isopropyl Alcohol)	ND	1.0	ND	0.41	
107-13-1	Acrylonitrile	ND	0.50	ND	0.23	
75-35-4	1,1-Dichloroethene	ND	0.50	ND	0.13	
75-09-2	Methylene Chloride	ND	0.50	ND	0.14	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.50	ND	0.16	
76-13-1	Trichlorotrifluoroethane	ND	0.50	ND	0.065	
75-15-0	Carbon Disulfide	ND	5.0	ND	1.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.50	ND	0.13	
75-34-3	1,1-Dichloroethane	ND	0.50	ND	0.12	
1634-04-4	Methyl tert-Butyl Ether	ND	0.50	ND	0.14	
108-05-4	Vinyl Acetate	ND	5.0	ND	1.4	
78-93-3	2-Butanone (MEK)	ND	5.0	ND	1.7	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
CAS Sample ID: P120324-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/24/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
156-59-2	cis-1,2-Dichloroethene	ND	0.50	ND	0.13	
141-78-6	Ethyl Acetate	ND	1.0	ND	0.28	
110-54-3	n-Hexane	ND	0.50	ND	0.14	
67-66-3	Chloroform	ND	0.50	ND	0.10	
109-99-9	Tetrahydrofuran (THF)	ND	0.50	ND	0.17	
71-55-6	1,1,1-Trichloroethane	ND	0.50	ND	0.092	
56-23-5	Carbon Tetrachloride	ND	0.50	ND	0.080	
110-82-7	Cyclohexane	ND	1.0	ND	0.29	
78-87-5	1,2-Dichloropropane	ND	0.50	ND	0.11	
75-27-4	Bromodichloromethane	ND	0.50	ND	0.075	
123-91-1	1,4-Dioxane	ND	0.50	ND	0.14	
80-62-6	Methyl Methacrylate	ND	1.0	ND	0.24	
142-82-5	n-Heptane	ND	0.50	ND	0.12	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-10-1	4-Methyl-2-pentanone	ND	0.50	ND	0.12	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-88-3	Toluene	ND	0.50	ND	0.13	
591-78-6	2-Hexanone	ND	0.50	ND	0.12	
124-48-1	Dibromochloromethane	ND	0.50	ND	0.059	
106-93-4	1,2-Dibromoethane	ND	0.50	ND	0.065	
123-86-4	n-Butyl Acetate	ND	0.50	ND	0.11	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
 CAS Sample ID: P120324-MB

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
111-65-9	n-Octane	ND	0.50	ND	0.11	
108-90-7	Chlorobenzene	ND	0.50	ND	0.11	
179601-23-1	m,p-Xylenes	ND	1.0	ND	0.23	
75-25-2	Bromoform	ND	0.50	ND	0.048	
100-42-5	Styrene	ND	0.50	ND	0.12	
95-47-6	o-Xylene	ND	0.50	ND	0.12	
111-84-2	n-Nonane	ND	0.50	ND	0.095	
98-82-8	Cumene	ND	0.50	ND	0.10	
80-56-8	alpha-Pinene	ND	0.50	ND	0.090	
103-65-1	n-Propylbenzene	ND	0.50	ND	0.10	
622-96-8	4-Ethyltoluene	ND	0.50	ND	0.10	
108-67-8	1,3,5-Trimethylbenzene	ND	0.50	ND	0.10	
95-63-6	1,2,4-Trimethylbenzene	ND	0.50	ND	0.10	
100-44-7	Benzyl Chloride	ND	0.50	ND	0.097	
541-73-1	1,3-Dichlorobenzene	ND	0.50	ND	0.083	
106-46-7	1,4-Dichlorobenzene	ND	0.50	ND	0.083	
95-50-1	1,2-Dichlorobenzene	ND	0.50	ND	0.083	
5989-27-5	d-Limonene	ND	0.50	ND	0.090	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.50	ND	0.052	
120-82-1	1,2,4-Trichlorobenzene	ND	0.50	ND	0.067	
87-68-3	Hexachlorobutadiene	ND	0.50	ND	0.047	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister(s)  
 Test Notes:

Date(s) Collected: 3/5 - 3/13/12  
 Date(s) Received: 3/20/12  
 Date(s) Analyzed: 3/24/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		Percent Recovered	Percent Recovered	Percent Recovered		
Method Blank	P120324-MB	102	100	99	70-130	
Lab Control Sample	P120324-LCS	102	100	100	70-130	
51-1F-030512	P1201074-001	103	99	98	70-130	
04-1F-031312	P1201074-002	102	99	98	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.

Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
 CAS Sample ID: P120324-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
115-07-1	Propene	198	183	92	54-137	
75-71-8	Dichlorodifluoromethane (CFC 12)	196	156	80	55-112	
74-87-3	Chloromethane	190	155	82	66-122	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	200	154	77	64-113	
106-99-0	1,3-Butadiene	204	182	89	74-142	
74-83-9	Bromomethane	194	160	82	72-124	
75-00-3	Chloroethane	196	157	80	69-115	
64-17-5	Ethanol	928	800	86	67-127	
75-05-8	Acetonitrile	194	178	92	63-126	
107-02-8	Acrolein	198	164	83	62-127	
67-64-1	Acetone	1,010	833	82	67-106	
75-69-4	Trichlorofluoromethane	202	161	80	66-121	
67-63-0	2-Propanol (Isopropyl Alcohol)	382	304	80	56-112	
107-13-1	Acrylonitrile	198	183	92	78-128	
75-35-4	1,1-Dichloroethene	212	181	85	74-116	
75-09-2	Methylene Chloride	206	168	82	69-103	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	208	199	96	76-142	
76-13-1	Trichlorotrifluoroethane	206	175	85	69-118	
75-15-0	Carbon Disulfide	208	169	81	71-112	
156-60-5	trans-1,2-Dichloroethene	196	172	88	73-121	
75-34-3	1,1-Dichloroethane	200	176	88	71-118	
1634-04-4	Methyl tert-Butyl Ether	198	177	89	72-115	
108-05-4	Vinyl Acetate	952	902	95	51-150	
78-93-3	2-Butanone (MEK)	206	187	91	80-130	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
CAS Sample ID: P120324-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/24/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
156-59-2	cis-1,2-Dichloroethene	206	179	87	73-117	
141-78-6	Ethyl Acetate	398	357	90	79-126	
110-54-3	n-Hexane	198	163	82	68-109	
67-66-3	Chloroform	214	176	82	67-118	
109-99-9	Tetrahydrofuran (THF)	202	175	87	57-130	
71-55-6	1,1,1-Trichloroethane	198	169	85	70-116	
56-23-5	Carbon Tetrachloride	202	178	88	68-123	
110-82-7	Cyclohexane	390	323	83	73-111	
78-87-5	1,2-Dichloropropane	198	166	84	74-114	
75-27-4	Bromodichloromethane	198	172	87	75-120	
123-91-1	1,4-Dioxane	200	173	87	74-120	
80-62-6	Methyl Methacrylate	400	361	90	80-124	
142-82-5	n-Heptane	196	171	87	75-114	
10061-01-5	cis-1,3-Dichloropropene	188	167	89	79-120	
108-10-1	4-Methyl-2-pentanone	204	183	90	79-128	
10061-02-6	trans-1,3-Dichloropropene	210	193	92	83-131	
108-88-3	Toluene	202	167	83	64-115	
591-78-6	2-Hexanone	222	196	88	73-120	
124-48-1	Dibromochloromethane	206	186	90	72-137	
106-93-4	1,2-Dibromoethane	200	169	85	70-126	
123-86-4	n-Butyl Acetate	220	191	87	65-127	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.



## LABORATORY CONTROL SAMPLE SUMMARY

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
 CAS Sample ID: P120324-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/24/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
111-65-9	n-Octane	200	182	91	72-118	
108-90-7	Chlorobenzene	202	165	82	65-117	
179601-23-1	m,p-Xylenes	392	336	86	68-119	
75-25-2	Bromoform	208	177	85	79-150	
100-42-5	Styrene	200	177	89	74-127	
95-47-6	o-Xylene	194	164	85	68-118	
111-84-2	n-Nonane	196	168	86	72-116	
98-82-8	Cumene	190	161	85	68-119	
80-56-8	alpha-Pinene	186	163	88	70-123	
103-65-1	n-Propylbenzene	192	160	83	69-119	
622-96-8	4-Ethyltoluene	198	165	83	68-121	
108-67-8	1,3,5-Trimethylbenzene	200	169	85	67-118	
95-63-6	1,2,4-Trimethylbenzene	194	168	87	66-122	
100-44-7	Benzyl Chloride	200	192	96	73-144	
541-73-1	1,3-Dichlorobenzene	200	155	78	64-122	
106-46-7	1,4-Dichlorobenzene	206	153	74	65-125	
95-50-1	1,2-Dichlorobenzene	198	154	78	63-128	
5989-27-5	d-Limonene	200	182	91	72-126	
96-12-8	1,2-Dibromo-3-chloropropane	196	171	87	72-139	
120-82-1	1,2,4-Trichlorobenzene	196	157	80	65-139	
87-68-3	Hexachlorobutadiene	202	161	80	58-137	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 51-1F-030512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
CAS Sample ID: P1201074-001

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01472

Date Collected: 3/5/12  
Date Received: 3/20/12  
Date Analyzed: 3/28/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.07      Final Pressure (psig): 3.59

Canister Dilution Factor: 1.72

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.043	ND	0.017	
107-06-2	1,2-Dichloroethane	<b>0.20</b>	0.043	<b>0.050</b>	0.011	
71-43-2	Benzene	<b>3.5</b>	0.13	<b>1.1</b>	0.040	
79-01-6	Trichloroethene	<b>0.47</b>	0.043	<b>0.088</b>	0.0080	
79-00-5	1,1,2-Trichloroethane	ND	0.17	ND	0.032	
127-18-4	Tetrachloroethene	<b>0.13</b>	0.043	<b>0.020</b>	0.0063	
100-41-4	Ethylbenzene	<b>2.2</b>	0.17	<b>0.52</b>	0.040	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.043	ND	0.0063	
91-20-3	Naphthalene	<b>0.21</b>	0.17	<b>0.039</b>	0.033	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 04-1F-031312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
 CAS Sample ID: P1201074-002

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01828

Date Collected: 3/13/12  
 Date Received: 3/20/12  
 Date Analyzed: 3/28/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.19      Final Pressure (psig): 3.58

Canister Dilution Factor: 1.74

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.044	ND	0.017	
107-06-2	1,2-Dichloroethane	<b>0.13</b>	0.044	<b>0.033</b>	0.011	
71-43-2	Benzene	<b>0.38</b>	0.13	<b>0.12</b>	0.041	
79-01-6	Trichloroethene	ND	0.044	ND	0.0081	
79-00-5	1,1,2-Trichloroethane	ND	0.17	ND	0.032	
127-18-4	Tetrachloroethene	<b>0.28</b>	0.044	<b>0.041</b>	0.0064	
100-41-4	Ethylbenzene	ND	0.17	ND	0.040	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.044	ND	0.0063	
91-20-3	Naphthalene	<b>0.19</b>	0.17	<b>0.036</b>	0.033	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
 CAS Sample ID: P120328-MB

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/28/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result		MRL		Data Qualifier
		$\mu\text{g}/\text{m}^3$	$\mu\text{g}/\text{m}^3$	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.025	ND	0.0098	
107-06-2	1,2-Dichloroethane	ND	0.025	ND	0.0062	
71-43-2	Benzene	ND	0.075	ND	0.023	
79-01-6	Trichloroethene	ND	0.025	ND	0.0047	
79-00-5	1,1,2-Trichloroethane	ND	0.10	ND	0.018	
127-18-4	Tetrachloroethene	ND	0.025	ND	0.0037	
100-41-4	Ethylbenzene	ND	0.10	ND	0.023	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.025	ND	0.0036	
91-20-3	Naphthalene	ND	0.10	ND	0.019	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister(s)  
 Test Notes:

Date(s) Collected: 3/5 - 3/13/12  
 Date(s) Received: 3/20/12  
 Date(s) Analyzed: 3/28/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		% Recovered	% Recovered	% Recovered		
Method Blank	P120328-MB	100	113	113	70-130	
Lab Control Sample	P120328-LCS	100	112	116	70-130	
51-1F-030512	P1201074-001	98	111	115	70-130	
04-1F-031312	P1201074-002	98	111	119	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.

Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201074  
CAS Sample ID: P120328-LCS

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/28/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
75-01-4	Vinyl Chloride	3.88	3.19	82	56-127	
107-06-2	1,2-Dichloroethane	4.00	3.19	80	51-140	
71-43-2	Benzene	3.96	3.65	92	56-125	
79-01-6	Trichloroethene	3.88	3.08	79	51-127	
79-00-5	1,1,2-Trichloroethane	3.92	3.10	79	49-137	
127-18-4	Tetrachloroethene	3.68	3.06	83	58-134	
100-41-4	Ethylbenzene	3.96	3.26	82	56-126	
79-34-5	1,1,2,2-Tetrachloroethane	3.84	3.15	82	53-148	
91-20-3	Naphthalene	3.44	3.48	101	25-149	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

Response Factor Report MS08

Method Path : J:\MS08\Methods\  
 Method File : M8011412A.M  
 Title : Massachusetts APH  
 Last Update : Tue Jan 31 15:28:17 2012  
 Response Via : Initial Calibration

Calibration Files

0.5 =01141202.D 1 =01141203.D 5 =01141204.D 25 =01141205.D  
 50 =01141206.D 100 =01141207.D

Compound	0.5	1	5	25	50	100	Avg	%RSD
1) I Bromochloromethane...	-----ISTD-----							
2) T Bromochloromet...	5.258	5.274	5.259	5.250	5.256	5.271	5.261	0.17
3) UN Isopentane							0.000	-1.00
4) UN n-Hexane							0.000	-1.00
5) S 1,2-Dichloroet...	1.611	1.608	1.551	1.598	1.567	1.556	1.582	1.70
6) T 1,2-Dichloroet...	4.656	4.658	4.495	4.586	4.518	4.500	4.569	1.65
7) I 1,4-Difluorobenzen...	-----ISTD-----							
8) T 1,4-Difluorobe...	2.208	2.210	2.217	2.178	2.149	2.066	2.171	2.66
9) UN Cyclohexane							0.000	-1.00
10) UN 2,3-Dimethylpe...							0.000	-1.00
11) UN n-Heptane							0.000	-1.00
12) S Toluene-d8 (SS2)	1.109	1.103	1.110	1.107	1.109	1.099	1.106	0.38
13) T Toluene-d8 (TIC)	3.041	3.035	3.049	3.051	3.050	3.036	3.044	0.24
14) UN n-Octane							0.000	-1.00
15) T C5-C8 Aliphati...	2.542	1.269	0.255	0.051	0.026	0.013	0.693	E1 148.16
16) I Chlorobenzene-d5 (...)	-----ISTD-----							
17) T Chlorobenzene-...	2.907	2.900	2.906	2.916	2.929	2.947	2.917	0.60
18) UN 2,3-Dimethylhe...							0.000	-1.00
19) UN n-Nonane							0.000	-1.00
20) S p-Bromofluorob...	0.431	0.430	0.420	0.427	0.421	0.420	0.425	1.26
21) T p-Bromofluorob...	2.281	2.274	2.217	2.243	2.230	2.227	2.245	1.17
22) UN Isopropylbenzene							0.000	-1.00
23) UN 1-Methyl-3-eth...							0.000	-1.00
24) UN 1,3,5-Trimethy...							0.000	-1.00
25) UN n-Decane							0.000	-1.00
26) UN p-Isopropyltol...							0.000	-1.00
27) UN 1,2,3-Trimethy...							0.000	-1.00
28) UN Butylcyclohexane							0.000	-1.00
29) UN n-Undecane							0.000	-1.00
30) UN n-Dodecane							0.000	-1.00
31) T C9-C12 Aliphati...	1.899	0.947	0.185	0.037	0.019	0.009	0.516	E1 148.66
32) T C9-C10 Aromati...							0.000	-1.00
33) T C9-C10 Aromati...							0.000	-1.00
34) C5C8_TIC_1	1.741	0.869	0.174	0.035	0.017	0.009	0.474	E2 148.17
35) C5C8_TIC_2	1.741	0.869	0.174	0.035	0.017	0.009	0.474	E2 148.17
36) C5C8_TIC_3	1.741	0.869	0.174	0.035	0.017	0.009	0.474	E2 148.17
37) C5C8_TIC_4	1.741	0.869	0.174	0.035	0.017	0.009	0.474	E2 148.17
38) C9C12_TIC_1	1.140	0.569	0.111	0.022	0.011	0.006	0.310	E2 148.66
39) C9C12_TIC_2	1.140	0.569	0.111	0.022	0.011	0.006	0.310	E2 148.66
40) C9C12_TIC_3	1.140	0.569	0.111	0.022	0.011	0.006	0.310	E2 148.66
41) C9C12_TIC_4	1.140	0.569	0.111	0.022	0.011	0.006	0.310	E2 148.66
42) C9C10_TIC_1							0.000	-1.00
43) C9C10_TIC_2							0.000	-1.00

*Com* 1/31/12

Response Factor Report MS08

Method Path : J:\MS08\Methods\  
Method File : M8011412A.M  
Title : Massachusetts APH  
(#) = Out of Range

*Em 1/31/12*



Massachusetts APH  
Hydrocarbon Ranges

ICAL Method: M8011412A.M

ICAL Date: 1/14/12

Instrument ID: MS08

areas

masses

	areas				masses							
	0.5	1	5	25	50	100	0.5	1	5	25	50	100
<b>C5-C8 Aliphatics</b>												
Isopentane	60478	112606	488243	2480105	4710733	9222808	0.520	1.04	5.20	26.0	52.0	104
n-Hexane	59953	110631	507126	2577865	4931102	9678199	0.490	0.98	4.90	24.5	49.0	98
Cyclohexane	63203	115230	548910	2874726	5524402	10890911	0.500	1.00	5.00	25.0	50.0	100
2,3-Dimethylpentane	68648	127222	585183	2998017	5742104	11319597	0.520	1.04	5.20	26.0	52.0	104
n-Heptane	66143	120286	558074	2884817	5509801	10873046	0.490	0.98	4.90	24.5	49.0	98
n-Octane	76169	137829	632891	3286017	6205035	12092514	0.505	1.01	5.05	25.3	50.5	101
<b>area sum:</b>	<b>394594</b>	<b>723804</b>	<b>3320427</b>	<b>17101547</b>	<b>32623177</b>	<b>64077075</b>	<b>3.025</b>	<b>6.050</b>	<b>30.25</b>	<b>151.30</b>	<b>302.50</b>	<b>605.00</b>

**C9-C12 Aliphatics**

	0.5	1	5	25	50	100	0.5	1	5	25	50	100
2,3-Dimethylheptane	78780	147724	682518	3501487	6691428	13148840	0.485	0.97	4.85	24.3	48.5	97
n-Nonane	77464	140680	662221	3417128	6555165	12940625	0.485	0.97	4.85	24.3	48.5	97
n-Decane	80593	147724	705259	3651028	7021424	13870530	0.490	0.98	4.90	24.5	49.0	98
Butylcyclohexane	88915	166782	776903	4026931	7749936	15415340	0.495	0.99	4.95	24.8	49.5	99
n-Undecane	80867	152557	719618	3772331	7263077	14385881	0.480	0.96	4.80	24.0	48.0	96
n-Dodecane	70045	143746	687255	3673524	7115012	14041583	0.500	1.00	5.00	25.0	50.0	100
<b>area sum:</b>	<b>476664</b>	<b>899213</b>	<b>4233774</b>	<b>22042429</b>	<b>42396042</b>	<b>83802799</b>	<b>2.935</b>	<b>5.870</b>	<b>29.35</b>	<b>146.90</b>	<b>293.50</b>	<b>587.00</b>

**C9-C10 Aromatics**

	0.5	1	5	25	50	100	0.5	1	5	25	50	100
Isopropylbenzene	10658	19169	87439	448287	849124	1695737	0.485	0.97	4.85	24.3	48.5	97
3-Ethyltoluene	11805	20564	97823	504510	977020	1973178	0.470	0.94	4.70	23.5	47.0	94
1,3,5-Trimethylbenzene	15759	29260	134073	700133	1353305	2726630	0.490	0.98	4.90	24.5	49.0	98
1,2,3-Trimethylbenzene	17195	31328	143473	757946	1493373	3063059	0.495	0.99	4.95	24.8	49.5	99
p-isopropyltoluene	9490	18116	83401	440962	872799	1814314	0.460	0.92	4.60	23.0	46.0	92
<b>area sum:</b>	<b>64907</b>	<b>118437</b>	<b>546209</b>	<b>2851838</b>	<b>5545621</b>	<b>11272918</b>	<b>2.400</b>	<b>4.800</b>	<b>24.00</b>	<b>120.10</b>	<b>240.00</b>	<b>480.00</b>

*em* 1/31/12

Massachusetts APH  
Hydrocarbon Ranges

ICAL: M8011412A.M

ICAL Date: 1/14/12

Instrument ID: MS08

areas

Internal Standards (TIC)	0.5	1	5	25	50	100
Bromochloromethane (IS1)	1312793	1307309	1293586	1289283	1284194	1283817
1,4-Difluorobenzene (IS2)	2547079	2541159	2512498	2475255	2422119	2352159
Chlorobenzene-d5 (IS3)	2929189	2913374	2880258	2895080	2878314	2908396

Internal Standards (EIC)

Bromochloromethane (IS1)	249696	247899	245985	245585	244327	243561
1,4-Difluorobenzene (IS2)	1153657	1149947	1133233	1136223	1127203	1138730
Chlorobenzene-d5 (IS3)	1007702	1004638	991212	992925	982812	986950

Surrogates (TIC)

1,2-Dichloroethane-d4	1162497	1154617	1105630	1126201	1103958	1095945
Toluene-d8	3508018	3490493	3455382	3466688	3437986	3457323
p-Bromofluorobenzene	2298220	2284797	2197239	2227396	2192159	2197999

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RRFs

C5-C8 Aliphatics	0.5	1	5	25	50	100	RRF <sub>avg</sub>	%RSD
	2.8268	2.6009	2.4215	2.4870	2.3919	2.3252	2.509	7.24
C9-C12 Aliphatics	0.5	1	5	25	50	100	RRF <sub>avg</sub>	%RSD
	4.0291	3.8120	3.6383	3.7780	3.6744	3.6163	3.758	4.09
C9-C10 Aromatics	0.5	1	5	25	50	100	RRF <sub>avg</sub>	%RSD
	0.6709	0.6140	0.5740	0.5979	0.5878	0.5949	0.607	5.63

em 1/31/12

**Massachusetts APH**  
**Continuing Calibration Verification Check Sheet**

Data File Name: **03241202.D**  
 Data File Path: **J:\MS08\Data\2012\_03\24\**  
 Operator: **EM**  
 Date Acquired: **3/24/12 1:17**  
 Acq. Method File: **TO15.M**  
 Sample Name: **25ng TO-15/MAPH CCV STD**  
 Misc Info: **S25-03151201/S25-03161202**  
 Instrument Name: **MS08**

Enter RRFs from current ICAL!

	<u>Internal Standards</u>	<u>RT</u>	<u>Area</u>
7)	1,4-Difluorobenzene (IS2)	15.89	1097014
16)	Chlorobenzene-d5 (IS3)	21.69	957299

	<u>C5-C8 Aliphatics</u>	<u>RT</u>	<u>Area</u>	<u>RRF</u>	<u>ng</u>	<u>% D</u>	<u>LCL</u>	<u>UCL</u>	<u>Pass/Fai</u>
3)	Isopentane	8.19	2331141	2.445	147.5	-2.53	-30	30	Pass
4)	n-Hexane	13.09	2450350						
9)	Cyclohexane	15.82	2698497						
10)	2,3-Dimethylpentane	16.17	2863333	<b>Spike</b>	<b>ICAL</b>				
11)	n-Heptane	17.35	2745668	<u>Amt (ng)</u>	<u>RRF</u>				
14)	n-Octane	20.69	<u>3146714</u>	151.30	2.509				
			<b>16235703</b>						

	<u>C9-C12 Aliphatics</u>	<u>RT</u>	<u>Area</u>	<u>RRF</u>	<u>ng</u>	<u>% D</u>	<u>LCL</u>	<u>UCL</u>	<u>Pass/Fai</u>
18)	2,3-Dimethylheptane	22.31	3331254	3.712	145.1	-1.23	-30	30	Pass
19)	n-Nonane	23.30	3243877						
25)	n-Decane	25.27	3468097						
28)	Butylcyclohexane	25.99	3790882	<b>Spike</b>	<b>ICAL</b>				
29)	n-Undecane	26.77	3564701	<u>Amt (ng)</u>	<u>RRF</u>				
30)	n-Dodecane	28.00	<u>3481141</u>	146.90	3.758				
			<b>20879952</b>						

	<u>C9-C10 Aromatics</u>	<u>RT</u>	<u>Area</u>	<u>RRF</u>	<u>ng</u>	<u>% D</u>	<u>LCL</u>	<u>UCL</u>	<u>Pass/Fai</u>
22)	Isopropylbenzene	23.78	435563	0.589	116.6	-2.92	-30	30	Pass
23)	1-Methyl-3-ethylbenzene	24.52	480909						
24)	1,3,5-Trimethylbenzene	24.66	662082						
26)	p-Isopropyltoluene	25.68	417535	<b>Spike</b>	<b>ICAL</b>				
27)	1,2,3-Trimethylbenzene	25.68	<u>713901</u>	<u>Amt (ng)</u>	<u>RRF</u>				
			<b>2709990</b>	120.1	0.607				

EM 3/24/12

Response Factor Report MS08

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Mon Dec 12 09:45:45 2011  
 Response Via : Initial Calibration

Calibration Files

0.1 =12091118.D 0.2 =12091119.D 0.5 =12091120.D 1.0 =12091121.D 5.0 =12091122.D 25 =12091123.D  
 50 =12091124.D 100 =12091125.D

Compound	0.1	0.2	0.5	1.0	5.0	25	50	100	Avg	%RSD
1) IR Bromochloromethane...										
2) T Propene	2.487	1.965	1.952	1.692	2.194	2.303	2.185	2.036	2.102	11.62
3) T Dichlorodifluo...	3.545	2.805	2.940	2.748	2.432	2.361	2.278	2.199	2.664	16.73
4) T Chloromethane	3.290	2.811	2.875	2.512	2.176	2.286	1.977	1.664	2.449	21.62
5) T 1,2-Dichloro-1...	1.901	1.661	1.575	1.458	1.359	1.333	1.200	1.190	1.460	16.68
6) T Vinyl Chloride	2.906	2.590	2.631	2.374	2.145	2.129	1.950	1.939	2.333	15.07
7) T 1,3-Butadiene	2.145	1.788	1.790	1.718	1.581	1.689	1.621	1.616	1.743	10.33
8) T Bromomethane	1.687	1.328	1.349	1.212	1.094	1.142	1.111	1.092	1.252	16.21
9) T Chloroethane	1.671	1.259	1.333	1.193	1.101	1.091	1.046	1.044	1.217	17.29
10) T Ethanol	2.040	1.345	1.186	1.085	1.022	1.096	1.076	1.064	1.239	27.34
11) T Acetonitrile	4.365	2.973	2.723	2.546	2.480	2.562	2.482	2.468	2.825	22.85
12) T Acrolein	1.328	0.931	0.907	0.861	0.772	0.814	0.789	0.779	0.897	20.48
13) T Acetone	1.645	1.316	1.249	1.129	0.994	1.015	0.981	0.964	1.161	20.26
14) T Trichlorofluor...	2.776	2.385	2.496	2.272	2.097	2.135	2.042	2.011	2.277	11.59
15) T 2-Propanol (Is...	4.463	4.536	4.168	2.920	2.763	2.577	2.672	3.443	26.09	26.09
16) T Acrylonitrile	2.701	1.907	1.878	1.783	1.745	1.813	1.765	1.738	1.916	16.85
17) T 1,1-Dichloroet...	1.458	1.418	1.452	1.263	1.206	1.223	1.172	1.152	1.293	9.95
18) T 2-Methyl-2-Pro...	4.464	3.710	3.939	3.616	3.558	3.584			3.812	9.14
19) T Methylene Chlo...			1.958	1.578	1.294	1.273	1.220	1.201	1.421	20.88
20) T 3-Chloro-1-pro...	2.347	2.260	2.345	2.220	2.150	2.296	2.230	2.191	2.255	3.15
21) T Trichlorotrifl...	1.348	1.244	1.299	1.193	1.089	1.098	1.062	1.035	1.171	10.01
22) T Carbon Disulfide	6.210	5.678	5.057	4.654	4.699	4.488	4.424	5.030	13.40	13.40
23) T trans-1,2-Dich...	2.436	2.237	2.206	2.142	1.951	1.953	1.883	1.835	2.080	10.00
24) T 1,1-Dichloroet...	2.896	2.757	2.747	2.500	2.307	2.379	2.288	2.236	2.514	10.07
25) T Methyl tert-Bu...	5.128	4.611	4.650	4.307	4.119	4.185	4.042	3.984	4.378	8.93
26) T Vinyl Acetate	0.430	0.368	0.423	0.399	0.393	0.410	0.402	0.398	0.403	4.70
27) T 2-Butanone (MEK)	1.265	1.003	1.038	0.972	0.927	0.964	0.928	0.794	0.986	13.56
28) T cis-1,2-Dichlo...	2.442	2.056	2.101	1.938	1.826	1.836	1.764	1.720	1.960	12.05
29) T Diisopropyl Ether	1.446	1.223	1.323	1.262	1.171	1.188	1.153	1.148	1.239	8.26
30) T Ethyl Acetate	0.611	0.554	0.573	0.545	0.517	0.535	0.520	0.516	0.546	6.02
31) T n-Hexane	3.541	3.006	3.181	2.805	2.582	2.633	2.545	2.543	2.855	12.70

Response Factor Report MS08

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M

Title	EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)										
32) T	Chloroform	2.679	2.408	2.409	2.204	1.983	2.054	1.974	1.937	2.206	12.18
33) S	1,2-Dichloroet...	1.586	1.577	1.582	1.590	1.535	1.533	1.520	1.504	1.553	2.19
34) T	Tetrahydrofura...	1.274	0.932	1.100	0.938	0.910	0.879	0.830	0.821	0.961	15.96
35) T	Ethyl tert-But...	1.999	1.832	2.033	1.818	1.732	1.757	1.708	1.687	1.821	7.18
36) T	1,2-Dichloroet...	2.099	1.992	1.976	1.806	1.683	1.700	1.599	1.531	1.798	11.39
37) IR	1,4-Difluorobenzen...	-----ISTD-----									
38) T	1,1,1-Trichloro...	0.531	0.465	0.486	0.453	0.419	0.434	0.413	0.403	0.451	9.46
39) T	Isopropyl Acetate	0.227	0.205	0.220	0.206	0.194	0.201	0.192	0.188	0.204	6.71
40) T	1-Butanol		0.348	0.319	0.300	0.320	0.335	0.324	0.318	0.323	4.66
41) T	Benzene	1.633	1.385	1.359	1.227	1.136	1.158	1.101	1.070	1.258	15.11
42) T	Carbon Tetrach...	0.394	0.356	0.369	0.340	0.334	0.352	0.337	0.333	0.352	6.03
43) T	Cyclohexane	0.666	0.576	0.588	0.539	0.492	0.500	0.479	0.469	0.539	12.59
44) T	tert-Amyl Meth...	1.092	0.926	0.998	0.937	0.900	0.918	0.881	0.858	0.939	7.95
45) T	1,2-Dichloropr...	0.397	0.361	0.368	0.324	0.312	0.316	0.300	0.293	0.334	11.12
46) T	Bromodichlorom...	0.458	0.387	0.392	0.378	0.362	0.378	0.362	0.353	0.384	8.57
47) T	Trichloroethene	0.466	0.375	0.379	0.344	0.323	0.328	0.316	0.309	0.355	14.62
48) T	1,4-Dioxane	0.305	0.282	0.288	0.248	0.245	0.246	0.237	0.230	0.260	10.55
49) T	2,2,4-Trimethy...	1.964	1.629	1.686	1.546	1.447	1.461	1.387	1.337	1.557	12.98
50) T	Methyl Methacr...	0.156	0.139	0.150	0.135	0.136	0.142	0.136	0.135	0.141	5.53
51) T	n-Heptane	0.427	0.373	0.385	0.364	0.341	0.344	0.332	0.324	0.361	9.35
52) T	cis-1,3-Dichlo...	0.571	0.502	0.521	0.491	0.482	0.500	0.481	0.471	0.502	6.33
53) T	4-Methyl-2-pen...	0.359	0.295	0.308	0.303	0.297	0.303	0.289	0.284	0.305	7.68
54) T	trans-1,3-Dich...	0.476	0.413	0.443	0.421	0.429	0.453	0.437	0.427	0.437	4.59
55) T	1,1,2-Trichlor...	0.340	0.301	0.310	0.284	0.276	0.285	0.271	0.265	0.292	8.44
56) IR	Chlorobenzene-d5	-----ISTD-----									
57) S	Toluene-d8 (SS2)	2.305	2.306	2.305	2.309	2.306	2.300	2.310	2.304	2.306	0.14
58) T	Toluene	3.623	2.920	3.025	2.738	2.571	2.574	2.473	2.421	2.793	14.19
59) T	2-Hexanone	2.049	1.429	1.515	1.403	1.395	1.422	1.367	1.337	1.490	15.57
60) T	Dibromochlorom...	0.696	0.618	0.636	0.636	0.612	0.648	0.629	0.625	0.638	4.14
61) T	1,2-Dibromoethane	0.831	0.730	0.718	0.658	0.641	0.662	0.638	0.628	0.688	9.97
62) T	n-Butyl Acetate	2.338	1.850	1.857	1.721	1.704	1.765	1.737	1.737	1.839	11.40
63) T	n-Octane	0.785	0.652	0.714	0.643	0.602	0.611	0.580	0.572	0.645	11.28
64) T	Tetrachloroethene	1.090	0.904	0.904	0.826	0.767	0.779	0.756	0.755	0.848	13.63
65) T	Chlorobenzene	2.375	1.868	1.886	1.744	1.610	1.637	1.574	1.548	1.780	15.30
66) T	Ethylbenzene	3.846	3.171	3.188	2.956	2.803	2.848	2.743	2.680	3.030	12.51
67) T	m- & p-Xylenes	3.081	2.499	2.611	2.392	2.216	2.273	2.170	2.112	2.419	13.08
68) T	Bromoform	0.622	0.494	0.556	0.541	0.551	0.598	0.589	0.594	0.568	7.19
69) T	Styrene	2.478	1.895	1.937	1.800	1.735	1.803	1.726	1.675	1.881	13.64
70) T	o-Xylene	3.279	2.708	2.726	2.479	2.331	2.373	2.283	2.224	2.550	13.65

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M

Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

71) T	n-Nonane	2.005	1.753	1.847	1.675	1.581	1.597	1.518	1.470	1.681	10.69
72) T	1,1,2,2-Tetrac...	1.348	1.164	1.262	1.150	1.083	1.121	1.085	1.066	1.160	8.48
73) S	Bromofluoroben...	0.774	0.771	0.779	0.781	0.775	0.778	0.782	0.780	0.777	0.50
74) T	Cumene	4.522	3.495	3.524	3.169	2.962	3.002	2.884	2.831	3.299	16.98
75) T	alpha-Pinene	1.782	1.636	1.734	1.601	1.537	1.578	1.518	1.498	1.611	6.35
76) T	n-Propylbenzene	5.441	4.229	4.306	3.924	3.710	3.754	3.553	3.394	4.039	16.01
77) T	3-Ethyltoluene	4.091	3.373	3.461	3.038	2.964	2.981	2.893	2.845	3.206	13.14
78) T	4-Ethyltoluene	4.003	3.202	3.301	3.151	2.837	2.983	2.810	2.776	3.133	12.84
79) T	1,3,5-Trimethy...	3.324	2.662	2.723	2.537	2.378	2.426	2.344	2.403	2.600	12.43
80) T	alpha-Methylst...	1.742	1.386	1.449	1.371	1.390	1.402	1.354	1.399	1.436	8.80
81) T	2-Ethyltoluene	4.143	3.362	3.430	3.169	2.975	3.048	2.919	2.985	3.254	12.43
82) T	1,2,4-Trimethy...	3.356	2.663	2.791	2.551	2.433	2.513	2.454	2.548	2.664	11.37
83) T	n-Decane	1.969	1.702	1.844	1.705	1.582	1.626	1.554	1.609	1.699	8.39
84) T	Benzyl Chloride	2.373	1.648	1.772	1.688	1.799	2.058	2.054	2.190	1.948	13.35
85) T	1,3-Dichlorobe...	2.498	1.658	1.649	1.455	1.377	1.409	1.374	1.436	1.607	23.49
86) T	1,4-Dichlorobe...	2.655	1.785	1.618	1.492	1.380	1.428	1.388	1.449	1.649	25.99
87) T	sec-Butylbenzene	4.527	3.773	3.816	3.471	3.264	3.328	3.214	3.176	3.571	12.80
88) T	4-Isopropyltol...	4.426	3.674	3.683	3.416	3.254	3.356	3.253	3.366	3.554	10.98
89) T	1,2,3-Trimethy...	3.266	2.650	2.852	2.588	2.471	2.573	2.506	2.619	2.691	9.63
90) T	1,2-Dichlorobe...	2.173	1.589	1.545	1.435	1.332	1.387	1.370	1.415	1.531	17.89
91) T	d-Limonene	1.124	0.998	1.106	1.069	1.064	1.072	1.040	1.065	1.067	3.58
92) T	1,2-Dibromo-3-...	0.563	0.488	0.501	0.486	0.494	0.525	0.532	0.542	0.517	5.50
93) T	n-Undecane	1.922	1.637	1.755	1.671	1.608	1.630	1.608	1.592	1.678	6.63
94) T	1,2,4-Trichlor...	1.335	1.165	1.062	1.017	1.071	1.071	1.086	1.118	1.122	9.32
95) T	Naphthalene	4.848	3.923	3.638	3.497	3.721	3.755	3.783	3.881	11.50	11.50
96) T	n-Dodecane	1.866	1.599	1.669	1.586	1.545	1.575	1.553	1.509	1.613	6.99
97) T	Hexachlorobuta...	0.841	0.749	0.730	0.673	0.635	0.664	0.663	0.676	0.704	9.50
98) T	Cyclohexanone	1.172	0.979	1.021	0.969	0.952	0.962	0.931	0.903	0.986	8.37
99) T	tert-Butylbenzene	3.486	2.799	2.880	2.658	2.449	2.533	2.453	2.528	2.723	12.71
100) T	n-Butylbenzene	3.596	2.904	2.958	2.670	2.572	2.625	2.613	2.593	2.817	12.33

(#) = Out of Range

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\24\  
 Data File : 03241201.D  
 Acq On : 24 Mar 2012 00:34  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 24 08:11:33 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	IR Bromochloromethane (IS1)	1.000	1.000	0.0	83	-0.01
2	T Propene	2.102	1.911	9.1	69	0.00
3	T Dichlorodifluoromethane (CF	2.664	2.198	17.5	77	0.01
4	T Chloromethane	2.449	2.091	14.6	76	0.00
5	T 1,2-Dichloro-1,1,2,2-tetra	1.460	1.164	20.3	73	0.00
6	T Vinyl Chloride	2.333	1.919	17.7	75	0.00
7	T 1,3-Butadiene	1.743	1.594	8.5	78	0.00
8	T Bromomethane	1.252	1.056	15.7	77	0.00
9	T Chloroethane	1.217	1.019	16.3	78	0.00
10	T Ethanol	1.239	1.162	6.2	88	-0.07
11	T Acetonitrile	2.825	2.548	9.8	83	-0.05
12	T Acrolein	0.897	0.744	17.1	76	-0.02
13	T Acetone	1.161	0.998	14.0	82	-0.04
14	T Trichlorofluoromethane	2.277	2.018	11.4	79	0.00
15	T 2-Propanol (Isopropanol)	3.443	3.218	6.5	97	-0.05
16	T Acrylonitrile	1.916	1.760	8.1	81	-0.03
17	T 1,1-Dichloroethene	1.293	1.152	10.9	78	0.00
18	T 2-Methyl-2-Propanol (tert-B	3.812	3.168	16.9	74	-0.03
19	T Methylene Chloride	1.421	1.188	16.4	78	-0.01
20	T 3-Chloro-1-propene (Allyl C	2.255	2.246	0.4	81	-0.01
21	T Trichlorotrifluoroethane	1.171	1.027	12.3	78	0.00
22	T Carbon Disulfide	5.030	4.402	12.5	78	0.00
23	T trans-1,2-Dichloroethene	2.080	1.877	9.8	80	0.00
24	T 1,1-Dichloroethane	2.514	2.309	8.2	81	-0.01
25	T Methyl tert-Butyl Ether	4.378	4.054	7.4	81	0.00
26	T Vinyl Acetate	0.403	0.395	2.0	80	-0.03
27	T 2-Butanone (MEK)	0.986	0.934	5.3	81	-0.02
28	T cis-1,2-Dichloroethene	1.960	1.765	9.9	80	-0.01
29	T Diisopropyl Ether	1.239	1.143	7.7	80	0.00
30	T Ethyl Acetate	0.546	0.519	4.9	81	-0.02
31	T n-Hexane	2.855	2.469	13.5	78	0.00
32	T Chloroform	2.206	1.955	11.4	79	-0.03
33	S 1,2-Dichloroethane-d4 (SS1)	1.553	1.606	-3.4	87	-0.01
34	T Tetrahydrofuran (THF)	0.961	0.863	10.2	82	0.00
35	T Ethyl tert-Butyl Ether	1.821	1.683	7.6	80	0.00
36	T 1,2-Dichloroethane	1.798	1.636	9.0	80	-0.01
37	IR 1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	86	0.00
38	T 1,1,1-Trichloroethane	0.451	0.399	11.5	79	0.00

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\24\  
 Data File : 03241201.D  
 Acq On : 24 Mar 2012 00:34  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 24 08:11:33 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
39 T	Isopropyl Acetate	0.204	0.187	8.3	80	-0.02
40 T	1-Butanol	0.323	0.343	-6.2	88	-0.05
41 T	Benzene	1.258	1.072	14.8	80	-0.01
42 T	Carbon Tetrachloride	0.352	0.318	9.7	78	-0.01
43 T	Cyclohexane	0.539	0.465	13.7	80	-0.01
44 T	tert-Amyl Methyl Ether	0.939	0.851	9.4	80	0.00
45 T	1,2-Dichloropropane	0.334	0.293	12.3	80	-0.01
46 T	Bromodichloromethane	0.384	0.350	8.9	80	0.00
47 T	Trichloroethene	0.355	0.296	16.6	78	-0.01
48 T	1,4-Dioxane	0.260	0.235	9.6	82	-0.02
49 T	2,2,4-Trimethylpentane (Iso	1.557	1.379	11.4	81	0.00
50 T	Methyl Methacrylate	0.141	0.132	6.4	81	-0.02
51 T	n-Heptane	0.361	0.325	10.0	81	-0.01
52 T	cis-1,3-Dichloropropene	0.502	0.468	6.8	81	0.00
53 T	4-Methyl-2-pentanone	0.305	0.289	5.2	82	-0.02
54 T	trans-1,3-Dichloropropene	0.437	0.422	3.4	80	0.00
55 T	1,1,2-Trichloroethane	0.292	0.259	11.3	78	-0.01
56 IR	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	86	0.00
57 S	Toluene-d8 (SS2)	2.306	2.302	0.2	86	0.00
58 T	Toluene	2.793	2.399	14.1	80	0.00
59 T	2-Hexanone	1.490	1.370	8.1	83	-0.01
60 T	Dibromochloromethane	0.638	0.588	7.8	78	0.00
61 T	1,2-Dibromoethane	0.688	0.601	12.6	78	0.00
62 T	n-Butyl Acetate	1.839	1.689	8.2	82	-0.01
63 T	n-Octane	0.645	0.606	6.0	85	0.00
64 T	Tetrachloroethene	0.848	0.706	16.7	78	0.00
65 T	Chlorobenzene	1.780	1.511	15.1	79	0.00
66 T	Ethylbenzene	3.030	2.603	14.1	78	0.00
67 T	m- & p-Xylenes	2.419	2.148	11.2	81	-0.01
68 T	Bromoform	0.568	0.547	3.7	78	0.00
69 T	Styrene	1.881	1.716	8.8	82	-0.01
70 T	o-Xylene	2.550	2.246	11.9	81	-0.01
71 T	n-Nonane	1.681	1.511	10.1	81	0.00
72 T	1,1,2,2-Tetrachloroethane	1.160	1.028	11.4	79	-0.01
73 S	Bromofluorobenzene (SS3)	0.777	0.795	-2.3	88	0.00
74 T	Cumene	3.299	2.937	11.0	84	0.00
75 T	alpha-Pinene	1.611	1.440	10.6	78	0.00
76 T	n-Propylbenzene	4.039	3.517	12.9	80	0.00



Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\24\  
 Data File : 03241201.D  
 Acq On : 24 Mar 2012 00:34  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-02271204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 24 08:11:33 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
77 T	3-Ethyltoluene	3.206	2.756	14.0	79	0.00
78 T	4-Ethyltoluene	3.133	2.746	12.4	79	-0.01
79 T	1,3,5-Trimethylbenzene	2.600	2.260	13.1	80	-0.01
80 T	alpha-Methylstyrene	1.436	1.296	9.7	79	-0.01
81 T	2-Ethyltoluene	3.254	2.808	13.7	79	-0.01
82 T	1,2,4-Trimethylbenzene	2.664	2.335	12.3	80	-0.01
83 T	n-Decane	1.699	1.527	10.1	81	-0.01
84 T	Benzyl Chloride	1.948	2.052	-5.3	86	-0.02
85 T	1,3-Dichlorobenzene	1.607	1.288	19.9	79	-0.01
86 T	1,4-Dichlorobenzene	1.649	1.298	21.3	78	-0.01
87 T	sec-Butylbenzene	3.571	3.106	13.0	80	-0.01
88 T	4-Isopropyltoluene (p-Cymen)	3.554	3.122	12.2	80	-0.01
89 T	1,2,3-Trimethylbenzene	2.691	2.373	11.8	79	0.00
90 T	1,2-Dichlorobenzene	1.531	1.255	18.0	78	-0.01
91 T	d-Limonene	1.067	0.988	7.4	79	-0.01
92 T	1,2-Dibromo-3-Chloropropane	0.517	0.478	7.5	78	-0.01
93 T	n-Undecane	1.678	1.536	8.5	81	0.00
94 T	1,2,4-Trichlorobenzene	1.122	0.967	13.8	77	-0.01
95 T	Naphthalene	3.881	3.454	11.0	80	0.00
96 T	n-Dodecane	1.613	1.520	5.8	83	0.00
97 T	Hexachlorobutadiene	0.704	0.595	15.5	77	0.00
98 T	Cyclohexanone	0.986	0.960	2.6	86	-0.02
99 T	tert-Butylbenzene	2.723	2.380	12.6	81	-0.01
100 T	n-Butylbenzene	2.817	2.441	13.3	80	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

Response Factor Report MS07

Method : J:\Ms07\METHODS\X7021712B.M (RTE Integrator)  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
 1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	Avg	%RSD
1) I Bromochloromethan	3.784	3.527	3.467	3.822	3.279	2.986	2.875	3.155	2.920	3.313	10.83
2) T Dichlorodifluorom		1.470	0.956	1.190	0.904	0.778	0.569	0.897	0.827	0.949	28.81
3) T Chloromethane	2.802	2.749	2.588	3.044	2.597	2.406	2.314	2.588	2.446	2.615	8.57
4) T Vinyl Chloride	1.894	1.496	1.211	1.613	1.335	1.208	1.105	1.433	1.367	1.407	17.11
5) T Bromomethane	1.511	1.416	1.357	1.517	1.295	1.237	1.195	1.312	1.224	1.340	8.92
6) T Chloroethane		1.548	1.214	1.122	1.122	1.080	1.193	1.090	1.208	1.208	14.52
7) T Acetone	3.051	2.887	2.831	3.145	2.692	2.565	2.447	2.691	2.536	2.761	8.59
8) T Trichlorofluorome	1.591	1.440	1.366	1.547	1.322	1.229	1.214	1.385	1.295	1.376	9.52
9) T 1,1-Dichloroethen		2.134	1.795	1.866	1.536	1.426	1.404	1.494	1.411	1.633	16.43
10) T Methylene Chlorid	1.408	1.325	1.295	1.376	1.169	1.081	1.062	1.170	1.097	1.220	10.87
11) T Trichlorotrifluor	1.972	1.710	1.589	1.762	1.505	1.406	1.393	1.512	1.439	1.588	12.18
12) T trans-1,2-Dichlor	3.265	3.020	3.016	3.383	2.925	2.756	2.681	2.937	2.841	2.980	7.60
13) T 1,1-Dichloroethan	5.168	4.740	4.762	5.045	4.335	4.137	4.166	4.690	4.471	4.613	7.90
14) T Methyl tert-Butyl	1.885	1.626	1.630	1.776	1.524	1.434	1.388	1.575	1.482	1.591	10.09
15) T cis-1,2-Dichloroe		3.041	3.257	2.586	2.390	2.280	2.524	2.374	2.374	2.636	14.03
16) T Chloroform	2.000	2.002	2.032	2.018	1.989	1.973	1.960	1.938	1.932	1.983	1.75
17) S 1,2-Dichloroethan	2.582	2.460	2.375	2.604	2.222	2.090	1.988	2.222	2.094	2.293	9.73
18) T 1,2-Dichloroethan	2.587	2.431	2.340	2.573	2.191	2.059	2.007	2.278	2.150	2.291	9.18
19) T 1,1,1-Trichloroet		7.536	7.534	6.055	5.665	5.665	5.395	5.903	5.468	6.222	14.88
20) T Benzene	2.280	1.842	1.769	1.950	1.676	1.579	1.549	1.771	1.692	1.790	12.41
21) T Carbon Tetrachlor											
22) I 1,4-Difluorobenze	0.417	0.397	0.373	0.409	0.354	0.331	0.319	0.358	0.336	0.366	9.68
23) T 1,2-Dichloropropa	0.491	0.464	0.436	0.489	0.420	0.398	0.394	0.447	0.416	0.439	8.26
24) T Bromodichlorometh	0.418	0.379	0.349	0.390	0.324	0.305	0.287	0.326	0.307	0.343	12.96
25) T Trichloroethene	0.324	0.290	0.270	0.278	0.242	0.231	0.227	0.258	0.243	0.263	11.99
26) T 1,4-Dioxane	0.567	0.523	0.508	0.540	0.499	0.484	0.488	0.559	0.522	0.521	5.67
27) T cis-1,3-Dichlorop	0.457	0.434	0.431	0.444	0.419	0.409	0.417	0.505	0.482	0.444	7.22
28) T trans-1,3-Dichlor											

(#) = Out of Range ### Number of calibration levels exceeded format ###  
 X7021712B.M Wed Mar 28 08:50:39 2012

Response Factor Report MS07

Method : J:\MS07\METHODS\X7021712B.M (RTE Integrator)  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
 1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	AVG	%RSD
29) T 1,1,2-Trichloroet	0.315	0.296	0.286	0.305	0.263	0.248	0.242	0.274	0.257	0.276	9.32
30) S Toluene-d8 (SS2)	1.053	1.071	1.054	1.047	1.051	1.052	1.071	1.060	1.049	1.056	0.84
31) T Toluene	1.782	1.541	1.420	1.485	1.269	1.198	1.177	1.307	1.186	1.374	14.76
32) T 1,2-Dibromoethane	0.379	0.362	0.345	0.362	0.323	0.307	0.304	0.351	0.331	0.340	7.65
33) T Tetrachloroethene	0.423	0.394	0.371	0.400	0.341	0.320	0.311	0.347	0.326	0.359	11.03
-----ISTD-----											
34) I Chlorobenzene-d5											
35) T Chlorobenzene	4.581	4.172	4.017	4.308	3.688	3.469	3.311	3.705	3.345	3.844	11.68
36) T Ethylbenzene	7.604	7.082	6.707	7.235	6.297	6.070	5.915	6.682	5.878	6.608	9.30
37) T m,p-Xylene	5.861	5.347	5.132	5.536	4.930	4.777	4.719	5.268	4.558	5.125	8.24
38) T o-Xylene	6.505	5.826	5.526	5.889	5.188	4.995	4.950	5.594	4.932	5.490	9.67
39) T 1,1,2,2-Tetrachlo	2.655	2.674	2.710	2.646	2.437	2.371	2.298	2.761	2.498	2.561	6.40
40) S Bromofluorobenzene	1.712	1.701	1.695	1.697	1.683	1.680	1.670	1.642	1.593	1.675	2.20
41) T 1,3-Dichlorobenze	3.304	3.261	3.222	3.204	2.835	2.719	2.608	3.138	2.846	3.015	8.71
42) T 1,4-Dichlorobenze	3.470	3.303	3.208	3.208	2.803	2.671	2.579	3.174	2.891	3.034	10.13
43) T 1,2-Dichlorobenze	3.271	3.123	3.112	3.026	2.662	2.536	2.418	3.053	2.772	2.886	10.31
44) T 1,2,4-Trichlorobe	2.184	2.118	2.169	1.753	1.641	1.608	1.575	2.139	1.978	1.907	13.66
45) T Naphthalene	6.857	6.355	6.704	5.143	5.107	5.172	5.297	8.126	7.328	6.232	17.82
46) T Hexachlorobutadie	1.486	1.423	1.403	1.217	1.055	1.013	0.975	1.227	1.145	1.216	15.40

(#) = Out of Range ### Number of calibration levels exceeded format ###  
 X7021712B.M Wed Mar 28 08:50:39 2012

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_03\28\  
 Data File : 03281203.D  
 Acq On : 28 Mar 2012 7:50 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00069  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Mar 28 08:50:09 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane (IS1)	1.000	1.000	0.0	93	0.00
2 T	Dichlorodifluoromethane (CF	3.313	2.572	22.4	73	0.00
3 T	Chloromethane	0.949	0.733	22.8	76	0.00
4 T	Vinyl Chloride	2.615	2.066	21.0	74	0.00
5 T	Bromomethane	1.407	1.205	14.4	84	0.00
6 T	Chloroethane	1.340	1.060	20.9	77	0.00
7 T	Acetone	1.208	1.079	10.7	83	0.00
8 T	Trichlorofluoromethane	2.761	2.164	21.6	75	0.00
9 T	1,1-Dichloroethene	1.376	1.146	16.7	81	0.00
10 T	Methylene Chloride	1.633	1.283	21.4	78	0.00
11 T	Trichlorotrifluoroethane	1.220	1.053	13.7	84	0.00
12 T	trans-1,2-Dichloroethene	1.588	1.270	20.0	79	0.00
13 T	1,1-Dichloroethane	2.980	2.388	19.9	76	0.00
14 T	Methyl tert-Butyl Ether	4.613	3.616	21.6	78	0.00
15 T	cis-1,2-Dichloroethene	1.591	1.323	16.8	81	0.00
16 T	Chloroform	2.636	2.156	18.2	78	0.00
17 S	1,2-Dichloroethane-d4 (SS1)	1.983	1.932	2.6	91	0.00
18 T	1,2-Dichloroethane	2.293	1.727	24.7	73	0.00
19 T	1,1,1-Trichloroethane	2.291	1.845	19.5	79	0.00
20 T	Benzene	6.222	5.286	15.0	82	0.00
21 T	Carbon Tetrachloride	1.790	1.409	21.3	79	0.00
22 I	1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	102	0.00
23 T	1,2-Dichloropropane	0.366	0.270	26.2	78	0.00
24 T	Bromodichloromethane	0.439	0.322	26.7	78	0.00
25 T	Trichloroethene	0.343	0.258	24.8	82	0.00
26 T	1,4-Dioxane	0.263	0.197	25.1	83	0.00
27 T	cis-1,3-Dichloropropene	0.521	0.390	25.1	80	0.00
28 T	trans-1,3-Dichloropropene	0.444	0.337	24.1	82	0.00
29 T	1,1,2-Trichloroethane	0.276	0.208	24.6	81	0.00
30 S	Toluene-d8 (SS2)	1.056	1.185	-12.2	115	0.00
31 T	Toluene	1.374	1.045	23.9	84	0.00
32 T	1,2-Dibromoethane	0.340	0.264	22.4	84	0.00
33 T	Tetrachloroethene	0.359	0.286	20.3	86	0.00
34 I	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	98	0.00
35 T	Chlorobenzene	3.844	3.214	16.4	86	0.00
36 T	Ethylbenzene	6.608	5.251	20.5	82	0.00
37 T	m,p-Xylene	5.125	4.240	17.3	85	0.00

*Handwritten signature/initials: KR 3/28/12*

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_03\28\  
 Data File : 03281203.D  
 Acq On : 28 Mar 2012 7:50 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00069  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Mar 28 08:50:09 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
38 T	o-Xylene	5.490	4.614	16.0	87	0.00
39 T	1,1,2,2-Tetrachloroethane	2.561	2.098	18.1	85	0.00
40 S	Bromofluorobenzene (SS3)	1.675	1.968	-17.5	115	0.00
41 T	1,3-Dichlorobenzene	3.015	2.714	10.0	94	0.00
42 T	1,4-Dichlorobenzene	3.034	2.668	12.1	94	0.00
43 T	1,2-Dichlorobenzene	2.886	2.597	10.0	96	0.00
44 T	1,2,4-Trichlorobenzene	1.907	1.795	5.9	108	0.00
45 T	Naphthalene	6.232	6.258	-0.4	120	0.00
46 T	Hexachlorobutadiene	1.216	1.030	15.3	96	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*KR 3/28/12*

## LABORATORY REPORT

April 5, 2012

Christopher Cote  
Montana DEQ  
1100 N. Last Chance Gulch  
Helena, MT 59601

### **RE: Background Indoor Air Study**

Dear Christopher:

Enclosed are the results of the samples submitted to our laboratory on March 22, 2012. For your reference, these analyses have been assigned our service request number P1201121.

All analyses were performed according to our laboratory's NELAP and DoD-ELAP-approved quality assurance program. The test results meet requirements of the current NELAP and DoD-ELAP standards, where applicable, and except as noted in the laboratory case narrative provided. For a specific list of NELAP and DoD-ELAP-accredited analytes, refer to the certifications section at [www.caslab.com](http://www.caslab.com). Results are intended to be considered in their entirety and apply only to the samples analyzed and reported herein.

Columbia Analytical Services, Inc. is certified by the California Department of Health Services, NELAP Laboratory Certificate No. 02115CA; Arizona Department of Health Services, Certificate No. AZ0694; Florida Department of Health, NELAP Certification E871020; New Jersey Department of Environmental Protection, NELAP Laboratory Certification ID #CA009; New York State Department of Health, NELAP NY Lab ID No: 11221; Oregon Environmental Laboratory Accreditation Program, NELAP ID: CA20007; The American Industrial Hygiene Association, Laboratory #101661; United States Department of Defense Environmental Laboratory Accreditation Program (DoD-ELAP), Certificate No. L11-203; Pennsylvania Registration No. 68-03307; TX Commission of Environmental Quality, NELAP ID T104704413-11-2; Minnesota Department of Health, NELAP Certificate No. 362188; Washington State Department of Ecology, ELAP Lab ID: C946, State of Utah Department of Health, NELAP Certificate No. CA015272011-1; Los Angeles Department of Building and Safety, Approval No: TA00001. Each of the certifications listed above have an explicit Scope of Accreditation that applies to specific matrices/methods/analytes; therefore, please contact me for information corresponding to a particular certification.

If you have any questions, please call me at (805) 526-7161.

Respectfully submitted,

**Columbia Analytical Services, Inc.**

Kate Aguilera  
Project Manager

Client: Montana DEQ  
Project: Background Indoor Air Study

CAS Project No: P1201121

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## CASE NARRATIVE

The samples were received intact under chain of custody on March 22, 2012 and were stored in accordance with the analytical method requirements. Please refer to the sample acceptance check form for additional information. The results reported herein are applicable only to the condition of the samples at the time of sample receipt.

### Air-Phase Petroleum Hydrocarbons (APH) Analysis

The samples were analyzed for total aliphatic and aromatic gasoline range hydrocarbons by gas chromatography/mass spectrometry according to the Method for the Determination of Air-Phase Petroleum Hydrocarbons (APH), Massachusetts Department of Environmental Protection, Revision 1, December, 2009.

### Volatile Organic Compound Analysis

The samples were also analyzed in Scan and SIM mode for selected volatile organic compounds in accordance with EPA Method TO-15 from the Compendium of Methods for the Determination of Toxic Organic Compounds in Ambient Air, Second Edition (EPA/625/R-96/010b), January, 1999. The analytical system was comprised of a gas chromatograph / mass spectrometer (GC/MS) interfaced to a whole-air preconcentrator.

The reporting limit is elevated sample “28-1F-031912”. The chromatogram indicated the presence of non-target background components. The sample was diluted in order to prevent damage to the instrument and to achieve optimal resolution of the target analytes.

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*The results of analyses are given in the attached laboratory report. All results are intended to be considered in their entirety, and Columbia Analytical Services, Inc. (CAS) is not responsible for utilization of less than the complete report.*

*Use of Columbia Analytical Services, Inc. (CAS) Name. Client shall not use CAS's name or trademark in any marketing or reporting materials, press releases or in any other manner (“Materials”) whatsoever and shall not attribute to CAS any test result, tolerance or specification derived from CAS's data (“Attribution”) without CAS's prior written consent, which may be withheld by CAS for any reason in its sole discretion. To request CAS's consent, Client shall provide copies of the proposed Materials or Attribution and describe in writing Client's proposed use of such Materials or Attribution. If CAS has not provided written approval of the Materials or Attribution within ten (10) days of receipt from Client, Client's request to use CAS's name or trademark in any Materials or Attribution shall be deemed denied. CAS may, in its discretion, reasonably charge Client for its time in reviewing Materials or Attribution requests. Client acknowledges and agrees that the unauthorized use of CAS's name or trademark may cause CAS to incur irreparable harm for which the recovery of money damages will be inadequate. Accordingly, Client acknowledges and agrees that a violation shall justify preliminary injunctive relief. For questions contact the laboratory.*

DETAIL SUMMARY REPORT

Client: Montana DEQ  
 Project ID: Background Indoor Air Study

Service Request: P1201121

Date Received: 3/22/2012  
 Time Received: 10:00

Client Sample ID	Lab Code	Matrix	Date Collected	Time Collected	Container ID	Pi1 (psig)	Pf1 (psig)	MA APH 1.0 - MA VOC PH Can	TO-15 - VOC Cans	TO-15 - VOC SIM
19-1F-031412	P1201121-001	Air	3/15/2012	18:15	AC01817	-3.81	3.51	X	X	X
53-1F-031412	P1201121-002	Air	3/15/2012	18:15	AC01478	-1.95	3.66	X	X	X
21-1F-031512	P1201121-003	Air	3/16/2012	12:45	AC01566	-4.65	3.73	X	X	X
23-1F-031512	P1201121-004	Air	3/16/2012	17:51	AC00988	-4.67	3.62	X	X	X
17-1F-031712	P1201121-005	Air	3/16/2012	15:00	AC01790	-2.83	3.61	X	X	X
20-1F-031712	P1201121-006	Air	3/18/2012	11:56	AC01305	-1.78	3.86	X	X	X
28-1F-031912	P1201121-007	Air	3/20/2012	15:05	AC00721	-4.99	3.65	X	X	X
22-1F-031912	P1201121-008	Air	3/20/2012	12:10	AC01171	-2.96	3.64	X	X	X
26-1F-031912	P1201121-009	Air	3/20/2012	09:50	AC01783	-3.72	3.54	X	X	X



**Air - Chain of Custody Record & Analytical Service Request**

**Requested Turnaround Time in Business Days (Surcharges) please circle**  
 1 Day (100%) 2 Day (75%) 3 Day (50%) 4 Day (35%) 5 Day (25%) 10 Day-Standard

CAS Project No. **P1201121**

Company Name & Address (Reporting Information)  
 MT Dept Env. Quality  
 PO Box 200901  
 1100 N. East Chace Sault  
 Helena MT 59620-0901

Project Manager: **Chris Cote**

Phone: 406-841-5078 Fax: 406-841-5050

Email Address for Result Reporting: **cote2@mt.gov; Martidbj@cdm.com**

Project Name: **Background Indoor Air Study**

Project Number: \_\_\_\_\_

P.O. # / Billing Information: **Reference to Background Indoor Air Study**

Sampler (Print & Sign): **Laura Alvey + Alan Schiff**

Client Sample ID	Laboratory ID Number	Date Collected	Time Collected	Canister ID (Bar code # - AC, SC, etc.)	Flow Controller ID (Bar code # - FC #)	Canister Start Pressure "Hg	Canister End Pressure "Hg/psig	Sample Volume	CAS Contact:	Analysis Method	Comments
19-1F-031412	①-375	3/15/12	18:15	FLA01817	FLA00327	-28.7	-7.0		X	TO-15	Analyze
53-1F-031412	②-154	3/15/12	18:15	FLA01478	FLA00412	-28.3	-4.0		X	TO-15 SIM	Project-specific list of compounds to project-specific reagents
21-1F-031512	③-458	3/16/12	12:45	FLA01566	FLA00463	-28.0	-10.0				
23-1F-031512	④-460	3/16/12	17:51	FLA00988	FLA00467	-27.6	-8.4				
17-1F-031512	⑤-272	3/16/12	15:00	FLA01790	FLA00069	-28.0	-4.5				
20-1F-031712	⑥-169	3/18/12	11:56	FLA01305	FLA00159	-28.5	-3.3				
28-1F-031912	⑦-990	3/20/12	15:05	FLA00721	FLA00478	-27.5	-9.0				
22-1F-031912	⑧-195	3/20/12	12:10	FLA01171	FLA00411	-27.5	-5.0				
26-1F-031912	⑨-370	3/20/12	9:50	FLA01783	FLA00401	-29.0	-7.5				
	-250										
	1867										

**Report Tier Levels - please select**

Tier I - Results (Default if not specified) \_\_\_\_\_

Tier II (Results + QC Summaries) \_\_\_\_\_

Tier III (Results + QC & Calibration Summaries)  X

Tier IV (Data Validation Package) 10% Surcharge \_\_\_\_\_

EDD required (Yes/No)  Yes  No

Type: **CVS Excel format**

Relinquished by: (Signature) \_\_\_\_\_ Date: 3/21/12 Time: 9:00 am

Received by: (Signature) \_\_\_\_\_ Date: 3/21/12 Time: 9:00 am

Cooler / Blank Temperature \_\_\_\_\_ °C

**Sample Acceptance Check Form**

Client: Montana DEQ Work order: P1201121

Project: Background Indoor Air Study

Sample(s) received on: 3/22/12 Date opened: 3/22/12 by: MZAMORA

**Note:** This form is used for all samples received by CAS. The use of this form for custody seals is strictly meant to indicate presence/absence and not as an indication of compliance or nonconformity. Thermal preservation and pH will only be evaluated either at the request of the client and/or as required by the method/SOP.

- |  | <u>Yes</u>                          | <u>No</u>                           | <u>N/A</u>                          |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| 1 Were <b>sample containers</b> properly marked with client sample ID?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 2 Container(s) <b>supplied by CAS</b> ?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 3 Did <b>sample containers</b> arrive in good condition?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 4 Were <b>chain-of-custody</b> papers used and filled out?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 5 Did <b>sample container labels</b> and/or tags agree with custody papers?                                      | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 6 Was <b>sample volume</b> received adequate for analysis?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 7 Are samples within specified holding times?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 8 Was proper <b>temperature</b> (thermal preservation) of cooler at receipt adhered to?                          | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 9 Was a <b>trip blank</b> received?  | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| 10 Were <b>custody seals</b> on outside of cooler/Box?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| Location of seal(s)? _____ Sealing Lid?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were signature and date included?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were seals intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were custody seals on outside of sample container?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| Location of seal(s)? _____ Sealing Lid?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were signature and date included?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were seals intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 11 Do containers have appropriate <b>preservation</b> , according to method/SOP or Client specified information? | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Is there a client indication that the submitted samples are <b>pH</b> preserved?                                 | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were <b>VOA vials</b> checked for presence/absence of air bubbles?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Does the client/method/SOP require that the analyst check the sample pH and <u>if necessary</u> alter it?        | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 12 <b>Tubes:</b> Are the tubes capped and intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Do they contain moisture?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 13 <b>Badges:</b> Are the badges properly capped and intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Are dual bed badges separated and individually capped and intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |

Lab Sample ID	Container Description	Required pH *	Received pH	Adjusted pH	VOA Headspace (Presence/Absence)	Receipt / Preservation Comments
P1201121-001.01	6.0 L Ambient Can					
P1201121-002.01	6.0 L Ambient Can					
P1201121-003.01	6.0 L Ambient Can					
P1201121-004.01	6.0 L Ambient Can					
P1201121-005.01	6.0 L Ambient Can					
P1201121-006.01	6.0 L Ambient Can					
P1201121-007.01	6.0 L Ambient Can					
P1201121-008.01	6.0 L Ambient Can					

Explain any discrepancies: (include lab sample ID numbers): \_\_\_\_\_



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 19-1F-031412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-001

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01817

Date Collected: 3/15/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.81      Final Pressure (psig): 3.51

Canister Dilution Factor: 1.67

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	41	33	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	17	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	8.4	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 53-1F-031412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-002

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01478

Date Collected: 3/15/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.95      Final Pressure (psig): 3.66

Canister Dilution Factor: 1.44

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	51	29	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	20	14	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	7.2	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 21-1F-031512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-003

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01566

Date Collected: 3/16/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.65      Final Pressure (psig): 3.73

Canister Dilution Factor: 1.83

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	84	37	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	67	18	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	15	9.2	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 23-1F-031512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-004

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00988

Date Collected: 3/16/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.67      Final Pressure (psig): 3.62

Canister Dilution Factor: 1.83

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	73	37	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	21	18	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	9.2	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 17-1F-031712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-005

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01790

Date Collected: 3/16/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.83      Final Pressure (psig): 3.61

Canister Dilution Factor: 1.54

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	190	31	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	27	15	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	10	7.7	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 20-1F-031712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-006

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01305

Date Collected: 3/18/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.78      Final Pressure (psig): 3.86

Canister Dilution Factor: 1.44

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	81	29	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	74	14	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	13	7.2	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 28-1F-031912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-007

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00721

Date Collected: 3/20/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 0.30 Liter(s)

Initial Pressure (psig): -4.99      Final Pressure (psig): 3.65

Canister Dilution Factor: 1.89

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	130	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	<b>190</b>	63	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	32	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 22-1F-031912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-008

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01171

Date Collected: 3/20/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/28/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.96      Final Pressure (psig): 3.64

Canister Dilution Factor: 1.56

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	41	31	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	39	16	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	7.8	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 26-1F-031912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-009

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01783

Date Collected: 3/20/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/28/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.72      Final Pressure (psig): 3.54

Canister Dilution Factor: 1.66

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	33	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	17	17	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	8.3	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P120327-MB

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	20	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	10	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	5.0	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

LABORATORY CONTROL SAMPLE SUMMARY

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**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P120327-LCS

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: NA Liter(s)

Compound	Spike Amount ng	Result ng	% Recovery	CAS	Data Qualifier
				Acceptance Limits	
C5 - C8 Aliphatic Hydrocarbons	24.5	22.8	93	70-130	
C9 - C12 Aliphatic Hydrocarbons	25.0	23.3	93	70-130	
C9 - C10 Aromatic Hydrocarbons	48.5	41.5	86	70-130	

LABORATORY DUPLICATE SUMMARY RESULTS

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**Client:** Montana DEQ

**Client Sample ID:** 53-1F-031412

**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121

CAS Sample ID: P1201121-002DUP

Test Code: Massachusetts APH, Revision 1, December 2009

Date Collected: 3/15/12

Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8

Date Received: 3/22/12

Analyst: Elsa Moctezuma

Date Analyzed: 3/27/12

Sampling Media: 6.0 L Summa Canister

Volume(s) Analyzed: 1.00 Liter(s)

Test Notes:

Container ID: AC01478

Initial Pressure (psig): -1.95      Final Pressure (psig): 3.66

Canister Dilution Factor: 1.44

Compound	Sample Result µg/m <sup>3</sup>	Duplicate Sample Result µg/m <sup>3</sup>	Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	50.9	50.2	50.55	1	30	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	19.6	19.8	19.7	1	30	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	ND	-	-	30	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 19-1F-031412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-001

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01817

Date Collected: 3/15/12  
Date Received: 3/22/12  
Date Analyzed: 3/27/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.81 Final Pressure (psig): 3.51

Canister Dilution Factor: 1.67

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	2.4	0.84	1.4	0.49	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.0	0.84	0.40	0.17	
74-87-3	Chloromethane	ND	0.84	ND	0.40	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.84	ND	0.12	
106-99-0	1,3-Butadiene	ND	0.84	ND	0.38	
74-83-9	Bromomethane	ND	0.84	ND	0.22	
75-00-3	Chloroethane	ND	0.84	ND	0.32	
64-17-5	Ethanol	180	8.4	97	4.4	
75-05-8	Acetonitrile	ND	0.84	ND	0.50	
107-02-8	Acrolein	ND	3.3	ND	1.5	
67-64-1	Acetone	23	8.4	9.5	3.5	
75-69-4	Trichlorofluoromethane	1.3	0.84	0.23	0.15	
67-63-0	2-Propanol (Isopropyl Alcohol)	12	1.7	4.7	0.68	
107-13-1	Acrylonitrile	ND	0.84	ND	0.38	
75-35-4	1,1-Dichloroethene	ND	0.84	ND	0.21	
75-09-2	Methylene Chloride	17	0.84	5.0	0.24	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.84	ND	0.27	
76-13-1	Trichlorotrifluoroethane	ND	0.84	ND	0.11	
75-15-0	Carbon Disulfide	ND	8.4	ND	2.7	
156-60-5	trans-1,2-Dichloroethene	ND	0.84	ND	0.21	
75-34-3	1,1-Dichloroethane	ND	0.84	ND	0.21	
1634-04-4	Methyl tert-Butyl Ether	ND	0.84	ND	0.23	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 19-1F-031412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-001

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01817

**Date Collected:** 3/15/12  
**Date Received:** 3/22/12  
**Date Analyzed:** 3/27/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -3.81      Final Pressure (psig): 3.51

Canister Dilution Factor: 1.67

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	8.4	ND	2.4	
78-93-3	2-Butanone (MEK)	ND	8.4	ND	2.8	
156-59-2	cis-1,2-Dichloroethene	ND	0.84	ND	0.21	
141-78-6	Ethyl Acetate	<b>16</b>	1.7	<b>4.4</b>	0.46	
110-54-3	n-Hexane	ND	0.84	ND	0.24	
67-66-3	Chloroform	ND	0.84	ND	0.17	
109-99-9	Tetrahydrofuran (THF)	ND	0.84	ND	0.28	
71-55-6	1,1,1-Trichloroethane	ND	0.84	ND	0.15	
56-23-5	Carbon Tetrachloride	ND	0.84	ND	0.13	
110-82-7	Cyclohexane	ND	1.7	ND	0.49	
78-87-5	1,2-Dichloropropane	ND	0.84	ND	0.18	
75-27-4	Bromodichloromethane	ND	0.84	ND	0.12	
123-91-1	1,4-Dioxane	ND	0.84	ND	0.23	
80-62-6	Methyl Methacrylate	ND	1.7	ND	0.41	
142-82-5	n-Heptane	ND	0.84	ND	0.20	
10061-01-5	cis-1,3-Dichloropropene	ND	0.84	ND	0.18	
108-10-1	4-Methyl-2-pentanone	ND	0.84	ND	0.20	
10061-02-6	trans-1,3-Dichloropropene	ND	0.84	ND	0.18	
108-88-3	Toluene	<b>5.7</b>	0.84	<b>1.5</b>	0.22	
591-78-6	2-Hexanone	ND	0.84	ND	0.20	
124-48-1	Dibromochloromethane	ND	0.84	ND	0.098	
106-93-4	1,2-Dibromoethane	ND	0.84	ND	0.11	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 19-1F-031412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-001

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01817

Date Collected: 3/15/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.81 Final Pressure (psig): 3.51

Canister Dilution Factor: 1.67

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	3.3	0.84	0.70	0.18	
111-65-9	n-Octane	ND	0.84	ND	0.18	
108-90-7	Chlorobenzene	ND	0.84	ND	0.18	
179601-23-1	m,p-Xylenes	1.9	1.7	0.45	0.38	
75-25-2	Bromoform	ND	0.84	ND	0.081	
100-42-5	Styrene	ND	0.84	ND	0.20	
95-47-6	o-Xylene	ND	0.84	ND	0.19	
111-84-2	n-Nonane	1.4	0.84	0.27	0.16	
98-82-8	Cumene	ND	0.84	ND	0.17	
80-56-8	alpha-Pinene	5.1	0.84	0.92	0.15	
103-65-1	n-Propylbenzene	ND	0.84	ND	0.17	
622-96-8	4-Ethyltoluene	ND	0.84	ND	0.17	
108-67-8	1,3,5-Trimethylbenzene	ND	0.84	ND	0.17	
95-63-6	1,2,4-Trimethylbenzene	ND	0.84	ND	0.17	
100-44-7	Benzyl Chloride	ND	0.84	ND	0.16	
541-73-1	1,3-Dichlorobenzene	ND	0.84	ND	0.14	
106-46-7	1,4-Dichlorobenzene	ND	0.84	ND	0.14	
95-50-1	1,2-Dichlorobenzene	ND	0.84	ND	0.14	
5989-27-5	d-Limonene	39	0.84	7.0	0.15	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.84	ND	0.086	
120-82-1	1,2,4-Trichlorobenzene	ND	0.84	ND	0.11	
87-68-3	Hexachlorobutadiene	ND	0.84	ND	0.078	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 53-1F-031412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-002

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01478

Date Collected: 3/15/12  
Date Received: 3/22/12  
Date Analyzed: 3/27/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.95 Final Pressure (psig): 3.66

Canister Dilution Factor: 1.44

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	ND	0.72	ND	0.42	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.0	0.72	0.40	0.15	
74-87-3	Chloromethane	ND	0.72	ND	0.35	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.72	ND	0.10	
106-99-0	1,3-Butadiene	ND	0.72	ND	0.33	
74-83-9	Bromomethane	ND	0.72	ND	0.19	
75-00-3	Chloroethane	ND	0.72	ND	0.27	
64-17-5	Ethanol	200	7.2	110	3.8	
75-05-8	Acetonitrile	ND	0.72	ND	0.43	
107-02-8	Acrolein	ND	2.9	ND	1.3	
67-64-1	Acetone	24	7.2	10	3.0	
75-69-4	Trichlorofluoromethane	1.3	0.72	0.24	0.13	
67-63-0	2-Propanol (Isopropyl Alcohol)	ND	1.4	ND	0.59	
107-13-1	Acrylonitrile	ND	0.72	ND	0.33	
75-35-4	1,1-Dichloroethene	ND	0.72	ND	0.18	
75-09-2	Methylene Chloride	19	0.72	5.4	0.21	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.72	ND	0.23	
76-13-1	Trichlorotrifluoroethane	ND	0.72	ND	0.094	
75-15-0	Carbon Disulfide	ND	7.2	ND	2.3	
156-60-5	trans-1,2-Dichloroethene	ND	0.72	ND	0.18	
75-34-3	1,1-Dichloroethane	ND	0.72	ND	0.18	
1634-04-4	Methyl tert-Butyl Ether	ND	0.72	ND	0.20	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 53-1F-031412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-002

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01478

**Date Collected:** 3/15/12  
**Date Received:** 3/22/12  
**Date Analyzed:** 3/27/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -1.95      Final Pressure (psig): 3.66

Canister Dilution Factor: 1.44

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	7.2	ND	2.0	
78-93-3	2-Butanone (MEK)	ND	7.2	ND	2.4	
156-59-2	cis-1,2-Dichloroethene	ND	0.72	ND	0.18	
141-78-6	Ethyl Acetate	16	1.4	4.5	0.40	
110-54-3	n-Hexane	ND	0.72	ND	0.20	
67-66-3	Chloroform	ND	0.72	ND	0.15	
109-99-9	Tetrahydrofuran (THF)	ND	0.72	ND	0.24	
71-55-6	1,1,1-Trichloroethane	ND	0.72	ND	0.13	
56-23-5	Carbon Tetrachloride	ND	0.72	ND	0.11	
110-82-7	Cyclohexane	ND	1.4	ND	0.42	
78-87-5	1,2-Dichloropropane	ND	0.72	ND	0.16	
75-27-4	Bromodichloromethane	ND	0.72	ND	0.11	
123-91-1	1,4-Dioxane	ND	0.72	ND	0.20	
80-62-6	Methyl Methacrylate	ND	1.4	ND	0.35	
142-82-5	n-Heptane	ND	0.72	ND	0.18	
10061-01-5	cis-1,3-Dichloropropene	ND	0.72	ND	0.16	
108-10-1	4-Methyl-2-pentanone	ND	0.72	ND	0.18	
10061-02-6	trans-1,3-Dichloropropene	ND	0.72	ND	0.16	
108-88-3	Toluene	5.5	0.72	1.5	0.19	
591-78-6	2-Hexanone	ND	0.72	ND	0.18	
124-48-1	Dibromochloromethane	ND	0.72	ND	0.085	
106-93-4	1,2-Dibromoethane	ND	0.72	ND	0.094	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 53-1F-031412  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201121  
**CAS Sample ID:** P1201121-002

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01478

**Date Collected:** 3/15/12  
**Date Received:** 3/22/12  
**Date Analyzed:** 3/27/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -1.95      Final Pressure (psig): 3.66

Canister Dilution Factor: 1.44

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	2.6	0.72	0.56	0.15	
111-65-9	n-Octane	ND	0.72	ND	0.15	
108-90-7	Chlorobenzene	ND	0.72	ND	0.16	
179601-23-1	m,p-Xylenes	1.8	1.4	0.42	0.33	
75-25-2	Bromoform	ND	0.72	ND	0.070	
100-42-5	Styrene	ND	0.72	ND	0.17	
95-47-6	o-Xylene	ND	0.72	ND	0.17	
111-84-2	n-Nonane	1.4	0.72	0.27	0.14	
98-82-8	Cumene	ND	0.72	ND	0.15	
80-56-8	alpha-Pinene	5.2	0.72	0.93	0.13	
103-65-1	n-Propylbenzene	ND	0.72	ND	0.15	
622-96-8	4-Ethyltoluene	ND	0.72	ND	0.15	
108-67-8	1,3,5-Trimethylbenzene	ND	0.72	ND	0.15	
95-63-6	1,2,4-Trimethylbenzene	ND	0.72	ND	0.15	
100-44-7	Benzyl Chloride	ND	0.72	ND	0.14	
541-73-1	1,3-Dichlorobenzene	ND	0.72	ND	0.12	
106-46-7	1,4-Dichlorobenzene	ND	0.72	ND	0.12	
95-50-1	1,2-Dichlorobenzene	ND	0.72	ND	0.12	
5989-27-5	d-Limonene	47	0.72	8.5	0.13	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.72	ND	0.075	
120-82-1	1,2,4-Trichlorobenzene	ND	0.72	ND	0.097	
87-68-3	Hexachlorobutadiene	ND	0.72	ND	0.068	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 21-1F-031512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-003

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01566

Date Collected: 3/16/12  
Date Received: 3/22/12  
Date Analyzed: 3/27/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.65      Final Pressure (psig): 3.73

Canister Dilution Factor: 1.83

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	10	0.92	6.1	0.53	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.2	0.92	0.44	0.19	
74-87-3	Chloromethane	ND	0.92	ND	0.44	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.92	ND	0.13	
106-99-0	1,3-Butadiene	ND	0.92	ND	0.41	
74-83-9	Bromomethane	ND	0.92	ND	0.24	
75-00-3	Chloroethane	ND	0.92	ND	0.35	
64-17-5	Ethanol	750	9.2	400	4.9	
75-05-8	Acetonitrile	ND	0.92	ND	0.55	
107-02-8	Acrolein	3.9	3.7	1.7	1.6	
67-64-1	Acetone	59	9.2	25	3.9	
75-69-4	Trichlorofluoromethane	2.3	0.92	0.41	0.16	
67-63-0	2-Propanol (Isopropyl Alcohol)	53	1.8	22	0.74	
107-13-1	Acrylonitrile	ND	0.92	ND	0.42	
75-35-4	1,1-Dichloroethene	ND	0.92	ND	0.23	
75-09-2	Methylene Chloride	2.0	0.92	0.59	0.26	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.92	ND	0.29	
76-13-1	Trichlorotrifluoroethane	ND	0.92	ND	0.12	
75-15-0	Carbon Disulfide	ND	9.2	ND	2.9	
156-60-5	trans-1,2-Dichloroethene	ND	0.92	ND	0.23	
75-34-3	1,1-Dichloroethane	ND	0.92	ND	0.23	
1634-04-4	Methyl tert-Butyl Ether	ND	0.92	ND	0.25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 21-1F-031512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-003

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01566

Date Collected: 3/16/12  
Date Received: 3/22/12  
Date Analyzed: 3/27/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.65 Final Pressure (psig): 3.73

Canister Dilution Factor: 1.83

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	9.2	ND	2.6	
78-93-3	2-Butanone (MEK)	ND	9.2	ND	3.1	
156-59-2	cis-1,2-Dichloroethene	ND	0.92	ND	0.23	
141-78-6	Ethyl Acetate	<b>13</b>	1.8	<b>3.5</b>	0.51	
110-54-3	n-Hexane	<b>1.3</b>	0.92	<b>0.37</b>	0.26	
67-66-3	Chloroform	ND	0.92	ND	0.19	
109-99-9	Tetrahydrofuran (THF)	<b>1.4</b>	0.92	<b>0.47</b>	0.31	
71-55-6	1,1,1-Trichloroethane	ND	0.92	ND	0.17	
56-23-5	Carbon Tetrachloride	ND	0.92	ND	0.15	
110-82-7	Cyclohexane	ND	1.8	ND	0.53	
78-87-5	1,2-Dichloropropane	ND	0.92	ND	0.20	
75-27-4	Bromodichloromethane	ND	0.92	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.92	ND	0.25	
80-62-6	Methyl Methacrylate	ND	1.8	ND	0.45	
142-82-5	n-Heptane	<b>4.1</b>	0.92	<b>1.0</b>	0.22	
10061-01-5	cis-1,3-Dichloropropene	ND	0.92	ND	0.20	
108-10-1	4-Methyl-2-pentanone	<b>2.1</b>	0.92	<b>0.52</b>	0.22	
10061-02-6	trans-1,3-Dichloropropene	ND	0.92	ND	0.20	
108-88-3	Toluene	<b>6.0</b>	0.92	<b>1.6</b>	0.24	
591-78-6	2-Hexanone	ND	0.92	ND	0.22	
124-48-1	Dibromochloromethane	ND	0.92	ND	0.11	
106-93-4	1,2-Dibromoethane	ND	0.92	ND	0.12	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 21-1F-031512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-003

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01566

Date Collected: 3/16/12  
Date Received: 3/22/12  
Date Analyzed: 3/27/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.65 Final Pressure (psig): 3.73

Canister Dilution Factor: 1.83

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	2.1	0.92	0.45	0.19	
111-65-9	n-Octane	1.1	0.92	0.23	0.20	
108-90-7	Chlorobenzene	ND	0.92	ND	0.20	
179601-23-1	m,p-Xylenes	11	1.8	2.6	0.42	
75-25-2	Bromoform	ND	0.92	ND	0.089	
100-42-5	Styrene	ND	0.92	ND	0.21	
95-47-6	o-Xylene	3.6	0.92	0.83	0.21	
111-84-2	n-Nonane	1.0	0.92	0.20	0.17	
98-82-8	Cumene	ND	0.92	ND	0.19	
80-56-8	alpha-Pinene	5.6	0.92	1.0	0.16	
103-65-1	n-Propylbenzene	ND	0.92	ND	0.19	
622-96-8	4-Ethyltoluene	ND	0.92	ND	0.19	
108-67-8	1,3,5-Trimethylbenzene	0.93	0.92	0.19	0.19	
95-63-6	1,2,4-Trimethylbenzene	2.9	0.92	0.58	0.19	
100-44-7	Benzyl Chloride	ND	0.92	ND	0.18	
541-73-1	1,3-Dichlorobenzene	ND	0.92	ND	0.15	
106-46-7	1,4-Dichlorobenzene	ND	0.92	ND	0.15	
95-50-1	1,2-Dichlorobenzene	ND	0.92	ND	0.15	
5989-27-5	d-Limonene	45	0.92	8.0	0.16	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.92	ND	0.095	
120-82-1	1,2,4-Trichlorobenzene	ND	0.92	ND	0.12	
87-68-3	Hexachlorobutadiene	ND	0.92	ND	0.086	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 23-1F-031512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-004

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00988

Date Collected: 3/16/12  
Date Received: 3/22/12  
Date Analyzed: 3/27 - 3/28/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.050 Liter(s)

Initial Pressure (psig): -4.67 Final Pressure (psig): 3.62

Canister Dilution Factor: 1.83

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	4.0	0.92	2.3	0.53	
75-71-8	Dichlorodifluoromethane (CFC 12)	1.9	0.92	0.39	0.19	
74-87-3	Chloromethane	ND	0.92	ND	0.44	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.92	ND	0.13	
106-99-0	1,3-Butadiene	ND	0.92	ND	0.41	
74-83-9	Bromomethane	ND	0.92	ND	0.24	
75-00-3	Chloroethane	ND	0.92	ND	0.35	
64-17-5	Ethanol	1,400	180	750	97	D
75-05-8	Acetonitrile	ND	0.92	ND	0.55	
107-02-8	Acrolein	ND	3.7	ND	1.6	
67-64-1	Acetone	42	9.2	18	3.9	
75-69-4	Trichlorofluoromethane	1.1	0.92	0.19	0.16	
67-63-0	2-Propanol (Isopropyl Alcohol)	27	1.8	11	0.74	
107-13-1	Acrylonitrile	ND	0.92	ND	0.42	
75-35-4	1,1-Dichloroethene	ND	0.92	ND	0.23	
75-09-2	Methylene Chloride	2.1	0.92	0.62	0.26	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.92	ND	0.29	
76-13-1	Trichlorotrifluoroethane	ND	0.92	ND	0.12	
75-15-0	Carbon Disulfide	ND	9.2	ND	2.9	
156-60-5	trans-1,2-Dichloroethene	ND	0.92	ND	0.23	
75-34-3	1,1-Dichloroethane	ND	0.92	ND	0.23	
1634-04-4	Methyl tert-Butyl Ether	ND	0.92	ND	0.25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 23-1F-031512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-004

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00988

Date Collected: 3/16/12  
Date Received: 3/22/12  
Date Analyzed: 3/27 - 3/28/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.050 Liter(s)

Initial Pressure (psig): -4.67 Final Pressure (psig): 3.62

Canister Dilution Factor: 1.83

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	9.2	ND	2.6	
78-93-3	2-Butanone (MEK)	ND	9.2	ND	3.1	
156-59-2	cis-1,2-Dichloroethene	ND	0.92	ND	0.23	
141-78-6	Ethyl Acetate	7.7	1.8	2.1	0.51	
110-54-3	n-Hexane	ND	0.92	ND	0.26	
67-66-3	Chloroform	3.0	0.92	0.62	0.19	
109-99-9	Tetrahydrofuran (THF)	ND	0.92	ND	0.31	
71-55-6	1,1,1-Trichloroethane	ND	0.92	ND	0.17	
56-23-5	Carbon Tetrachloride	ND	0.92	ND	0.15	
110-82-7	Cyclohexane	ND	1.8	ND	0.53	
78-87-5	1,2-Dichloropropane	ND	0.92	ND	0.20	
75-27-4	Bromodichloromethane	ND	0.92	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.92	ND	0.25	
80-62-6	Methyl Methacrylate	ND	1.8	ND	0.45	
142-82-5	n-Heptane	ND	0.92	ND	0.22	
10061-01-5	cis-1,3-Dichloropropene	ND	0.92	ND	0.20	
108-10-1	4-Methyl-2-pentanone	ND	0.92	ND	0.22	
10061-02-6	trans-1,3-Dichloropropene	ND	0.92	ND	0.20	
108-88-3	Toluene	2.6	0.92	0.69	0.24	
591-78-6	2-Hexanone	ND	0.92	ND	0.22	
124-48-1	Dibromochloromethane	ND	0.92	ND	0.11	
106-93-4	1,2-Dibromoethane	ND	0.92	ND	0.12	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 23-1F-031512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-004

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00988

Date Collected: 3/16/12  
Date Received: 3/22/12  
Date Analyzed: 3/27 - 3/28/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.050 Liter(s)

Initial Pressure (psig): -4.67 Final Pressure (psig): 3.62

Canister Dilution Factor: 1.83

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.92	ND	0.19	
111-65-9	n-Octane	ND	0.92	ND	0.20	
108-90-7	Chlorobenzene	ND	0.92	ND	0.20	
179601-23-1	m,p-Xylenes	ND	1.8	ND	0.42	
75-25-2	Bromoform	ND	0.92	ND	0.089	
100-42-5	Styrene	1.1	0.92	0.25	0.21	
95-47-6	o-Xylene	ND	0.92	ND	0.21	
111-84-2	n-Nonane	ND	0.92	ND	0.17	
98-82-8	Cumene	ND	0.92	ND	0.19	
80-56-8	alpha-Pinene	2.0	0.92	0.35	0.16	
103-65-1	n-Propylbenzene	ND	0.92	ND	0.19	
622-96-8	4-Ethyltoluene	ND	0.92	ND	0.19	
108-67-8	1,3,5-Trimethylbenzene	ND	0.92	ND	0.19	
95-63-6	1,2,4-Trimethylbenzene	ND	0.92	ND	0.19	
100-44-7	Benzyl Chloride	ND	0.92	ND	0.18	
541-73-1	1,3-Dichlorobenzene	ND	0.92	ND	0.15	
106-46-7	1,4-Dichlorobenzene	ND	0.92	ND	0.15	
95-50-1	1,2-Dichlorobenzene	ND	0.92	ND	0.15	
5989-27-5	d-Limonene	20	0.92	3.6	0.16	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.92	ND	0.095	
120-82-1	1,2,4-Trichlorobenzene	ND	0.92	ND	0.12	
87-68-3	Hexachlorobutadiene	ND	0.92	ND	0.086	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 17-1F-031712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-005

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01790

**Date Collected:** 3/16/12  
**Date Received:** 3/22/12  
**Date Analyzed:** 3/27 - 3/28/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.050 Liter(s)

Initial Pressure (psig): -2.83      Final Pressure (psig): 3.61

Canister Dilution Factor: 1.54

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	11	0.77	6.2	0.45	
75-71-8	Dichlorodifluoromethane (CFC 12)	1.9	0.77	0.39	0.16	
74-87-3	Chloromethane	ND	0.77	ND	0.37	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.77	ND	0.11	
106-99-0	1,3-Butadiene	ND	0.77	ND	0.35	
74-83-9	Bromomethane	ND	0.77	ND	0.20	
75-00-3	Chloroethane	ND	0.77	ND	0.29	
64-17-5	Ethanol	1,100	150	560	82	D
75-05-8	Acetonitrile	ND	0.77	ND	0.46	
107-02-8	Acrolein	ND	3.1	ND	1.3	
67-64-1	Acetone	68	7.7	29	3.2	
75-69-4	Trichlorofluoromethane	1.1	0.77	0.19	0.14	
67-63-0	2-Propanol (Isopropyl Alcohol)	62	1.5	25	0.63	
107-13-1	Acrylonitrile	ND	0.77	ND	0.35	
75-35-4	1,1-Dichloroethene	ND	0.77	ND	0.19	
75-09-2	Methylene Chloride	5.1	0.77	1.5	0.22	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.77	ND	0.25	
76-13-1	Trichlorotrifluoroethane	ND	0.77	ND	0.10	
75-15-0	Carbon Disulfide	ND	7.7	ND	2.5	
156-60-5	trans-1,2-Dichloroethene	ND	0.77	ND	0.19	
75-34-3	1,1-Dichloroethane	ND	0.77	ND	0.19	
1634-04-4	Methyl tert-Butyl Ether	ND	0.77	ND	0.21	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 17-1F-031712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-005

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01790

Date Collected: 3/16/12  
Date Received: 3/22/12  
Date Analyzed: 3/27 - 3/28/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.050 Liter(s)

Initial Pressure (psig): -2.83 Final Pressure (psig): 3.61

Canister Dilution Factor: 1.54

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	7.7	ND	2.2	
78-93-3	2-Butanone (MEK)	ND	7.7	ND	2.6	
156-59-2	cis-1,2-Dichloroethene	ND	0.77	ND	0.19	
141-78-6	Ethyl Acetate	17	1.5	4.7	0.43	
110-54-3	n-Hexane	6.0	0.77	1.7	0.22	
67-66-3	Chloroform	2.4	0.77	0.50	0.16	
109-99-9	Tetrahydrofuran (THF)	1.9	0.77	0.66	0.26	
71-55-6	1,1,1-Trichloroethane	ND	0.77	ND	0.14	
56-23-5	Carbon Tetrachloride	1.5	0.77	0.23	0.12	
110-82-7	Cyclohexane	2.6	1.5	0.77	0.45	
78-87-5	1,2-Dichloropropane	ND	0.77	ND	0.17	
75-27-4	Bromodichloromethane	ND	0.77	ND	0.11	
123-91-1	1,4-Dioxane	ND	0.77	ND	0.21	
80-62-6	Methyl Methacrylate	ND	1.5	ND	0.38	
142-82-5	n-Heptane	3.3	0.77	0.81	0.19	
10061-01-5	cis-1,3-Dichloropropene	ND	0.77	ND	0.17	
108-10-1	4-Methyl-2-pentanone	1.1	0.77	0.28	0.19	
10061-02-6	trans-1,3-Dichloropropene	ND	0.77	ND	0.17	
108-88-3	Toluene	28	0.77	7.6	0.20	
591-78-6	2-Hexanone	ND	0.77	ND	0.19	
124-48-1	Dibromochloromethane	ND	0.77	ND	0.090	
106-93-4	1,2-Dibromoethane	ND	0.77	ND	0.10	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 17-1F-031712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-005

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01790

Date Collected: 3/16/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/27 - 3/28/12  
 Volume(s) Analyzed: 1.00 Liter(s)  
 0.050 Liter(s)

Initial Pressure (psig): -2.83      Final Pressure (psig): 3.61

Canister Dilution Factor: 1.54

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	4.0	0.77	0.85	0.16	
111-65-9	n-Octane	0.85	0.77	0.18	0.16	
108-90-7	Chlorobenzene	ND	0.77	ND	0.17	
179601-23-1	m,p-Xylenes	8.4	1.5	1.9	0.35	
75-25-2	Bromoform	ND	0.77	ND	0.075	
100-42-5	Styrene	2.9	0.77	0.68	0.18	
95-47-6	o-Xylene	2.8	0.77	0.64	0.18	
111-84-2	n-Nonane	ND	0.77	ND	0.15	
98-82-8	Cumene	ND	0.77	ND	0.16	
80-56-8	alpha-Pinene	26	0.77	4.7	0.14	
103-65-1	n-Propylbenzene	ND	0.77	ND	0.16	
622-96-8	4-Ethyltoluene	ND	0.77	ND	0.16	
108-67-8	1,3,5-Trimethylbenzene	ND	0.77	ND	0.16	
95-63-6	1,2,4-Trimethylbenzene	1.8	0.77	0.37	0.16	
100-44-7	Benzyl Chloride	ND	0.77	ND	0.15	
541-73-1	1,3-Dichlorobenzene	ND	0.77	ND	0.13	
106-46-7	1,4-Dichlorobenzene	ND	0.77	ND	0.13	
95-50-1	1,2-Dichlorobenzene	ND	0.77	ND	0.13	
5989-27-5	d-Limonene	21	0.77	3.7	0.14	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.77	ND	0.080	
120-82-1	1,2,4-Trichlorobenzene	ND	0.77	ND	0.10	
87-68-3	Hexachlorobutadiene	ND	0.77	ND	0.072	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 20-1F-031712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-006

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01305

Date Collected: 3/18/12  
Date Received: 3/22/12  
Date Analyzed: 3/27 - 3/28/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.050 Liter(s)

Initial Pressure (psig): -1.78 Final Pressure (psig): 3.86

Canister Dilution Factor: 1.44

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	12	0.72	7.0	0.42	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.2	0.72	0.45	0.15	
74-87-3	Chloromethane	ND	0.72	ND	0.35	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.72	ND	0.10	
106-99-0	1,3-Butadiene	ND	0.72	ND	0.33	
74-83-9	Bromomethane	ND	0.72	ND	0.19	
75-00-3	Chloroethane	ND	0.72	ND	0.27	
64-17-5	Ethanol	780	140	410	76	D
75-05-8	Acetonitrile	ND	0.72	ND	0.43	
107-02-8	Acrolein	ND	2.9	ND	1.3	
67-64-1	Acetone	41	7.2	17	3.0	
75-69-4	Trichlorofluoromethane	3.6	0.72	0.64	0.13	
67-63-0	2-Propanol (Isopropyl Alcohol)	75	1.4	31	0.59	
107-13-1	Acrylonitrile	ND	0.72	ND	0.33	
75-35-4	1,1-Dichloroethene	ND	0.72	ND	0.18	
75-09-2	Methylene Chloride	1.0	0.72	0.30	0.21	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.72	ND	0.23	
76-13-1	Trichlorotrifluoroethane	ND	0.72	ND	0.094	
75-15-0	Carbon Disulfide	ND	7.2	ND	2.3	
156-60-5	trans-1,2-Dichloroethene	ND	0.72	ND	0.18	
75-34-3	1,1-Dichloroethane	ND	0.72	ND	0.18	
1634-04-4	Methyl tert-Butyl Ether	ND	0.72	ND	0.20	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 20-1F-031712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-006

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01305

Date Collected: 3/18/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/27 - 3/28/12  
 Volume(s) Analyzed: 1.00 Liter(s)  
 0.050 Liter(s)

Initial Pressure (psig): -1.78      Final Pressure (psig): 3.86

Canister Dilution Factor: 1.44

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	7.2	ND	2.0	
78-93-3	2-Butanone (MEK)	ND	7.2	ND	2.4	
156-59-2	cis-1,2-Dichloroethene	ND	0.72	ND	0.18	
141-78-6	Ethyl Acetate	<b>3.1</b>	1.4	<b>0.87</b>	0.40	
110-54-3	n-Hexane	<b>2.2</b>	0.72	<b>0.62</b>	0.20	
67-66-3	Chloroform	ND	0.72	ND	0.15	
109-99-9	Tetrahydrofuran (THF)	<b>0.86</b>	0.72	<b>0.29</b>	0.24	
71-55-6	1,1,1-Trichloroethane	ND	0.72	ND	0.13	
56-23-5	Carbon Tetrachloride	ND	0.72	ND	0.11	
110-82-7	Cyclohexane	ND	1.4	ND	0.42	
78-87-5	1,2-Dichloropropane	ND	0.72	ND	0.16	
75-27-4	Bromodichloromethane	ND	0.72	ND	0.11	
123-91-1	1,4-Dioxane	ND	0.72	ND	0.20	
80-62-6	Methyl Methacrylate	ND	1.4	ND	0.35	
142-82-5	n-Heptane	<b>1.2</b>	0.72	<b>0.30</b>	0.18	
10061-01-5	cis-1,3-Dichloropropene	ND	0.72	ND	0.16	
108-10-1	4-Methyl-2-pentanone	<b>1.0</b>	0.72	<b>0.25</b>	0.18	
10061-02-6	trans-1,3-Dichloropropene	ND	0.72	ND	0.16	
108-88-3	Toluene	<b>11</b>	0.72	<b>3.0</b>	0.19	
591-78-6	2-Hexanone	ND	0.72	ND	0.18	
124-48-1	Dibromochloromethane	ND	0.72	ND	0.085	
106-93-4	1,2-Dibromoethane	ND	0.72	ND	0.094	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 20-1F-031712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-006

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01305

Date Collected: 3/18/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/27 - 3/28/12  
 Volume(s) Analyzed: 1.00 Liter(s)  
 0.050 Liter(s)

Initial Pressure (psig): -1.78 Final Pressure (psig): 3.86

Canister Dilution Factor: 1.44

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.72	ND	0.15	
111-65-9	n-Octane	<b>0.89</b>	0.72	<b>0.19</b>	0.15	
108-90-7	Chlorobenzene	ND	0.72	ND	0.16	
179601-23-1	m,p-Xylenes	<b>8.7</b>	1.4	<b>2.0</b>	0.33	
75-25-2	Bromoform	ND	0.72	ND	0.070	
100-42-5	Styrene	ND	0.72	ND	0.17	
95-47-6	o-Xylene	<b>2.5</b>	0.72	<b>0.59</b>	0.17	
111-84-2	n-Nonane	<b>3.0</b>	0.72	<b>0.57</b>	0.14	
98-82-8	Cumene	ND	0.72	ND	0.15	
80-56-8	alpha-Pinene	<b>4.7</b>	0.72	<b>0.84</b>	0.13	
103-65-1	n-Propylbenzene	ND	0.72	ND	0.15	
622-96-8	4-Ethyltoluene	ND	0.72	ND	0.15	
108-67-8	1,3,5-Trimethylbenzene	ND	0.72	ND	0.15	
95-63-6	1,2,4-Trimethylbenzene	<b>1.8</b>	0.72	<b>0.37</b>	0.15	
100-44-7	Benzyl Chloride	ND	0.72	ND	0.14	
541-73-1	1,3-Dichlorobenzene	ND	0.72	ND	0.12	
106-46-7	1,4-Dichlorobenzene	ND	0.72	ND	0.12	
95-50-1	1,2-Dichlorobenzene	ND	0.72	ND	0.12	
5989-27-5	d-Limonene	<b>230</b>	14	<b>41</b>	2.6	<b>D</b>
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.72	ND	0.075	
120-82-1	1,2,4-Trichlorobenzene	ND	0.72	ND	0.097	
87-68-3	Hexachlorobutadiene	ND	0.72	ND	0.068	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 28-1F-031912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-007

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00721

Date Collected: 3/20/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 0.30 Liter(s)

Initial Pressure (psig): -4.99 Final Pressure (psig): 3.65

Canister Dilution Factor: 1.89

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	47	3.2	27	1.8	
75-71-8	Dichlorodifluoromethane (CFC 12)	ND	3.2	ND	0.64	
74-87-3	Chloromethane	ND	3.2	ND	1.5	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	3.2	ND	0.45	
106-99-0	1,3-Butadiene	ND	3.2	ND	1.4	
74-83-9	Bromomethane	ND	3.2	ND	0.81	
75-00-3	Chloroethane	ND	3.2	ND	1.2	
64-17-5	Ethanol	660	32	350	17	
75-05-8	Acetonitrile	ND	3.2	ND	1.9	
107-02-8	Acrolein	ND	13	ND	5.5	
67-64-1	Acetone	49	32	20	13	
75-69-4	Trichlorofluoromethane	ND	3.2	ND	0.56	
67-63-0	2-Propanol (Isopropyl Alcohol)	170	6.3	68	2.6	
107-13-1	Acrylonitrile	ND	3.2	ND	1.5	
75-35-4	1,1-Dichloroethene	ND	3.2	ND	0.79	
75-09-2	Methylene Chloride	ND	3.2	ND	0.91	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	3.2	ND	1.0	
76-13-1	Trichlorotrifluoroethane	ND	3.2	ND	0.41	
75-15-0	Carbon Disulfide	ND	32	ND	10	
156-60-5	trans-1,2-Dichloroethene	ND	3.2	ND	0.79	
75-34-3	1,1-Dichloroethane	ND	3.2	ND	0.78	
1634-04-4	Methyl tert-Butyl Ether	ND	3.2	ND	0.87	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 28-1F-031912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-007

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00721

Date Collected: 3/20/12  
Date Received: 3/22/12  
Date Analyzed: 3/27/12  
Volume(s) Analyzed: 0.30 Liter(s)

Initial Pressure (psig): -4.99 Final Pressure (psig): 3.65

Canister Dilution Factor: 1.89

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	32	ND	8.9	
78-93-3	2-Butanone (MEK)	ND	32	ND	11	
156-59-2	cis-1,2-Dichloroethene	ND	3.2	ND	0.79	
141-78-6	Ethyl Acetate	ND	6.3	ND	1.7	
110-54-3	n-Hexane	ND	3.2	ND	0.89	
67-66-3	Chloroform	<b>4.0</b>	3.2	<b>0.82</b>	0.65	
109-99-9	Tetrahydrofuran (THF)	ND	3.2	ND	1.1	
71-55-6	1,1,1-Trichloroethane	ND	3.2	ND	0.58	
56-23-5	Carbon Tetrachloride	ND	3.2	ND	0.50	
110-82-7	Cyclohexane	ND	6.3	ND	1.8	
78-87-5	1,2-Dichloropropane	ND	3.2	ND	0.68	
75-27-4	Bromodichloromethane	ND	3.2	ND	0.47	
123-91-1	1,4-Dioxane	ND	3.2	ND	0.87	
80-62-6	Methyl Methacrylate	ND	6.3	ND	1.5	
142-82-5	n-Heptane	ND	3.2	ND	0.77	
10061-01-5	cis-1,3-Dichloropropene	ND	3.2	ND	0.69	
108-10-1	4-Methyl-2-pentanone	ND	3.2	ND	0.77	
10061-02-6	trans-1,3-Dichloropropene	ND	3.2	ND	0.69	
108-88-3	Toluene	<b>5.0</b>	3.2	<b>1.3</b>	0.84	
591-78-6	2-Hexanone	ND	3.2	ND	0.77	
124-48-1	Dibromochloromethane	ND	3.2	ND	0.37	
106-93-4	1,2-Dibromoethane	ND	3.2	ND	0.41	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 28-1F-031912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-007

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00721

Date Collected: 3/20/12  
Date Received: 3/22/12  
Date Analyzed: 3/27/12  
Volume(s) Analyzed: 0.30 Liter(s)

Initial Pressure (psig): -4.99      Final Pressure (psig): 3.65

Canister Dilution Factor: 1.89

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	3.2	ND	0.66	
111-65-9	n-Octane	ND	3.2	ND	0.67	
108-90-7	Chlorobenzene	ND	3.2	ND	0.68	
179601-23-1	m,p-Xylenes	ND	6.3	ND	1.5	
75-25-2	Bromoform	ND	3.2	ND	0.30	
100-42-5	Styrene	ND	3.2	ND	0.74	
95-47-6	o-Xylene	ND	3.2	ND	0.73	
111-84-2	n-Nonane	ND	3.2	ND	0.60	
98-82-8	Cumene	ND	3.2	ND	0.64	
80-56-8	alpha-Pinene	ND	3.2	ND	0.57	
103-65-1	n-Propylbenzene	ND	3.2	ND	0.64	
622-96-8	4-Ethyltoluene	ND	3.2	ND	0.64	
108-67-8	1,3,5-Trimethylbenzene	ND	3.2	ND	0.64	
95-63-6	1,2,4-Trimethylbenzene	ND	3.2	ND	0.64	
100-44-7	Benzyl Chloride	ND	3.2	ND	0.61	
541-73-1	1,3-Dichlorobenzene	ND	3.2	ND	0.52	
106-46-7	1,4-Dichlorobenzene	ND	3.2	ND	0.52	
95-50-1	1,2-Dichlorobenzene	ND	3.2	ND	0.52	
5989-27-5	d-Limonene	<b>120</b>	3.2	<b>21</b>	0.57	
96-12-8	1,2-Dibromo-3-chloropropane	ND	3.2	ND	0.33	
120-82-1	1,2,4-Trichlorobenzene	ND	3.2	ND	0.42	
87-68-3	Hexachlorobutadiene	ND	3.2	ND	0.30	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 22-1F-031912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-008

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01171

Date Collected: 3/20/12  
Date Received: 3/22/12  
Date Analyzed: 3/28/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.96 Final Pressure (psig): 3.64

Canister Dilution Factor: 1.56

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	4.1	0.78	2.4	0.45	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.0	0.78	0.41	0.16	
74-87-3	Chloromethane	ND	0.78	ND	0.38	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.78	ND	0.11	
106-99-0	1,3-Butadiene	ND	0.78	ND	0.35	
74-83-9	Bromomethane	ND	0.78	ND	0.20	
75-00-3	Chloroethane	ND	0.78	ND	0.30	
64-17-5	Ethanol	130	7.8	71	4.1	
75-05-8	Acetonitrile	ND	0.78	ND	0.46	
107-02-8	Acrolein	ND	3.1	ND	1.4	
67-64-1	Acetone	29	7.8	12	3.3	
75-69-4	Trichlorofluoromethane	1.2	0.78	0.21	0.14	
67-63-0	2-Propanol (Isopropyl Alcohol)	24	1.6	10	0.63	
107-13-1	Acrylonitrile	ND	0.78	ND	0.36	
75-35-4	1,1-Dichloroethene	ND	0.78	ND	0.20	
75-09-2	Methylene Chloride	8.5	0.78	2.5	0.22	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.78	ND	0.25	
76-13-1	Trichlorotrifluoroethane	ND	0.78	ND	0.10	
75-15-0	Carbon Disulfide	ND	7.8	ND	2.5	
156-60-5	trans-1,2-Dichloroethene	ND	0.78	ND	0.20	
75-34-3	1,1-Dichloroethane	ND	0.78	ND	0.19	
1634-04-4	Methyl tert-Butyl Ether	ND	0.78	ND	0.22	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 22-1F-031912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-008

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01171

**Date Collected:** 3/20/12  
**Date Received:** 3/22/12  
**Date Analyzed:** 3/28/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -2.96      Final Pressure (psig): 3.64

Canister Dilution Factor: 1.56

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	7.8	ND	2.2	
78-93-3	2-Butanone (MEK)	ND	7.8	ND	2.6	
156-59-2	cis-1,2-Dichloroethene	ND	0.78	ND	0.20	
141-78-6	Ethyl Acetate	ND	1.6	ND	0.43	
110-54-3	n-Hexane	<b>1.8</b>	0.78	<b>0.50</b>	0.22	
67-66-3	Chloroform	ND	0.78	ND	0.16	
109-99-9	Tetrahydrofuran (THF)	ND	0.78	ND	0.26	
71-55-6	1,1,1-Trichloroethane	ND	0.78	ND	0.14	
56-23-5	Carbon Tetrachloride	ND	0.78	ND	0.12	
110-82-7	Cyclohexane	ND	1.6	ND	0.45	
78-87-5	1,2-Dichloropropane	ND	0.78	ND	0.17	
75-27-4	Bromodichloromethane	ND	0.78	ND	0.12	
123-91-1	1,4-Dioxane	ND	0.78	ND	0.22	
80-62-6	Methyl Methacrylate	ND	1.6	ND	0.38	
142-82-5	n-Heptane	<b>0.91</b>	0.78	<b>0.22</b>	0.19	
10061-01-5	cis-1,3-Dichloropropene	ND	0.78	ND	0.17	
108-10-1	4-Methyl-2-pentanone	ND	0.78	ND	0.19	
10061-02-6	trans-1,3-Dichloropropene	ND	0.78	ND	0.17	
108-88-3	Toluene	<b>4.5</b>	0.78	<b>1.2</b>	0.21	
591-78-6	2-Hexanone	ND	0.78	ND	0.19	
124-48-1	Dibromochloromethane	ND	0.78	ND	0.092	
106-93-4	1,2-Dibromoethane	ND	0.78	ND	0.10	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 22-1F-031912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-008

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01171

Date Collected: 3/20/12  
Date Received: 3/22/12  
Date Analyzed: 3/28/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.96 Final Pressure (psig): 3.64

Canister Dilution Factor: 1.56

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.78	ND	0.16	
111-65-9	n-Octane	ND	0.78	ND	0.17	
108-90-7	Chlorobenzene	ND	0.78	ND	0.17	
179601-23-1	m,p-Xylenes	ND	1.6	ND	0.36	
75-25-2	Bromoform	ND	0.78	ND	0.075	
100-42-5	Styrene	ND	0.78	ND	0.18	
95-47-6	o-Xylene	ND	0.78	ND	0.18	
111-84-2	n-Nonane	<b>2.8</b>	0.78	<b>0.54</b>	0.15	
98-82-8	Cumene	ND	0.78	ND	0.16	
80-56-8	alpha-Pinene	<b>0.98</b>	0.78	<b>0.18</b>	0.14	
103-65-1	n-Propylbenzene	ND	0.78	ND	0.16	
622-96-8	4-Ethyltoluene	ND	0.78	ND	0.16	
108-67-8	1,3,5-Trimethylbenzene	ND	0.78	ND	0.16	
95-63-6	1,2,4-Trimethylbenzene	<b>0.96</b>	0.78	<b>0.19</b>	0.16	
100-44-7	Benzyl Chloride	ND	0.78	ND	0.15	
541-73-1	1,3-Dichlorobenzene	ND	0.78	ND	0.13	
106-46-7	1,4-Dichlorobenzene	ND	0.78	ND	0.13	
95-50-1	1,2-Dichlorobenzene	ND	0.78	ND	0.13	
5989-27-5	d-Limonene	<b>5.0</b>	0.78	<b>0.89</b>	0.14	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.78	ND	0.081	
120-82-1	1,2,4-Trichlorobenzene	ND	0.78	ND	0.11	
87-68-3	Hexachlorobutadiene	ND	0.78	ND	0.073	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 26-1F-031912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-009

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01783

Date Collected: 3/20/12  
Date Received: 3/22/12  
Date Analyzed: 3/28/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.72 Final Pressure (psig): 3.54

Canister Dilution Factor: 1.66

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	1.7	0.83	0.99	0.48	
75-71-8	Dichlorodifluoromethane (CFC 12)	1.9	0.83	0.39	0.17	
74-87-3	Chloromethane	ND	0.83	ND	0.40	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.83	ND	0.12	
106-99-0	1,3-Butadiene	ND	0.83	ND	0.38	
74-83-9	Bromomethane	ND	0.83	ND	0.21	
75-00-3	Chloroethane	ND	0.83	ND	0.31	
64-17-5	Ethanol	280	8.3	150	4.4	
75-05-8	Acetonitrile	ND	0.83	ND	0.49	
107-02-8	Acrolein	ND	3.3	ND	1.4	
67-64-1	Acetone	36	8.3	15	3.5	
75-69-4	Trichlorofluoromethane	5.3	0.83	0.95	0.15	
67-63-0	2-Propanol (Isopropyl Alcohol)	11	1.7	4.4	0.68	M1
107-13-1	Acrylonitrile	ND	0.83	ND	0.38	
75-35-4	1,1-Dichloroethene	ND	0.83	ND	0.21	
75-09-2	Methylene Chloride	38	0.83	11	0.24	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.83	ND	0.27	
76-13-1	Trichlorotrifluoroethane	ND	0.83	ND	0.11	
75-15-0	Carbon Disulfide	ND	8.3	ND	2.7	
156-60-5	trans-1,2-Dichloroethene	ND	0.83	ND	0.21	
75-34-3	1,1-Dichloroethane	ND	0.83	ND	0.21	
1634-04-4	Methyl tert-Butyl Ether	ND	0.83	ND	0.23	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

M1 = Matrix interference due to coelution with a non-target compound; results may be biased high.



## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 26-1F-031912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-009

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01783

**Date Collected:** 3/20/12  
**Date Received:** 3/22/12  
**Date Analyzed:** 3/28/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -3.72      Final Pressure (psig): 3.54

Canister Dilution Factor: 1.66

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	8.3	ND	2.4	
78-93-3	2-Butanone (MEK)	8.3	8.3	2.8	2.8	
156-59-2	cis-1,2-Dichloroethene	ND	0.83	ND	0.21	
141-78-6	Ethyl Acetate	3.5	1.7	0.97	0.46	
110-54-3	n-Hexane	0.89	0.83	0.25	0.24	
67-66-3	Chloroform	ND	0.83	ND	0.17	
109-99-9	Tetrahydrofuran (THF)	ND	0.83	ND	0.28	
71-55-6	1,1,1-Trichloroethane	ND	0.83	ND	0.15	
56-23-5	Carbon Tetrachloride	ND	0.83	ND	0.13	
110-82-7	Cyclohexane	ND	1.7	ND	0.48	
78-87-5	1,2-Dichloropropane	ND	0.83	ND	0.18	
75-27-4	Bromodichloromethane	ND	0.83	ND	0.12	
123-91-1	1,4-Dioxane	ND	0.83	ND	0.23	
80-62-6	Methyl Methacrylate	ND	1.7	ND	0.41	
142-82-5	n-Heptane	ND	0.83	ND	0.20	
10061-01-5	cis-1,3-Dichloropropene	ND	0.83	ND	0.18	
108-10-1	4-Methyl-2-pentanone	ND	0.83	ND	0.20	
10061-02-6	trans-1,3-Dichloropropene	ND	0.83	ND	0.18	
108-88-3	Toluene	2.1	0.83	0.57	0.22	
591-78-6	2-Hexanone	ND	0.83	ND	0.20	
124-48-1	Dibromochloromethane	ND	0.83	ND	0.097	
106-93-4	1,2-Dibromoethane	ND	0.83	ND	0.11	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 26-1F-031912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-009

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01783

Date Collected: 3/20/12  
Date Received: 3/22/12  
Date Analyzed: 3/28/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.72 Final Pressure (psig): 3.54

Canister Dilution Factor: 1.66

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.83	ND	0.17	
111-65-9	n-Octane	ND	0.83	ND	0.18	
108-90-7	Chlorobenzene	ND	0.83	ND	0.18	
179601-23-1	m,p-Xylenes	ND	1.7	ND	0.38	
75-25-2	Bromoform	ND	0.83	ND	0.080	
100-42-5	Styrene	ND	0.83	ND	0.20	
95-47-6	o-Xylene	ND	0.83	ND	0.19	
111-84-2	n-Nonane	ND	0.83	ND	0.16	
98-82-8	Cumene	ND	0.83	ND	0.17	
80-56-8	alpha-Pinene	1.2	0.83	0.21	0.15	
103-65-1	n-Propylbenzene	ND	0.83	ND	0.17	
622-96-8	4-Ethyltoluene	ND	0.83	ND	0.17	
108-67-8	1,3,5-Trimethylbenzene	ND	0.83	ND	0.17	
95-63-6	1,2,4-Trimethylbenzene	ND	0.83	ND	0.17	
100-44-7	Benzyl Chloride	ND	0.83	ND	0.16	
541-73-1	1,3-Dichlorobenzene	ND	0.83	ND	0.14	
106-46-7	1,4-Dichlorobenzene	ND	0.83	ND	0.14	
95-50-1	1,2-Dichlorobenzene	ND	0.83	ND	0.14	
5989-27-5	d-Limonene	3.2	0.83	0.58	0.15	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.83	ND	0.086	
120-82-1	1,2,4-Trichlorobenzene	ND	0.83	ND	0.11	
87-68-3	Hexachlorobutadiene	ND	0.83	ND	0.078	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P120327-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	ND	0.50	ND	0.29	
75-71-8	Dichlorodifluoromethane (CFC 12)	ND	0.50	ND	0.10	
74-87-3	Chloromethane	ND	0.50	ND	0.24	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.50	ND	0.072	
106-99-0	1,3-Butadiene	ND	0.50	ND	0.23	
74-83-9	Bromomethane	ND	0.50	ND	0.13	
75-00-3	Chloroethane	ND	0.50	ND	0.19	
64-17-5	Ethanol	ND	5.0	ND	2.7	
75-05-8	Acetonitrile	ND	0.50	ND	0.30	
107-02-8	Acrolein	ND	2.0	ND	0.87	
67-64-1	Acetone	ND	5.0	ND	2.1	
75-69-4	Trichlorofluoromethane	ND	0.50	ND	0.089	
67-63-0	2-Propanol (Isopropyl Alcohol)	ND	1.0	ND	0.41	
107-13-1	Acrylonitrile	ND	0.50	ND	0.23	
75-35-4	1,1-Dichloroethene	ND	0.50	ND	0.13	
75-09-2	Methylene Chloride	ND	0.50	ND	0.14	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.50	ND	0.16	
76-13-1	Trichlorotrifluoroethane	ND	0.50	ND	0.065	
75-15-0	Carbon Disulfide	ND	5.0	ND	1.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.50	ND	0.13	
75-34-3	1,1-Dichloroethane	ND	0.50	ND	0.12	
1634-04-4	Methyl tert-Butyl Ether	ND	0.50	ND	0.14	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P120327-MB

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/27/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	5.0	ND	1.4	
78-93-3	2-Butanone (MEK)	ND	5.0	ND	1.7	
156-59-2	cis-1,2-Dichloroethene	ND	0.50	ND	0.13	
141-78-6	Ethyl Acetate	ND	1.0	ND	0.28	
110-54-3	n-Hexane	ND	0.50	ND	0.14	
67-66-3	Chloroform	ND	0.50	ND	0.10	
109-99-9	Tetrahydrofuran (THF)	ND	0.50	ND	0.17	
71-55-6	1,1,1-Trichloroethane	ND	0.50	ND	0.092	
56-23-5	Carbon Tetrachloride	ND	0.50	ND	0.080	
110-82-7	Cyclohexane	ND	1.0	ND	0.29	
78-87-5	1,2-Dichloropropane	ND	0.50	ND	0.11	
75-27-4	Bromodichloromethane	ND	0.50	ND	0.075	
123-91-1	1,4-Dioxane	ND	0.50	ND	0.14	
80-62-6	Methyl Methacrylate	ND	1.0	ND	0.24	
142-82-5	n-Heptane	ND	0.50	ND	0.12	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-10-1	4-Methyl-2-pentanone	ND	0.50	ND	0.12	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-88-3	Toluene	ND	0.50	ND	0.13	
591-78-6	2-Hexanone	ND	0.50	ND	0.12	
124-48-1	Dibromochloromethane	ND	0.50	ND	0.059	
106-93-4	1,2-Dibromoethane	ND	0.50	ND	0.065	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P120327-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 3/27/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.50	ND	0.11	
111-65-9	n-Octane	ND	0.50	ND	0.11	
108-90-7	Chlorobenzene	ND	0.50	ND	0.11	
179601-23-1	m,p-Xylenes	ND	1.0	ND	0.23	
75-25-2	Bromoform	ND	0.50	ND	0.048	
100-42-5	Styrene	ND	0.50	ND	0.12	
95-47-6	o-Xylene	ND	0.50	ND	0.12	
111-84-2	n-Nonane	ND	0.50	ND	0.095	
98-82-8	Cumene	ND	0.50	ND	0.10	
80-56-8	alpha-Pinene	ND	0.50	ND	0.090	
103-65-1	n-Propylbenzene	ND	0.50	ND	0.10	
622-96-8	4-Ethyltoluene	ND	0.50	ND	0.10	
108-67-8	1,3,5-Trimethylbenzene	ND	0.50	ND	0.10	
95-63-6	1,2,4-Trimethylbenzene	ND	0.50	ND	0.10	
100-44-7	Benzyl Chloride	ND	0.50	ND	0.097	
541-73-1	1,3-Dichlorobenzene	ND	0.50	ND	0.083	
106-46-7	1,4-Dichlorobenzene	ND	0.50	ND	0.083	
95-50-1	1,2-Dichlorobenzene	ND	0.50	ND	0.083	
5989-27-5	d-Limonene	ND	0.50	ND	0.090	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.50	ND	0.052	
120-82-1	1,2,4-Trichlorobenzene	ND	0.50	ND	0.067	
87-68-3	Hexachlorobutadiene	ND	0.50	ND	0.047	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P120328-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/28/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	ND	0.50	ND	0.29	
75-71-8	Dichlorodifluoromethane (CFC 12)	ND	0.50	ND	0.10	
74-87-3	Chloromethane	ND	0.50	ND	0.24	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.50	ND	0.072	
106-99-0	1,3-Butadiene	ND	0.50	ND	0.23	
74-83-9	Bromomethane	ND	0.50	ND	0.13	
75-00-3	Chloroethane	ND	0.50	ND	0.19	
64-17-5	Ethanol	ND	5.0	ND	2.7	
75-05-8	Acetonitrile	ND	0.50	ND	0.30	
107-02-8	Acrolein	ND	2.0	ND	0.87	
67-64-1	Acetone	ND	5.0	ND	2.1	
75-69-4	Trichlorofluoromethane	ND	0.50	ND	0.089	
67-63-0	2-Propanol (Isopropyl Alcohol)	ND	1.0	ND	0.41	
107-13-1	Acrylonitrile	ND	0.50	ND	0.23	
75-35-4	1,1-Dichloroethene	ND	0.50	ND	0.13	
75-09-2	Methylene Chloride	ND	0.50	ND	0.14	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.50	ND	0.16	
76-13-1	Trichlorotrifluoroethane	ND	0.50	ND	0.065	
75-15-0	Carbon Disulfide	ND	5.0	ND	1.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.50	ND	0.13	
75-34-3	1,1-Dichloroethane	ND	0.50	ND	0.12	
1634-04-4	Methyl tert-Butyl Ether	ND	0.50	ND	0.14	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P120328-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/28/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	5.0	ND	1.4	
78-93-3	2-Butanone (MEK)	ND	5.0	ND	1.7	
156-59-2	cis-1,2-Dichloroethene	ND	0.50	ND	0.13	
141-78-6	Ethyl Acetate	ND	1.0	ND	0.28	
110-54-3	n-Hexane	ND	0.50	ND	0.14	
67-66-3	Chloroform	ND	0.50	ND	0.10	
109-99-9	Tetrahydrofuran (THF)	ND	0.50	ND	0.17	
71-55-6	1,1,1-Trichloroethane	ND	0.50	ND	0.092	
56-23-5	Carbon Tetrachloride	ND	0.50	ND	0.080	
110-82-7	Cyclohexane	ND	1.0	ND	0.29	
78-87-5	1,2-Dichloropropane	ND	0.50	ND	0.11	
75-27-4	Bromodichloromethane	ND	0.50	ND	0.075	
123-91-1	1,4-Dioxane	ND	0.50	ND	0.14	
80-62-6	Methyl Methacrylate	ND	1.0	ND	0.24	
142-82-5	n-Heptane	ND	0.50	ND	0.12	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-10-1	4-Methyl-2-pentanone	ND	0.50	ND	0.12	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-88-3	Toluene	ND	0.50	ND	0.13	
591-78-6	2-Hexanone	ND	0.50	ND	0.12	
124-48-1	Dibromochloromethane	ND	0.50	ND	0.059	
106-93-4	1,2-Dibromoethane	ND	0.50	ND	0.065	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P120328-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/28/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.50	ND	0.11	
111-65-9	n-Octane	ND	0.50	ND	0.11	
108-90-7	Chlorobenzene	ND	0.50	ND	0.11	
179601-23-1	m,p-Xylenes	ND	1.0	ND	0.23	
75-25-2	Bromoform	ND	0.50	ND	0.048	
100-42-5	Styrene	ND	0.50	ND	0.12	
95-47-6	o-Xylene	ND	0.50	ND	0.12	
111-84-2	n-Nonane	ND	0.50	ND	0.095	
98-82-8	Cumene	ND	0.50	ND	0.10	
80-56-8	alpha-Pinene	ND	0.50	ND	0.090	
103-65-1	n-Propylbenzene	ND	0.50	ND	0.10	
622-96-8	4-Ethyltoluene	ND	0.50	ND	0.10	
108-67-8	1,3,5-Trimethylbenzene	ND	0.50	ND	0.10	
95-63-6	1,2,4-Trimethylbenzene	ND	0.50	ND	0.10	
100-44-7	Benzyl Chloride	ND	0.50	ND	0.097	
541-73-1	1,3-Dichlorobenzene	ND	0.50	ND	0.083	
106-46-7	1,4-Dichlorobenzene	ND	0.50	ND	0.083	
95-50-1	1,2-Dichlorobenzene	ND	0.50	ND	0.083	
5989-27-5	d-Limonene	ND	0.50	ND	0.090	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.50	ND	0.052	
120-82-1	1,2,4-Trichlorobenzene	ND	0.50	ND	0.067	
87-68-3	Hexachlorobutadiene	ND	0.50	ND	0.047	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister(s)  
 Test Notes:

Date(s) Collected: 3/15 - 3/20/12  
 Date(s) Received: 3/22/12  
 Date(s) Analyzed: 3/27 - 3/28/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		Percent Recovered	Percent Recovered	Percent Recovered		
Method Blank	P120327-MB	105	99	97	70-130	
Method Blank	P120328-MB	103	99	100	70-130	
Lab Control Sample	P120327-LCS	106	98	97	70-130	
Lab Control Sample	P120328-LCS	103	99	100	70-130	
19-1F-031412	P1201121-001	106	98	99	70-130	
53-1F-031412	P1201121-002	106	97	99	70-130	
53-1F-031412	P1201121-002DUP	105	99	96	70-130	
21-1F-031512	P1201121-003	103	100	96	70-130	
23-1F-031512	P1201121-004	102	99	98	70-130	
17-1F-031712	P1201121-005	102	98	98	70-130	
20-1F-031712	P1201121-006	103	99	99	70-130	
28-1F-031912	P1201121-007	102	99	101	70-130	
22-1F-031912	P1201121-008	102	99	100	70-130	
26-1F-031912	P1201121-009	102	99	100	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.

Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.

LABORATORY CONTROL SAMPLE SUMMARY

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**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P120327-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/27/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
115-07-1	Propene	198	187	94	59-137	
75-71-8	Dichlorodifluoromethane (CFC 12)	196	157	80	63-115	
74-87-3	Chloromethane	190	153	81	59-124	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	200	153	77	65-113	
106-99-0	1,3-Butadiene	204	181	89	60-138	
74-83-9	Bromomethane	194	158	81	69-129	
75-00-3	Chloroethane	196	158	81	60-120	
64-17-5	Ethanol	928	777	84	58-121	
75-05-8	Acetonitrile	194	177	91	64-129	
107-02-8	Acrolein	198	163	82	54-127	
67-64-1	Acetone	1,010	830	82	59-114	
75-69-4	Trichlorofluoromethane	202	160	79	66-108	
67-63-0	2-Propanol (Isopropyl Alcohol)	382	293	77	50-113	
107-13-1	Acrylonitrile	198	183	92	72-135	
75-35-4	1,1-Dichloroethene	212	179	84	70-117	
75-09-2	Methylene Chloride	206	165	80	61-108	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	208	200	96	70-131	
76-13-1	Trichlorotrifluoroethane	206	171	83	70-113	
75-15-0	Carbon Disulfide	208	168	81	65-112	
156-60-5	trans-1,2-Dichloroethene	196	173	88	71-119	
75-34-3	1,1-Dichloroethane	200	177	89	71-116	
1634-04-4	Methyl tert-Butyl Ether	198	177	89	67-116	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result. Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P120327-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/27/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
108-05-4	Vinyl Acetate	952	876	92	59-142	
78-93-3	2-Butanone (MEK)	206	184	89	68-125	
156-59-2	cis-1,2-Dichloroethene	206	179	87	69-119	
141-78-6	Ethyl Acetate	398	355	89	63-130	
110-54-3	n-Hexane	198	164	83	57-120	
67-66-3	Chloroform	214	176	82	69-111	
109-99-9	Tetrahydrofuran (THF)	202	171	85	57-123	
71-55-6	1,1,1-Trichloroethane	198	168	85	73-119	
56-23-5	Carbon Tetrachloride	202	175	87	74-129	
110-82-7	Cyclohexane	390	320	82	70-113	
78-87-5	1,2-Dichloropropane	198	166	84	69-118	
75-27-4	Bromodichloromethane	198	171	86	75-124	
123-91-1	1,4-Dioxane	200	169	85	71-123	
80-62-6	Methyl Methacrylate	400	354	89	72-127	
142-82-5	n-Heptane	196	168	86	68-120	
10061-01-5	cis-1,3-Dichloropropene	188	166	88	71-130	
108-10-1	4-Methyl-2-pentanone	204	185	91	69-130	
10061-02-6	trans-1,3-Dichloropropene	210	192	91	76-133	
108-88-3	Toluene	202	161	80	67-111	
591-78-6	2-Hexanone	222	193	87	70-123	
124-48-1	Dibromochloromethane	206	176	85	75-129	
106-93-4	1,2-Dibromoethane	200	162	81	73-122	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P120327-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/27/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
123-86-4	n-Butyl Acetate	220	188	85	68-132	
111-65-9	n-Octane	200	176	88	68-116	
108-90-7	Chlorobenzene	202	157	78	69-113	
179601-23-1	m,p-Xylenes	392	323	82	70-116	
75-25-2	Bromoform	208	169	81	69-127	
100-42-5	Styrene	200	168	84	71-125	
95-47-6	o-Xylene	194	158	81	70-116	
111-84-2	n-Nonane	196	167	85	68-116	
98-82-8	Cumene	190	155	82	70-116	
80-56-8	alpha-Pinene	186	157	84	71-119	
103-65-1	n-Propylbenzene	192	155	81	71-119	
622-96-8	4-Ethyltoluene	198	159	80	71-119	
108-67-8	1,3,5-Trimethylbenzene	200	163	82	71-121	
95-63-6	1,2,4-Trimethylbenzene	194	163	84	73-127	
100-44-7	Benzyl Chloride	200	184	92	65-137	
541-73-1	1,3-Dichlorobenzene	200	148	74	68-123	
106-46-7	1,4-Dichlorobenzene	206	145	70	65-120	
95-50-1	1,2-Dichlorobenzene	198	147	74	67-121	
5989-27-5	d-Limonene	200	177	89	67-130	
96-12-8	1,2-Dibromo-3-chloropropane	196	162	83	72-133	
120-82-1	1,2,4-Trichlorobenzene	196	149	76	62-133	
87-68-3	Hexachlorobutadiene	202	152	75	60-128	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P120328-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/28/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
115-07-1	Propene	198	180	91	59-137	
75-71-8	Dichlorodifluoromethane (CFC 12)	196	154	79	63-115	
74-87-3	Chloromethane	190	147	77	59-124	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	200	152	76	65-113	
106-99-0	1,3-Butadiene	204	180	88	60-138	
74-83-9	Bromomethane	194	156	80	69-129	
75-00-3	Chloroethane	196	156	80	60-120	
64-17-5	Ethanol	928	755	81	58-121	
75-05-8	Acetonitrile	194	173	89	64-129	
107-02-8	Acrolein	198	158	80	54-127	
67-64-1	Acetone	1,010	805	80	59-114	
75-69-4	Trichlorofluoromethane	202	157	78	66-108	
67-63-0	2-Propanol (Isopropyl Alcohol)	382	289	76	50-113	
107-13-1	Acrylonitrile	198	179	90	72-135	
75-35-4	1,1-Dichloroethene	212	177	83	70-117	
75-09-2	Methylene Chloride	206	162	79	61-108	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	208	192	92	70-131	
76-13-1	Trichlorotrifluoroethane	206	170	83	70-113	
75-15-0	Carbon Disulfide	208	166	80	65-112	
156-60-5	trans-1,2-Dichloroethene	196	166	85	71-119	
75-34-3	1,1-Dichloroethane	200	171	86	71-116	
1634-04-4	Methyl tert-Butyl Ether	198	172	87	67-116	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result. Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P120328-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 3/28/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
108-05-4	Vinyl Acetate	952	857	90	59-142	
78-93-3	2-Butanone (MEK)	206	181	88	68-125	
156-59-2	cis-1,2-Dichloroethene	206	173	84	69-119	
141-78-6	Ethyl Acetate	398	346	87	63-130	
110-54-3	n-Hexane	198	161	81	57-120	
67-66-3	Chloroform	214	171	80	69-111	
109-99-9	Tetrahydrofuran (THF)	202	166	82	57-123	
71-55-6	1,1,1-Trichloroethane	198	164	83	73-119	
56-23-5	Carbon Tetrachloride	202	172	85	74-129	
110-82-7	Cyclohexane	390	314	81	70-113	
78-87-5	1,2-Dichloropropane	198	161	81	69-118	
75-27-4	Bromodichloromethane	198	165	83	75-124	
123-91-1	1,4-Dioxane	200	166	83	71-123	
80-62-6	Methyl Methacrylate	400	350	88	72-127	
142-82-5	n-Heptane	196	165	84	68-120	
10061-01-5	cis-1,3-Dichloropropene	188	162	86	71-130	
108-10-1	4-Methyl-2-pentanone	204	180	88	69-130	
10061-02-6	trans-1,3-Dichloropropene	210	186	89	76-133	
108-88-3	Toluene	202	158	78	67-111	
591-78-6	2-Hexanone	222	186	84	70-123	
124-48-1	Dibromochloromethane	206	176	85	75-129	
106-93-4	1,2-Dibromoethane	200	161	81	73-122	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P120328-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/28/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
123-86-4	n-Butyl Acetate	220	180	82	68-132	
111-65-9	n-Octane	200	172	86	68-116	
108-90-7	Chlorobenzene	202	157	78	69-113	
179601-23-1	m,p-Xylenes	392	319	81	70-116	
75-25-2	Bromoform	208	169	81	69-127	
100-42-5	Styrene	200	167	84	71-125	
95-47-6	o-Xylene	194	157	81	70-116	
111-84-2	n-Nonane	196	161	82	68-116	
98-82-8	Cumene	190	153	81	70-116	
80-56-8	alpha-Pinene	186	155	83	71-119	
103-65-1	n-Propylbenzene	192	153	80	71-119	
622-96-8	4-Ethyltoluene	198	157	79	71-119	
108-67-8	1,3,5-Trimethylbenzene	200	162	81	71-121	
95-63-6	1,2,4-Trimethylbenzene	194	161	83	73-127	
100-44-7	Benzyl Chloride	200	183	92	65-137	
541-73-1	1,3-Dichlorobenzene	200	149	75	68-123	
106-46-7	1,4-Dichlorobenzene	206	146	71	65-120	
95-50-1	1,2-Dichlorobenzene	198	148	75	67-121	
5989-27-5	d-Limonene	200	172	86	67-130	
96-12-8	1,2-Dibromo-3-chloropropane	196	164	84	72-133	
120-82-1	1,2,4-Trichlorobenzene	196	150	77	62-133	
87-68-3	Hexachlorobutadiene	202	153	76	60-128	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY DUPLICATE SUMMARY RESULTS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 53-1F-031412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-002DUP

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01478

Date Collected: 3/15/12  
Date Received: 3/22/12  
Date Analyzed: 3/27/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.95

Final Pressure (psig): 3.66

Canister Dilution Factor: 1.44

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
Propene	ND	ND	ND	ND	-	-	25	
Dichlorodifluoromethane (CFC 12)	1.96	0.396	1.91	0.386	1.935	<b>3</b>	25	
Chloromethane	ND	ND	ND	ND	-	-	25	
1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	ND	ND	ND	-	-	25	
1,3-Butadiene	ND	ND	ND	ND	-	-	25	
Bromomethane	ND	ND	ND	ND	-	-	25	
Chloroethane	ND	ND	ND	ND	-	-	25	
Ethanol	201	107	198	105	199.5	<b>2</b>	25	
Acetonitrile	ND	ND	ND	ND	-	-	25	
Acrolein	ND	ND	ND	ND	-	-	25	
Acetone	23.9	10.1	24.1	10.1	24	<b>0.8</b>	25	
Trichlorofluoromethane	1.33	0.238	1.28	0.228	1.305	<b>4</b>	25	
2-Propanol (Isopropyl Alcohol)	ND	ND	1.51	0.615	-	-	25	
Acrylonitrile	ND	ND	ND	ND	-	-	25	
1,1-Dichloroethene	ND	ND	ND	ND	-	-	25	
Methylene Chloride	18.6	5.37	18.5	5.33	18.55	<b>0.5</b>	25	
3-Chloro-1-propene (Allyl Chloride)	ND	ND	ND	ND	-	-	25	
Trichlorotrifluoroethane	ND	ND	ND	ND	-	-	25	
Carbon Disulfide	ND	ND	ND	ND	-	-	25	
trans-1,2-Dichloroethene	ND	ND	ND	ND	-	-	25	
1,1-Dichloroethane	ND	ND	ND	ND	-	-	25	
Methyl tert-Butyl Ether	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.



## LABORATORY DUPLICATE SUMMARY RESULTS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 53-1F-031412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-002DUP

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01478

**Date Collected:** 3/15/12  
**Date Received:** 3/22/12  
**Date Analyzed:** 3/27/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -1.95

Final Pressure (psig): 3.66

Canister Dilution Factor: 1.44

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
Vinyl Acetate	ND	ND	ND	ND	-	-	25	
2-Butanone (MEK)	ND	ND	ND	ND	-	-	25	
cis-1,2-Dichloroethene	ND	ND	ND	ND	-	-	25	
Ethyl Acetate	16.3	4.52	16.3	4.54	16.3	0	25	
n-Hexane	ND	ND	ND	ND	-	-	25	
Chloroform	ND	ND	ND	ND	-	-	25	
Tetrahydrofuran (THF)	ND	ND	ND	ND	-	-	25	
1,1,1-Trichloroethane	ND	ND	ND	ND	-	-	25	
Carbon Tetrachloride	ND	ND	ND	ND	-	-	25	
Cyclohexane	ND	ND	ND	ND	-	-	25	
1,2-Dichloropropane	ND	ND	ND	ND	-	-	25	
Bromodichloromethane	ND	ND	ND	ND	-	-	25	
1,4-Dioxane	ND	ND	ND	ND	-	-	25	
Methyl Methacrylate	ND	ND	ND	ND	-	-	25	
n-Heptane	ND	ND	ND	ND	-	-	25	
cis-1,3-Dichloropropene	ND	ND	ND	ND	-	-	25	
4-Methyl-2-pentanone	ND	ND	ND	ND	-	-	25	
trans-1,3-Dichloropropene	ND	ND	ND	ND	-	-	25	
Toluene	5.48	1.45	5.61	1.49	5.545	2	25	
2-Hexanone	ND	ND	ND	ND	-	-	25	
Dibromochloromethane	ND	ND	ND	ND	-	-	25	
1,2-Dibromoethane	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

LABORATORY DUPLICATE SUMMARY RESULTS

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 53-1F-031412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-002DUP

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01478

Date Collected: 3/15/12  
Date Received: 3/22/12  
Date Analyzed: 3/27/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.95

Final Pressure (psig): 3.66

Canister Dilution Factor: 1.44

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
n-Butyl Acetate	2.64	0.557	2.69	0.567	2.665	2	25	
n-Octane	ND	ND	ND	ND	-	-	25	
Chlorobenzene	ND	ND	ND	ND	-	-	25	
m,p-Xylenes	1.81	0.416	1.83	0.421	1.82	1	25	
Bromoform	ND	ND	ND	ND	-	-	25	
Styrene	ND	ND	ND	ND	-	-	25	
o-Xylene	ND	ND	ND	ND	-	-	25	
n-Nonane	1.42	0.271	1.49	0.284	1.455	5	25	
Cumene	ND	ND	ND	ND	-	-	25	
alpha-Pinene	5.15	0.925	5.31	0.953	5.23	3	25	
n-Propylbenzene	ND	ND	ND	ND	-	-	25	
4-Ethyltoluene	ND	ND	ND	ND	-	-	25	
1,3,5-Trimethylbenzene	ND	ND	ND	ND	-	-	25	
1,2,4-Trimethylbenzene	ND	ND	ND	ND	-	-	25	
Benzyl Chloride	ND	ND	ND	ND	-	-	25	
1,3-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
1,4-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
1,2-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
d-Limonene	47.3	8.49	48.4	8.68	47.85	2	25	
1,2-Dibromo-3-chloropropane	ND	ND	ND	ND	-	-	25	
1,2,4-Trichlorobenzene	ND	ND	ND	ND	-	-	25	
Hexachlorobutadiene	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 19-1F-031412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-001

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01817

Date Collected: 3/15/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/31/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.81      Final Pressure (psig): 3.51

Canister Dilution Factor: 1.67

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.042	ND	0.016	
107-06-2	1,2-Dichloroethane	<b>0.12</b>	0.042	<b>0.030</b>	0.010	
71-43-2	Benzene	<b>0.70</b>	0.13	<b>0.22</b>	0.039	
79-01-6	Trichloroethene	ND	0.042	ND	0.0078	
79-00-5	1,1,2-Trichloroethane	ND	0.17	ND	0.031	
127-18-4	Tetrachloroethene	<b>0.070</b>	0.042	<b>0.010</b>	0.0062	
100-41-4	Ethylbenzene	<b>0.76</b>	0.17	<b>0.18</b>	0.038	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.042	ND	0.0061	
91-20-3	Naphthalene	ND	0.17	ND	0.032	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 53-1F-031412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-002

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01478

Date Collected: 3/15/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/31/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.95      Final Pressure (psig): 3.66

Canister Dilution Factor: 1.44

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.036	ND	0.014	
107-06-2	1,2-Dichloroethane	<b>0.14</b>	0.036	<b>0.034</b>	0.0089	
71-43-2	Benzene	<b>0.72</b>	0.11	<b>0.22</b>	0.034	
79-01-6	Trichloroethene	ND	0.036	ND	0.0067	
79-00-5	1,1,2-Trichloroethane	ND	0.14	ND	0.026	
127-18-4	Tetrachloroethene	ND	0.036	ND	0.0053	
100-41-4	Ethylbenzene	<b>0.61</b>	0.14	<b>0.14</b>	0.033	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.036	ND	0.0052	
91-20-3	Naphthalene	ND	0.14	ND	0.027	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 21-1F-031512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-003

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01566

Date Collected: 3/16/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/31/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.65      Final Pressure (psig): 3.73

Canister Dilution Factor: 1.83

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.046	ND	0.018	
107-06-2	1,2-Dichloroethane	1.1	0.046	0.26	0.011	
71-43-2	Benzene	0.92	0.14	0.29	0.043	
79-01-6	Trichloroethene	1.3	0.046	0.24	0.0085	
79-00-5	1,1,2-Trichloroethane	ND	0.18	ND	0.034	
127-18-4	Tetrachloroethene	0.20	0.046	0.029	0.0067	
100-41-4	Ethylbenzene	2.9	0.18	0.67	0.042	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.046	ND	0.0067	
91-20-3	Naphthalene	0.39	0.18	0.074	0.035	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 23-1F-031512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-004

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00988

Date Collected: 3/16/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/31/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.67      Final Pressure (psig): 3.62

Canister Dilution Factor: 1.83

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.046	ND	0.018	
107-06-2	1,2-Dichloroethane	<b>0.71</b>	0.046	<b>0.17</b>	0.011	
71-43-2	Benzene	<b>0.46</b>	0.14	<b>0.14</b>	0.043	
79-01-6	Trichloroethene	ND	0.046	ND	0.0085	
79-00-5	1,1,2-Trichloroethane	ND	0.18	ND	0.034	
127-18-4	Tetrachloroethene	<b>0.21</b>	0.046	<b>0.031</b>	0.0067	
100-41-4	Ethylbenzene	<b>0.43</b>	0.18	<b>0.098</b>	0.042	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.046	ND	0.0067	
91-20-3	Naphthalene	ND	0.18	ND	0.035	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 17-1F-031712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P1201121-005

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01790

Date Collected: 3/16/12  
Date Received: 3/22/12  
Date Analyzed: 3/31/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.83      Final Pressure (psig): 3.61

Canister Dilution Factor: 1.54

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.039	ND	0.015	
107-06-2	1,2-Dichloroethane	<b>1.0</b>	0.039	<b>0.25</b>	0.0095	
71-43-2	Benzene	<b>4.0</b>	0.12	<b>1.3</b>	0.036	
79-01-6	Trichloroethene	<b>0.088</b>	0.039	<b>0.016</b>	0.0072	
79-00-5	1,1,2-Trichloroethane	ND	0.15	ND	0.028	
127-18-4	Tetrachloroethene	<b>0.17</b>	0.039	<b>0.024</b>	0.0057	
100-41-4	Ethylbenzene	<b>2.4</b>	0.15	<b>0.54</b>	0.035	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.039	ND	0.0056	
91-20-3	Naphthalene	<b>0.74</b>	0.15	<b>0.14</b>	0.029	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 20-1F-031712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-006

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01305

Date Collected: 3/18/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/31/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.78      Final Pressure (psig): 3.86

Canister Dilution Factor: 1.44

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.036	ND	0.014	
107-06-2	1,2-Dichloroethane	<b>0.085</b>	0.036	<b>0.021</b>	0.0089	
71-43-2	Benzene	<b>1.6</b>	0.11	<b>0.51</b>	0.034	
79-01-6	Trichloroethene	ND	0.036	ND	0.0067	
79-00-5	1,1,2-Trichloroethane	ND	0.14	ND	0.026	
127-18-4	Tetrachloroethene	<b>1.1</b>	0.036	<b>0.16</b>	0.0053	
100-41-4	Ethylbenzene	<b>2.3</b>	0.14	<b>0.53</b>	0.033	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.036	ND	0.0052	
91-20-3	Naphthalene	<b>0.32</b>	0.14	<b>0.062</b>	0.027	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 28-1F-031912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-007

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00721

Date Collected: 3/20/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/31/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.99      Final Pressure (psig): 3.65

Canister Dilution Factor: 1.89

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.047	ND	0.018	
107-06-2	1,2-Dichloroethane	<b>0.053</b>	0.047	<b>0.013</b>	0.012	
71-43-2	Benzene	<b>0.47</b>	0.14	<b>0.15</b>	0.044	
79-01-6	Trichloroethene	ND	0.047	ND	0.0088	
79-00-5	1,1,2-Trichloroethane	ND	0.19	ND	0.035	
127-18-4	Tetrachloroethene	<b>2.6</b>	0.047	<b>0.38</b>	0.0070	
100-41-4	Ethylbenzene	<b>0.33</b>	0.19	<b>0.076</b>	0.044	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.047	ND	0.0069	
91-20-3	Naphthalene	<b>0.44</b>	0.19	<b>0.083</b>	0.036	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 22-1F-031912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-008

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01171

Date Collected: 3/20/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/31/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.96      Final Pressure (psig): 3.64

Canister Dilution Factor: 1.56

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.039	ND	0.015	
107-06-2	1,2-Dichloroethane	<b>0.054</b>	0.039	<b>0.013</b>	0.0096	
71-43-2	Benzene	<b>0.62</b>	0.12	<b>0.20</b>	0.037	
79-01-6	Trichloroethene	<b>2.1</b>	0.039	<b>0.39</b>	0.0073	
79-00-5	1,1,2-Trichloroethane	ND	0.16	ND	0.029	
127-18-4	Tetrachloroethene	<b>0.048</b>	0.039	<b>0.0071</b>	0.0058	
100-41-4	Ethylbenzene	<b>0.25</b>	0.16	<b>0.057</b>	0.036	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.039	ND	0.0057	
91-20-3	Naphthalene	<b>0.18</b>	0.16	<b>0.035</b>	0.030	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 26-1F-031912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-009

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01783

Date Collected: 3/20/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/31/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.72      Final Pressure (psig): 3.54

Canister Dilution Factor: 1.66

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.042	ND	0.016	
107-06-2	1,2-Dichloroethane	<b>0.13</b>	0.042	<b>0.032</b>	0.010	
71-43-2	Benzene	<b>0.51</b>	0.12	<b>0.16</b>	0.039	
79-01-6	Trichloroethene	ND	0.042	ND	0.0077	
79-00-5	1,1,2-Trichloroethane	ND	0.17	ND	0.030	
127-18-4	Tetrachloroethene	<b>0.096</b>	0.042	<b>0.014</b>	0.0061	
100-41-4	Ethylbenzene	<b>0.53</b>	0.17	<b>0.12</b>	0.038	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.042	ND	0.0060	
91-20-3	Naphthalene	<b>0.28</b>	0.17	<b>0.053</b>	0.032	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P120331-MB

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/31/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.025	ND	0.0098	
107-06-2	1,2-Dichloroethane	ND	0.025	ND	0.0062	
71-43-2	Benzene	ND	0.075	ND	0.023	
79-01-6	Trichloroethene	ND	0.025	ND	0.0047	
79-00-5	1,1,2-Trichloroethane	ND	0.10	ND	0.018	
127-18-4	Tetrachloroethene	ND	0.025	ND	0.0037	
100-41-4	Ethylbenzene	ND	0.10	ND	0.023	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.025	ND	0.0036	
91-20-3	Naphthalene	ND	0.10	ND	0.019	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister(s)  
 Test Notes:

Date(s) Collected: 3/15 - 3/20/12  
 Date(s) Received: 3/22/12  
 Date(s) Analyzed: 3/31/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		% Recovered	% Recovered	% Recovered		
Method Blank	P120331-MB	102	113	111	70-130	
Lab Control Sample	P120331-LCS	103	113	115	70-130	
19-1F-031412	P1201121-001	103	115	105	70-130	
19-1F-031412	P1201121-001DUP	100	113	106	70-130	
53-1F-031412	P1201121-002	102	112	104	70-130	
21-1F-031512	P1201121-003	101	116	106	70-130	
23-1F-031512	P1201121-004	101	112	110	70-130	
17-1F-031712	P1201121-005	99	111	108	70-130	
20-1F-031712	P1201121-006	101	111	107	70-130	
28-1F-031912	P1201121-007	102	112	112	70-130	
22-1F-031912	P1201121-008	98	114	111	70-130	
26-1F-031912	P1201121-009	99	112	114	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.

LABORATORY CONTROL SAMPLE SUMMARY

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**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
CAS Sample ID: P120331-LCS

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 3/31/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
75-01-4	Vinyl Chloride	3.88	<b>3.18</b>	<b>82</b>	56-127	
107-06-2	1,2-Dichloroethane	4.00	<b>3.15</b>	<b>79</b>	51-140	
71-43-2	Benzene	3.96	<b>3.54</b>	<b>89</b>	56-125	
79-01-6	Trichloroethene	3.88	<b>2.93</b>	<b>76</b>	51-127	
79-00-5	1,1,2-Trichloroethane	3.92	<b>2.92</b>	<b>74</b>	49-137	
127-18-4	Tetrachloroethene	3.68	<b>2.84</b>	<b>77</b>	58-134	
100-41-4	Ethylbenzene	3.96	<b>3.05</b>	<b>77</b>	56-126	
79-34-5	1,1,2,2-Tetrachloroethane	3.84	<b>2.99</b>	<b>78</b>	53-148	
91-20-3	Naphthalene	3.44	<b>3.24</b>	<b>94</b>	25-149	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY DUPLICATE SUMMARY RESULTS

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**Client:** Montana DEQ  
**Client Sample ID:** 19-1F-031412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201121  
 CAS Sample ID: P1201121-001DUP

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01817

Date Collected: 3/15/12  
 Date Received: 3/22/12  
 Date Analyzed: 3/31/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.81      Final Pressure (psig): 3.51

Canister Dilution Factor: 1.67

CAS #	Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
		µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
75-01-4	Vinyl Chloride	ND	ND	ND	ND	-	-	25	
107-06-2	1,2-Dichloroethane	0.122	0.0301	0.120	0.0297	0.121	2	25	
71-43-2	Benzene	0.705	0.221	0.718	0.225	0.7115	2	25	
79-01-6	Trichloroethene	ND	ND	ND	ND	-	-	25	
79-00-5	1,1,2-Trichloroethane	ND	ND	ND	ND	-	-	25	
127-18-4	Tetrachloroethene	0.0701	0.0103	0.0690	0.0102	0.06955	2	25	
100-41-4	Ethylbenzene	0.764	0.176	0.778	0.179	0.771	2	25	
79-34-5	1,1,2,2-Tetrachloroethane	ND	ND	ND	ND	-	-	25	
91-20-3	Naphthalene	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

Massachusetts APH

Hydrocarbon Ranges

ICAL Method: M801412A.M

ICAL Date: 1/14/2012

Instrument ID: MS08

	areas				masses							
	0.5	1	5	25	50	100	0.5	1	5	25	50	100
<b>C5-C8 Aliphatics</b>												
Isopentane	60478	112606	488243	2480105	4710733	9222808	0.520	1.04	5.20	26.0	52.0	104
n-Hexane	59953	110631	507126	2577865	4931102	9678199	0.490	0.98	4.90	24.5	49.0	98
Cyclohexane	63203	115230	548910	2874726	5524402	10890911	0.500	1.00	5.00	25.0	50.0	100
2,3-Dimethylpentane	68648	127222	585183	2998017	5742104	11319597	0.520	1.04	5.20	26.0	52.0	104
n-Heptane	66143	120286	558074	2884817	5509801	10873046	0.490	0.98	4.90	24.5	49.0	98
n-Octane	76169	137829	632891	3286017	6205035	12092514	0.505	1.01	5.05	25.3	50.5	101
<b>area sum:</b>	<b>394594</b>	<b>723804</b>	<b>3320427</b>	<b>17101547</b>	<b>32623177</b>	<b>64077075</b>	<b>3.025</b>	<b>6.050</b>	<b>30.25</b>	<b>151.30</b>	<b>302.50</b>	<b>605.00</b>
<b>C9-C12 Aliphatics</b>												
2,3-Dimethylheptane	78780	147724	682518	3501487	6691428	13148840	0.485	0.97	4.85	24.3	48.5	97
n-Nonane	77464	140680	662221	3417128	6555165	12940625	0.485	0.97	4.85	24.3	48.5	97
n-Decane	80593	147724	705259	3651028	7021424	13870530	0.490	0.98	4.90	24.5	49.0	98
3-Butylcyclohexane	88915	166782	776903	4026931	7749936	15415340	0.495	0.99	4.95	24.8	49.5	99
n-Undecane	80867	152557	719618	3772331	7263077	14385881	0.480	0.96	4.80	24.0	48.0	96
n-Dodecane	70045	143746	687255	3673524	7115012	14041583	0.500	1.00	5.00	25.0	50.0	100
<b>area sum:</b>	<b>476664</b>	<b>899213</b>	<b>4233774</b>	<b>22042429</b>	<b>42396042</b>	<b>83802799</b>	<b>2.935</b>	<b>5.870</b>	<b>29.35</b>	<b>146.90</b>	<b>293.50</b>	<b>587.00</b>
<b>C9-C10 Aromatics</b>												
Isopropylbenzene	10658	19169	87439	448287	849124	1695737	0.485	0.97	4.85	24.3	48.5	97
3-Ethyltoluene	11805	20564	97823	504510	977020	1973178	0.470	0.94	4.70	23.5	47.0	94
1,3,5-Trimethylbenzene	15759	29260	134073	700133	1353305	2726630	0.490	0.98	4.90	24.5	49.0	98
1,2,3-Trimethylbenzene	17195	31328	143473	757946	1493373	3063059	0.495	0.99	4.95	24.8	49.5	99
p-Isopropyltoluene	9490	18116	83401	440962	872799	1814314	0.460	0.92	4.60	23.0	46.0	92
<b>area sum:</b>	<b>64907</b>	<b>118437</b>	<b>546209</b>	<b>2851838</b>	<b>5545621</b>	<b>11272918</b>	<b>2.400</b>	<b>4.800</b>	<b>24.00</b>	<b>120.10</b>	<b>240.00</b>	<b>480.00</b>



Massachusetts APH  
Hydrocarbon Ranges

ICAL: M8011412A.M

ICAL Date: 1/14/2012

Instrument ID: MS08

areas

	0.5	1	5	25	50	100
<u>Internal Standards (TIC)</u>						
Bromochloromethane (IS1)	1312793	1307309	1293586	1289283	1284194	1283817
1,4-Difluorobenzene (IS2)	2547079	2541159	2512498	2475255	2422119	2352159
Chlorobenzene-d5 (IS3)	2929189	2913374	2880258	2895080	2878314	2908396

Internal Standards (EIC)

Bromochloromethane (IS1)	249696	247899	245985	245585	244327	243561
1,4-Difluorobenzene (IS2)	1153657	1149947	1133233	1136223	1127203	1138730
Chlorobenzene-d5 (IS3)	1007702	1004638	991212	992925	982812	986950

Surrogates (TIC)

1,2-Dichloroethane-d4	1162497	1154617	1105630	1126201	1103958	1095945
Toluene-d8	3508018	3490493	3455382	3466688	3437986	3457323
o-Bromofluorobenzene	2298220	2284797	2197239	2227396	2192159	2197999

C5-C8 Aliphatics

	0.5	1	5	25	50	100	RRF <sub>avg</sub>	%RSD
	2.8268	2.6009	2.4215	2.4870	2.3919	2.3252	2.509	7.24

C9-C12 Aliphatics

	0.5	1	5	25	50	100	RRF <sub>avg</sub>	%RSD
	4.0291	3.8120	3.6383	3.7780	3.6744	3.6163	3.758	4.09

C9-C10 Aromatics

	0.5	1	5	25	50	100	RRF <sub>avg</sub>	%RSD
	0.6709	0.6140	0.5740	0.5979	0.5878	0.5949	0.607	5.63

**Massachusetts APH**  
**Continuing Calibration Verification Check Sheet**

Data File Name: 03271202.D  
 Data File Path: J:\MS08\Data\2012\_03\27\  
 Operator: EM  
 Date Acquired: 3/27/12 3:36  
 Acq. Method File: TO15.M  
 Sample Name: 25ng TO-15/MAPH CCV STD  
 Misc Info: S25-03151201/S25-03161202  
 Instrument Name: MS08

Enter RRFs from current ICAL!

Internal Standards	<u>RT</u>	<u>Area</u>						
7) 1,4-Difluorobenzene (IS2)	15.89	1117304						
16) Chlorobenzene-d5 (IS3)	21.69	978107						
C5-C8 Aliphatics	<u>RT</u>	<u>Area</u>	<u>RRF</u>	<u>ng</u>	<u>% D</u>	<u>LCL</u>	<u>UCL</u>	<u>Pass/Fai</u>
3) Isopentane	8.20	2419558	2.482	149.7	-1.08	-30	30	Pass
4) n-Hexane	13.09	2528433						
9) Cyclohexane	15.81	2791802						
10) 2,3-Dimethylpentane	16.17	2948565	Spike	ICAL				
11) n-Heptane	17.35	2834960	Amt (ng)	RRF				
14) n-Octane	20.69	<u>3258272</u>	151.30	2.509				
		<b>16781590</b>						
C9-C12 Aliphatics	<u>RT</u>	<u>Area</u>	<u>RRF</u>	<u>ng</u>	<u>% D</u>	<u>LCL</u>	<u>UCL</u>	<u>Pass/Fai</u>
18) 2,3-Dimethylheptane	22.31	3441244	3.743	146.3	-0.39	-30	30	Pass
19) n-Nonane	23.30	3354311						
25) n-Decane	25.27	3574971						
28) Butylcyclohexane	25.99	3888285	Spike	ICAL				
29) n-Undecane	26.77	3691884	Amt (ng)	RRF				
30) n-Dodecane	28.00	<u>3563475</u>	146.90	3.758				
		<b>21514170</b>						
C9-C10 Aromatics	<u>RT</u>	<u>Area</u>	<u>RRF</u>	<u>ng</u>	<u>% D</u>	<u>LCL</u>	<u>UCL</u>	<u>Pass/Fai</u>
22) Isopropylbenzene	23.78	435871	0.579	114.5	-4.64	-30	30	Pass
23) 1-Methyl-3-ethylbenzene	24.52	481596						
24) 1,3,5-Trimethylbenzene	24.66	665971						
26) p-Isopropyltoluene	25.68	416913	Spike	ICAL				
27) 1,2,3-Trimethylbenzene	25.68	<u>719386</u>	Amt (ng)	RRF				
		<b>2719737</b>	120.1	0.607				

EM 3/27/12

Response Factor Report MS08

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Mon Dec 12 09:45:45 2011  
 Response Via : Initial Calibration

Calibration Files

0.1 =12091118.D 0.2 =12091119.D 0.5 =12091120.D 1.0 =12091121.D 5.0 =12091122.D 25 =12091123.D  
 50 =12091124.D 100 =12091125.D

Compound	0.1	0.2	0.5	1.0	5.0	25	50	100	Avg	%RSD
1) IR Bromochloromethane...										
2) T Propene	2.487	1.965	1.952	1.692	2.194	2.303	2.185	2.036	2.102	11.62
3) T Dichlorodifluo...	3.545	2.805	2.940	2.748	2.432	2.361	2.278	2.199	2.664	16.73
4) T Chloromethane	3.290	2.811	2.875	2.512	2.176	2.286	1.977	1.664	2.449	21.62
5) T 1,2-Dichloro-1...	1.901	1.661	1.575	1.458	1.359	1.333	1.200	1.190	1.460	16.68
6) T Vinyl Chloride	2.906	2.590	2.631	2.374	2.145	2.129	1.950	1.939	2.333	15.07
7) T 1,3-Butadiene	2.145	1.788	1.790	1.718	1.581	1.689	1.621	1.616	1.743	10.33
8) T Bromomethane	1.687	1.328	1.349	1.212	1.094	1.142	1.111	1.092	1.252	16.21
9) T Chloroethane	1.671	1.259	1.333	1.193	1.101	1.091	1.046	1.044	1.217	17.29
10) T Ethanol	2.040	1.345	1.186	1.085	1.022	1.096	1.076	1.064	1.239	27.34
11) T Acetonitrile	4.365	2.973	2.723	2.546	2.480	2.562	2.482	2.468	2.825	22.85
12) T Acrolein	1.328	0.931	0.907	0.861	0.772	0.814	0.789	0.779	0.897	20.48
13) T Acetone	1.645	1.316	1.249	1.129	0.994	1.015	0.981	0.964	1.161	20.26
14) T Trichlorofluor...	2.776	2.385	2.496	2.272	2.097	2.135	2.042	2.011	2.277	11.59
15) T 2-Propanol (Is...		4.463	4.536	4.168	2.920	2.763	2.577	2.672	3.443	26.09
16) T Acrylonitrile	2.701	1.907	1.878	1.783	1.745	1.813	1.765	1.738	1.916	16.85
17) T 1,1-Dichloroet...	1.458	1.418	1.452	1.263	1.206	1.223	1.172	1.152	1.293	9.95
18) T 2-Methyl-2-Pro...	4.464	3.710	3.939	3.616	3.558	3.584			3.812	9.14
19) T Methylene Chlo...			1.958	1.578	1.294	1.273	1.220	1.201	1.421	20.88
20) T 3-Chloro-1-pro...	2.347	2.260	2.345	2.220	2.150	2.296	2.230	2.191	2.255	3.15
21) T Trichlorotrifl...	1.348	1.244	1.299	1.193	1.089	1.098	1.062	1.035	1.171	10.01
22) T Carbon Disulfide		6.210	5.678	5.057	4.654	4.699	4.488	4.424	5.030	13.40
23) T trans-1,2-Dich...	2.436	2.237	2.206	2.142	1.951	1.953	1.883	1.835	2.080	10.00
24) T 1,1-Dichloroet...	2.896	2.757	2.747	2.500	2.307	2.379	2.288	2.236	2.514	10.07
25) T Methyl tert-Bu...	5.128	4.611	4.650	4.307	4.119	4.185	4.042	3.984	4.378	8.93
26) T Vinyl Acetate	0.430	0.368	0.423	0.399	0.393	0.410	0.402	0.398	0.403	4.70
27) T 2-Butanone (MEK)	1.265	1.003	1.038	0.972	0.927	0.964	0.928	0.794	0.986	13.56
28) T cis-1,2-Dichlo...	2.442	2.056	2.101	1.938	1.826	1.836	1.764	1.720	1.960	12.05
29) T Diisopropyl Ether	1.446	1.223	1.323	1.262	1.171	1.188	1.153	1.148	1.239	8.26
30) T Ethyl Acetate	0.611	0.554	0.573	0.545	0.517	0.535	0.520	0.516	0.546	6.02
31) T n-Hexane	3.541	3.006	3.181	2.805	2.582	2.633	2.545	2.543	2.855	12.70

Response Factor Report MS08

Method	Path	J:\MS08\Methods\	EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)									
Method	File	R8120911A.M	ISTD									
Title	Title											
32) T	Chloroform	2.679	2.408	2.409	2.204	1.983	2.054	1.974	1.937	2.206	12.18	
33) S	1,2-Dichloroet...	1.586	1.577	1.582	1.590	1.535	1.533	1.520	1.504	1.553	2.19	
34) T	Tetrahydrofura...	1.274	0.932	1.100	0.938	0.910	0.879	0.830	0.821	0.961	15.96	
35) T	Ethyl tert-But...	1.999	1.832	2.033	1.818	1.732	1.757	1.708	1.687	1.821	7.18	
36) T	1,2-Dichloroet...	2.099	1.992	1.976	1.806	1.683	1.700	1.599	1.531	1.798	11.39	
37) IR	1,4-Difluorobenzen...	ISTD										
38) T	1,1,1-Trichloror...	0.531	0.465	0.486	0.453	0.419	0.434	0.413	0.403	0.451	9.46	
39) T	Isopropyl Acetate	0.227	0.205	0.220	0.206	0.194	0.201	0.192	0.188	0.204	6.71	
40) T	1-Butanol	0.348	0.319	0.300	0.320	0.335	0.324	0.318	0.323	0.323	4.66	
41) T	Benzene	1.633	1.385	1.359	1.227	1.136	1.158	1.101	1.070	1.258	15.11	
42) T	Carbon Tetrach...	0.394	0.356	0.369	0.340	0.334	0.352	0.337	0.333	0.352	6.03	
43) T	Cyclohexane	0.666	0.576	0.588	0.539	0.492	0.500	0.479	0.469	0.539	12.59	
44) T	tert-Amyl Meth...	1.092	0.926	0.998	0.937	0.900	0.918	0.881	0.858	0.939	7.95	
45) T	1,2-Dichloropr...	0.397	0.361	0.368	0.324	0.312	0.316	0.300	0.293	0.334	11.12	
46) T	Bromodichlorom...	0.458	0.387	0.392	0.378	0.362	0.378	0.362	0.353	0.384	8.57	
47) T	Trichloroethene	0.466	0.375	0.379	0.344	0.323	0.328	0.316	0.309	0.355	14.62	
48) T	1,4-Dioxane	0.305	0.282	0.288	0.248	0.245	0.246	0.237	0.230	0.260	10.55	
49) T	2,2,4-Trimethy...	1.964	1.629	1.686	1.546	1.447	1.461	1.387	1.337	1.557	12.98	
50) T	Methyl Methacr...	0.156	0.139	0.150	0.135	0.136	0.142	0.136	0.135	0.141	5.53	
51) T	n-Heptane	0.427	0.373	0.385	0.364	0.341	0.344	0.332	0.324	0.361	9.35	
52) T	cis-1,3-Dichlo...	0.571	0.502	0.521	0.491	0.482	0.500	0.481	0.471	0.502	6.33	
53) T	4-Methyl-2-pen...	0.359	0.295	0.308	0.303	0.297	0.303	0.289	0.284	0.305	7.68	
54) T	trans-1,3-Dich...	0.476	0.413	0.443	0.421	0.429	0.453	0.437	0.427	0.437	4.59	
55) T	1,1,2-Trichlor...	0.340	0.301	0.310	0.284	0.276	0.285	0.271	0.265	0.292	8.44	
56) IR	Chlorobenzene-d5 (...)	ISTD										
57) S	Toluene-d8 (SS2)	2.305	2.306	2.305	2.309	2.306	2.300	2.310	2.304	2.306	0.14	
58) T	Toluene	3.623	2.920	3.025	2.738	2.571	2.574	2.473	2.421	2.793	14.19	
59) T	2-Hexanone	2.049	1.429	1.515	1.403	1.395	1.422	1.367	1.337	1.490	15.57	
60) T	Dibromochlorom...	0.696	0.618	0.636	0.636	0.612	0.648	0.629	0.625	0.638	4.14	
61) T	1,2-Dibromoethane	0.831	0.730	0.718	0.658	0.641	0.662	0.638	0.628	0.688	9.97	
62) T	n-Butyl Acetate	2.338	1.850	1.857	1.721	1.704	1.765	1.737	1.737	1.839	11.40	
63) T	n-Octane	0.785	0.652	0.714	0.643	0.602	0.611	0.580	0.572	0.645	11.28	
64) T	Tetrachloroethene	1.090	0.904	0.904	0.826	0.767	0.779	0.756	0.755	0.848	13.63	
65) T	Chlorobenzene	2.375	1.868	1.886	1.744	1.610	1.637	1.574	1.548	1.780	15.30	
66) T	Ethylbenzene	3.846	3.171	3.188	2.956	2.803	2.848	2.743	2.680	3.030	12.51	
67) T	m- & p-Xylenes	3.081	2.499	2.611	2.392	2.216	2.273	2.170	2.112	2.419	13.08	
68) T	Bromoform	0.622	0.494	0.556	0.541	0.551	0.598	0.589	0.594	0.568	7.19	
69) T	Styrene	2.478	1.895	1.937	1.800	1.735	1.803	1.726	1.675	1.881	13.64	
70) T	o-Xylene	3.279	2.708	2.726	2.479	2.331	2.373	2.283	2.224	2.550	13.65	

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

71)	T	n-Nonane	2.005	1.753	1.847	1.675	1.581	1.597	1.518	1.470	1.681	10.69
72)	T	1,1,2,2-Tetrac...	1.348	1.164	1.262	1.150	1.083	1.121	1.085	1.066	1.160	8.48
73)	S	Bromofluoroben...	0.774	0.771	0.779	0.781	0.775	0.778	0.782	0.780	0.777	0.50
74)	T	Cumene	4.522	3.495	3.524	3.169	2.962	3.002	2.884	2.831	3.299	16.98
75)	T	alpha-Pinene	1.782	1.636	1.734	1.601	1.537	1.578	1.518	1.498	1.611	6.35
76)	T	n-Propylbenzene	5.441	4.229	4.306	3.924	3.710	3.754	3.553	3.394	4.039	16.01
77)	T	3-Ethyltoluene	4.091	3.373	3.461	3.038	2.964	2.981	2.893	2.845	3.206	13.14
78)	T	4-Ethyltoluene	4.003	3.202	3.301	3.151	2.837	2.983	2.810	2.776	3.133	12.84
79)	T	1,3,5-Trimethy...	3.324	2.662	2.723	2.537	2.378	2.426	2.344	2.403	2.600	12.43
80)	T	alpha-Methylst...	1.742	1.386	1.449	1.371	1.390	1.402	1.354	1.399	1.436	8.80
81)	T	2-Ethyltoluene	4.143	3.362	3.430	3.169	2.975	3.048	2.919	2.985	3.254	12.43
82)	T	1,2,4-Trimethy...	3.356	2.663	2.791	2.551	2.433	2.513	2.454	2.548	2.664	11.37
83)	T	n-Decane	1.969	1.702	1.844	1.705	1.582	1.626	1.554	1.609	1.699	8.39
84)	T	Benzyl Chloride	2.373	1.648	1.772	1.688	1.799	2.058	2.054	2.190	1.948	13.35
85)	T	1,3-Dichlorobe...	2.498	1.658	1.649	1.455	1.377	1.409	1.374	1.436	1.607	23.49
86)	T	1,4-Dichlorobe...	2.655	1.785	1.618	1.492	1.380	1.428	1.388	1.449	1.649	25.99
87)	T	sec-Butylbenzene	4.527	3.773	3.816	3.471	3.264	3.328	3.214	3.176	3.571	12.80
88)	T	4-Isopropyltol...	4.426	3.674	3.683	3.416	3.254	3.356	3.253	3.366	3.554	10.98
889)	T	1,2,3-Trimethy...	3.266	2.650	2.852	2.588	2.471	2.573	2.506	2.619	2.691	9.63
890)	T	1,2-Dichlorobe...	2.173	1.589	1.545	1.435	1.332	1.387	1.370	1.415	1.531	17.89
891)	T	d-Limonene	1.124	0.998	1.106	1.069	1.064	1.072	1.040	1.065	1.067	3.58
92)	T	1,2-Dibromo-3-...	0.563	0.488	0.501	0.486	0.494	0.525	0.532	0.542	0.517	5.50
93)	T	n-Undecane	1.922	1.637	1.755	1.671	1.608	1.630	1.608	1.592	1.678	6.63
94)	T	1,2,4-Trichlor...	1.335	1.165	1.062	1.017	1.017	1.071	1.086	1.118	1.122	9.32
95)	T	Naphthalene	4.848	3.923	3.638	3.497	3.721	3.755	3.783	3.881	11.50	11.50
96)	T	n-Dodecane	1.866	1.599	1.669	1.586	1.545	1.575	1.553	1.509	1.613	6.99
97)	T	Hexachlorobuta...	0.841	0.749	0.730	0.673	0.635	0.664	0.663	0.676	0.704	9.50
98)	T	Cyclohexanone	1.172	0.979	1.021	0.969	0.952	0.962	0.931	0.903	0.986	8.37
99)	T	tert-Butylbenzene	3.486	2.799	2.880	2.658	2.449	2.533	2.453	2.528	2.723	12.71
100)	T	n-Butylbenzene	3.596	2.904	2.958	2.670	2.572	2.625	2.613	2.593	2.817	12.33

(#) = Out of Range

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\27\  
 Data File : 03271201.D  
 Acq On : 27 Mar 2012 2:54  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 27 08:42:57 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	IR Bromochloromethane (IS1)	1.000	1.000	0.0	84	-0.01
2	T Propene	2.102	2.112	-0.5	77	0.00
3	T Dichlorodifluoromethane (CF	2.664	2.197	17.5	78	0.01
4	T Chloromethane	2.449	2.077	15.2	76	0.00
5	T 1,2-Dichloro-1,1,2,2-tetra	1.460	1.162	20.4	73	0.00
6	T Vinyl Chloride	2.333	1.938	16.9	77	0.00
7	T 1,3-Butadiene	1.743	1.605	7.9	80	0.00
8	T Bromomethane	1.252	1.046	16.5	77	0.00
9	T Chloroethane	1.217	1.028	15.5	79	0.00
10	T Ethanol	1.239	1.128	9.0	86	-0.07
11	T Acetonitrile	2.825	2.506	11.3	82	-0.05
12	T Acrolein	0.897	0.786	12.4	81	-0.02
13	T Acetone	1.161	0.998	14.0	83	-0.04
14	T Trichlorofluoromethane	2.277	2.021	11.2	80	0.00
15	T 2-Propanol (Isopropanol)	3.443	2.787	19.1	85	-0.05
16	T Acrylonitrile	1.916	1.761	8.1	82	-0.03
17	T 1,1-Dichloroethene	1.293	1.147	11.3	79	0.00
18	T 2-Methyl-2-Propanol (tert-B	3.812	3.111	18.4	73	-0.03
19	T Methylene Chloride	1.421	1.204	15.3	79	0.00
20	T 3-Chloro-1-propene (Allyl C	2.255	2.255	0.0	83	-0.01
21	T Trichlorotrifluoroethane	1.171	1.027	12.3	79	0.00
22	T Carbon Disulfide	5.030	4.445	11.6	80	0.00
23	T trans-1,2-Dichloroethene	2.080	1.902	8.6	82	-0.01
24	T 1,1-Dichloroethane	2.514	2.311	8.1	82	-0.01
25	T Methyl tert-Butyl Ether	4.378	4.075	6.9	82	0.00
26	T Vinyl Acetate	0.403	0.403	0.0	83	-0.03
27	T 2-Butanone (MEK)	0.986	0.939	4.8	82	-0.02
28	T cis-1,2-Dichloroethene	1.960	1.778	9.3	81	-0.01
29	T Diisopropyl Ether	1.239	1.147	7.4	81	0.00
30	T Ethyl Acetate	0.546	0.523	4.2	82	-0.02
31	T n-Hexane	2.855	2.532	11.3	81	0.00
32	T Chloroform	2.206	1.979	10.3	81	-0.03
33	S 1,2-Dichloroethane-d4 (SS1)	1.553	1.610	-3.7	88	-0.01
34	T Tetrahydrofuran (THF)	0.961	0.861	10.4	82	0.00
35	T Ethyl tert-Butyl Ether	1.821	1.702	6.5	81	0.00
36	T 1,2-Dichloroethane	1.798	1.659	7.7	82	-0.01
37	IR 1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	88	0.00
38	T 1,1,1-Trichloroethane	0.451	0.401	11.1	81	0.00

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\27\  
 Data File : 03271201.D  
 Acq On : 27 Mar 2012 2:54  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 27 08:42:57 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
39 T	Isopropyl Acetate	0.204	0.188	7.8	82	-0.02
40 T	1-Butanol	0.323	0.344	-6.5	91	-0.05
41 T	Benzene	1.258	1.074	14.6	82	-0.01
42 T	Carbon Tetrachloride	0.352	0.321	8.8	80	0.00
43 T	Cyclohexane	0.539	0.468	13.2	82	-0.01
44 T	tert-Amyl Methyl Ether	0.939	0.854	9.1	82	-0.01
45 T	1,2-Dichloropropane	0.334	0.297	11.1	83	-0.01
46 T	Bromodichloromethane	0.384	0.351	8.6	82	0.00
47 T	Trichloroethene	0.355	0.296	16.6	79	-0.01
48 T	1,4-Dioxane	0.260	0.235	9.6	84	-0.02
49 T	2,2,4-Trimethylpentane (Iso	1.557	1.394	10.5	84	0.00
50 T	Methyl Methacrylate	0.141	0.133	5.7	83	-0.02
51 T	n-Heptane	0.361	0.329	8.9	84	-0.01
52 T	cis-1,3-Dichloropropene	0.502	0.468	6.8	82	0.00
53 T	4-Methyl-2-pentanone	0.305	0.293	3.9	85	-0.02
54 T	trans-1,3-Dichloropropene	0.437	0.423	3.2	82	-0.01
55 T	1,1,2-Trichloroethane	0.292	0.262	10.3	81	-0.01
56 IR	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	89	0.00
57 S	Toluene-d8 (SS2)	2.306	2.282	1.0	89	0.00
58 T	Toluene	2.793	2.378	14.9	83	0.00
59 T	2-Hexanone	1.490	1.395	6.4	88	-0.02
60 T	Dibromochloromethane	0.638	0.582	8.8	80	0.00
61 T	1,2-Dibromoethane	0.688	0.597	13.2	81	-0.01
62 T	n-Butyl Acetate	1.839	1.730	5.9	88	-0.01
63 T	n-Octane	0.645	0.606	6.0	89	0.00
64 T	Tetrachloroethene	0.848	0.684	19.3	79	0.00
65 T	Chlorobenzene	1.780	1.486	16.5	81	0.00
66 T	Ethylbenzene	3.030	2.612	13.8	82	0.00
67 T	m- & p-Xylenes	2.419	2.142	11.5	84	-0.02
68 T	Bromoform	0.568	0.530	6.7	79	-0.01
69 T	Styrene	1.881	1.704	9.4	85	-0.01
70 T	o-Xylene	2.550	2.230	12.5	84	-0.01
71 T	n-Nonane	1.681	1.521	9.5	85	-0.01
72 T	1,1,2,2-Tetrachloroethane	1.160	1.023	11.8	82	-0.01
73 S	Bromofluorobenzene (SS3)	0.777	0.781	-0.5	90	0.00
74 T	Cumene	3.299	2.898	12.2	86	0.00
75 T	alpha-Pinene	1.611	1.456	9.6	83	0.00
76 T	n-Propylbenzene	4.039	3.497	13.4	83	0.00

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\27\  
 Data File : 03271201.D  
 Acq On : 27 Mar 2012 2:54  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 27 08:42:57 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
77 T	3-Ethyltoluene	3.206	2.733	14.8	82	-0.01
78 T	4-Ethyltoluene	3.133	2.715	13.3	81	-0.01
79 T	1,3,5-Trimethylbenzene	2.600	2.239	13.9	83	-0.01
80 T	alpha-Methylstyrene	1.436	1.311	8.7	84	-0.01
81 T	2-Ethyltoluene	3.254	2.785	14.4	82	-0.01
82 T	1,2,4-Trimethylbenzene	2.664	2.315	13.1	82	-0.01
83 T	n-Decane	1.699	1.534	9.7	84	-0.01
84 T	Benzyl Chloride	1.948	2.035	-4.5	88	-0.02
85 T	1,3-Dichlorobenzene	1.607	1.258	21.7	80	-0.01
86 T	1,4-Dichlorobenzene	1.649	1.263	23.4	79	-0.01
87 T	sec-Butylbenzene	3.571	3.086	13.6	83	-0.01
88 T	4-Isopropyltoluene (p-Cymen	3.554	3.091	13.0	82	-0.01
89 T	1,2,3-Trimethylbenzene	2.691	2.365	12.1	82	-0.01
90 T	1,2-Dichlorobenzene	1.531	1.235	19.3	80	-0.01
91 T	d-Limonene	1.067	1.028	3.7	86	-0.01
92 T	1,2-Dibromo-3-Chloropropane	0.517	0.472	8.7	80	-0.01
93 T	n-Undecane	1.678	1.542	8.1	85	0.00
94 T	1,2,4-Trichlorobenzene	1.122	0.935	16.7	78	-0.01
95 T	Naphthalene	3.881	3.346	13.8	80	-0.01
96 T	n-Dodecane	1.613	1.501	6.9	85	0.00
97 T	Hexachlorobutadiene	0.704	0.575	18.3	77	0.00
98 T	Cyclohexanone	0.986	0.980	0.6	91	-0.02
99 T	tert-Butylbenzene	2.723	2.344	13.9	83	-0.01
100 T	n-Butylbenzene	2.817	2.423	14.0	83	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*EM* 3/27/12



Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\28\  
 Data File : 03281201.D  
 Acq On : 28 Mar 2012 3:15  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 28 08:49:11 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	IR Bromochloromethane (IS1)	1.000	1.000	0.0	82	-0.01
2	T Propene	2.102	2.115	-0.6	75	0.00
3	T Dichlorodifluoromethane (CF	2.664	2.225	16.5	77	0.00
4	T Chloromethane	2.449	2.065	15.7	74	0.00
5	T 1,2-Dichloro-1,1,2,2-tetra	1.460	1.189	18.6	73	0.00
6	T Vinyl Chloride	2.333	1.960	16.0	76	0.00
7	T 1,3-Butadiene	1.743	1.627	6.7	79	0.00
8	T Bromomethane	1.252	1.038	17.1	75	0.00
9	T Chloroethane	1.217	1.044	14.2	79	0.00
10	T Ethanol	1.239	1.105	10.8	83	-0.07
11	T Acetonitrile	2.825	2.467	12.7	79	-0.05
12	T Acrolein	0.897	0.776	13.5	78	-0.02
13	T Acetone	1.161	0.986	15.1	80	-0.05
14	T Trichlorofluoromethane	2.277	2.004	12.0	77	0.00
15	T 2-Propanol (Isopropanol)	3.443	2.762	19.8	82	-0.05
16	T Acrylonitrile	1.916	1.730	9.7	78	-0.03
17	T 1,1-Dichloroethene	1.293	1.150	11.1	77	0.00
18	T 2-Methyl-2-Propanol (tert-B	3.812	2.960	22.4	68	-0.03
19	T Methylene Chloride	1.421	1.191	16.2	77	-0.01
20	T 3-Chloro-1-propene (Allyl C	2.255	2.222	1.5	80	-0.01
21	T Trichlorotrifluoroethane	1.171	1.026	12.4	77	0.00
22	T Carbon Disulfide	5.030	4.409	12.3	77	0.00
23	T trans-1,2-Dichloroethene	2.080	1.867	10.2	79	-0.01
24	T 1,1-Dichloroethane	2.514	2.273	9.6	79	-0.02
25	T Methyl tert-Butyl Ether	4.378	4.045	7.6	79	0.00
26	T Vinyl Acetate	0.403	0.398	1.2	80	-0.03
27	T 2-Butanone (MEK)	0.986	0.924	6.3	79	-0.02
28	T cis-1,2-Dichloroethene	1.960	1.740	11.2	78	-0.01
29	T Diisopropyl Ether	1.239	1.135	8.4	78	0.00
30	T Ethyl Acetate	0.546	0.515	5.7	79	-0.03
31	T n-Hexane	2.855	2.517	11.8	79	0.00
32	T Chloroform	2.206	1.934	12.3	77	-0.03
33	S 1,2-Dichloroethane-d4 (SS1)	1.553	1.583	-1.9	85	-0.01
34	T Tetrahydrofuran (THF)	0.961	0.842	12.4	79	0.00
35	T Ethyl tert-Butyl Ether	1.821	1.686	7.4	79	-0.01
36	T 1,2-Dichloroethane	1.798	1.603	10.8	77	-0.01
37	IR 1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	86	0.00
38	T 1,1,1-Trichloroethane	0.451	0.395	12.4	78	0.00

*Em* 3/28/12

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\28\  
 Data File : 03281201.D  
 Acq On : 28 Mar 2012 3:15  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 28 08:49:11 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
39 T	Isopropyl Acetate	0.204	0.186	8.8	79	-0.02
40 T	1-Butanol	0.323	0.338	-4.6	86	-0.05
41 T	Benzene	1.258	1.060	15.7	78	-0.01
42 T	Carbon Tetrachloride	0.352	0.315	10.5	77	-0.01
43 T	Cyclohexane	0.539	0.463	14.1	79	-0.01
44 T	tert-Amyl Methyl Ether	0.939	0.839	10.6	78	-0.01
45 T	1,2-Dichloropropane	0.334	0.287	14.1	78	-0.01
46 T	Bromodichloromethane	0.384	0.346	9.9	78	0.00
47 T	Trichloroethene	0.355	0.293	17.5	77	-0.01
48 T	1,4-Dioxane	0.260	0.230	11.5	80	-0.02
49 T	2,2,4-Trimethylpentane (Iso	1.557	1.363	12.5	80	-0.01
50 T	Methyl Methacrylate	0.141	0.131	7.1	79	-0.02
51 T	n-Heptane	0.361	0.320	11.4	80	-0.01
52 T	cis-1,3-Dichloropropene	0.502	0.459	8.6	79	0.00
53 T	4-Methyl-2-pentanone	0.305	0.288	5.6	82	-0.02
54 T	trans-1,3-Dichloropropene	0.437	0.412	5.7	78	0.00
55 T	1,1,2-Trichloroethane	0.292	0.255	12.7	77	-0.01
56 IR	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	86	0.00
57 S	Toluene-d8 (SS2)	2.306	2.271	1.5	85	0.00
58 T	Toluene	2.793	2.328	16.6	78	0.00
59 T	2-Hexanone	1.490	1.336	10.3	81	-0.02
60 T	Dibromochloromethane	0.638	0.576	9.7	77	0.00
61 T	1,2-Dibromoethane	0.688	0.587	14.7	77	-0.01
62 T	n-Butyl Acetate	1.839	1.646	10.5	80	-0.01
63 T	n-Octane	0.645	0.591	8.4	83	0.00
64 T	Tetrachloroethene	0.848	0.689	18.8	76	0.00
65 T	Chlorobenzene	1.780	1.472	17.3	78	0.00
66 T	Ethylbenzene	3.030	2.550	15.8	77	0.00
67 T	m- & p-Xylenes	2.419	2.094	13.4	79	-0.02
68 T	Bromoform	0.568	0.528	7.0	76	-0.01
69 T	Styrene	1.881	1.682	10.6	80	-0.01
70 T	o-Xylene	2.550	2.179	14.5	79	-0.01
71 T	n-Nonane	1.681	1.467	12.7	79	-0.01
72 T	1,1,2,2-Tetrachloroethane	1.160	1.004	13.4	77	-0.01
73 S	Bromofluorobenzene (SS3)	0.777	0.795	-2.3	88	0.00
74 T	Cumene	3.299	2.855	13.5	82	0.00
75 T	alpha-Pinene	1.611	1.440	10.6	79	0.00
76 T	n-Propylbenzene	4.039	3.424	15.2	79	0.00

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_03\28\  
 Data File : 03281201.D  
 Acq On : 28 Mar 2012 3:15  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03151201/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 28 08:49:11 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
77 T	3-Ethyltoluene	3.206	2.700	15.8	78	-0.01
78 T	4-Ethyltoluene	3.133	2.667	14.9	77	-0.01
79 T	1,3,5-Trimethylbenzene	2.600	2.215	14.8	79	-0.01
80 T	alpha-Methylstyrene	1.436	1.297	9.7	80	-0.01
81 T	2-Ethyltoluene	3.254	2.747	15.6	78	-0.01
82 T	1,2,4-Trimethylbenzene	2.664	2.270	14.8	78	-0.01
83 T	n-Decane	1.699	1.497	11.9	79	-0.01
84 T	Benzyl Chloride	1.948	1.984	-1.8	83	-0.02
85 T	1,3-Dichlorobenzene	1.607	1.248	22.3	76	-0.01
86 T	1,4-Dichlorobenzene	1.649	1.258	23.7	76	-0.02
87 T	sec-Butylbenzene	3.571	3.048	14.6	79	-0.01
88 T	4-Isopropyltoluene (p-Cymen	3.554	3.043	14.4	78	-0.01
89 T	1,2,3-Trimethylbenzene	2.691	2.322	13.7	78	-0.01
90 T	1,2-Dichlorobenzene	1.531	1.222	20.2	76	-0.01
91 T	d-Limonene	1.067	1.006	5.7	81	-0.01
92 T	1,2-Dibromo-3-Chloropropane	0.517	0.462	10.6	76	-0.01
93 T	n-Undecane	1.678	1.491	11.1	79	0.00
94 T	1,2,4-Trichlorobenzene	1.122	0.937	16.5	75	-0.01
95 T	Naphthalene	3.881	3.335	14.1	77	-0.01
96 T	n-Dodecane	1.613	1.472	8.7	81	0.00
97 T	Hexachlorobutadiene	0.704	0.582	17.3	76	0.00
98 T	Cyclohexanone	0.986	0.944	4.3	85	-0.02
99 T	tert-Butylbenzene	2.723	2.327	14.5	79	-0.01
100 T	n-Butylbenzene	2.817	2.362	16.2	78	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*EM 3/28/12*

Response Factor Report MS07

Method : J:\Ms07\METHODS\X7021712B.M (RTE Integrator)  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
 1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	Avg	%RSD
1) I Bromochloromethan	3.784	3.527	3.467	3.822	3.279	2.986	2.875	3.155	2.920	3.313	10.83
2) T Dichlorodifluorom		1.470	0.956	1.190	0.904	0.778	0.569	0.897	0.827	0.949	28.81
3) T Chloromethane	2.802	2.749	2.588	3.044	2.597	2.406	2.314	2.588	2.446	2.615	8.57
4) T Vinyl Chloride	1.894	1.496	1.211	1.613	1.335	1.208	1.105	1.433	1.367	1.407	17.11
5) T Bromomethane	1.511	1.416	1.357	1.517	1.295	1.237	1.195	1.312	1.224	1.340	8.92
6) T Chloroethane				1.548	1.214	1.122	1.080	1.193	1.090	1.208	14.52
7) T Acetone	3.051	2.887	2.831	3.145	2.692	2.565	2.447	2.691	2.536	2.761	8.59
8) T Trichlorofluorome	1.591	1.440	1.366	1.547	1.322	1.229	1.214	1.385	1.295	1.376	9.52
9) T 1,1-Dichloroethen	2.134	1.795	1.866	1.866	1.536	1.426	1.404	1.494	1.411	1.633	16.43
10) T Methylene Chlorid	1.408	1.325	1.295	1.376	1.169	1.081	1.062	1.170	1.097	1.220	10.87
11) T Trichlorotrifluor	1.972	1.710	1.589	1.762	1.505	1.406	1.393	1.512	1.439	1.588	12.18
12) T trans-1,2-Dichlor	3.265	3.020	3.016	3.383	2.925	2.756	2.681	2.937	2.841	2.980	7.60
13) T 1,1-Dichloroethan	5.168	4.740	4.762	5.045	4.335	4.137	4.166	4.690	4.471	4.613	7.90
14) T Methyl tert-Butyl	1.885	1.626	1.630	1.776	1.524	1.434	1.388	1.575	1.482	1.591	10.09
15) T cis-1,2-Dichloroe				3.041	2.586	2.390	2.280	2.524	2.374	2.636	14.03
16) T Chloroform	2.000	2.002	2.032	2.018	1.989	1.973	1.960	1.938	1.932	1.983	1.75
17) S 1,2-Dichloroethan	2.582	2.460	2.375	2.604	2.222	2.090	1.988	2.222	2.094	2.293	9.73
18) T 1,2-Dichloroethan	2.587	2.431	2.340	2.573	2.191	2.059	2.007	2.278	2.150	2.291	9.18
19) T 1,1,1-Trichloroet				7.536	6.055	5.665	5.395	5.903	5.468	6.222	14.88
20) T Benzene	2.280	1.842	1.769	1.950	1.676	1.579	1.549	1.771	1.692	1.790	12.41
21) T Carbon Tetrachlor											
22) I 1,4-Difluorobenze	0.417	0.397	0.373	0.409	0.354	0.331	0.319	0.358	0.336	0.366	9.68
23) T 1,2-Dichloropropa	0.491	0.464	0.436	0.489	0.420	0.398	0.394	0.447	0.416	0.439	8.26
24) T Bromodichlorometh	0.418	0.379	0.349	0.390	0.324	0.305	0.287	0.326	0.307	0.343	12.96
25) T Trichloroethene	0.324	0.290	0.270	0.278	0.242	0.231	0.227	0.258	0.243	0.263	11.99
26) T 1,4-Dioxane	0.567	0.523	0.508	0.540	0.499	0.484	0.488	0.559	0.522	0.521	5.67
27) T cis-1,3-Dichlorop	0.457	0.434	0.431	0.444	0.419	0.409	0.417	0.505	0.482	0.444	7.22
28) T trans-1,3-Dichlor											

(#) = Out of Range ### Number of calibration levels exceeded format ###  
 X7021712B.M Sat Mar 31 06:25:22 2012

Response Factor Report MS07

Method : J:\Ms07\METHODS\X7021712B.M (RTE Integrator)  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
 1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	Avg	%RSD
29) T 1,1,2-Trichloroet	0.315	0.296	0.286	0.305	0.263	0.248	0.242	0.274	0.257	0.276	9.32
30) S Toluene-d8 (SS2)	1.053	1.071	1.054	1.047	1.051	1.052	1.071	1.060	1.049	1.056	0.84
31) T Toluene	1.782	1.541	1.420	1.485	1.269	1.198	1.177	1.307	1.186	1.374	14.76
32) T 1,2-Dibromoethane	0.379	0.362	0.345	0.362	0.323	0.307	0.304	0.351	0.331	0.340	7.65
33) T Tetrachloroethene	0.423	0.394	0.371	0.400	0.341	0.320	0.311	0.347	0.326	0.359	11.03
-----ISTD-----											
34) I Chlorobenzene-d5											
35) T Chlorobenzene	4.581	4.172	4.017	4.308	3.688	3.469	3.311	3.705	3.345	3.844	11.68
36) T Ethylbenzene	7.604	7.082	6.707	7.235	6.297	6.070	5.915	6.682	5.878	6.608	9.30
37) T m,p-Xylene	5.861	5.347	5.132	5.536	4.930	4.777	4.719	5.268	4.558	5.125	8.24
38) T o-Xylene	6.505	5.826	5.526	5.889	5.188	4.995	4.950	5.594	4.932	5.490	9.67
39) T 1,1,2,2-Tetrachlo	2.655	2.674	2.710	2.646	2.437	2.371	2.298	2.761	2.498	2.561	6.40
40) S Bromofluorobenzene	1.712	1.701	1.695	1.697	1.683	1.680	1.670	1.642	1.593	1.675	2.20
41) T 1,3-Dichlorobenze	3.304	3.261	3.222	3.204	2.835	2.719	2.608	3.138	2.846	3.015	8.71
42) T 1,4-Dichlorobenze	3.470	3.303	3.208	3.208	2.803	2.671	2.579	3.174	2.891	3.034	10.13
43) T 1,2-Dichlorobenze	3.271	3.123	3.112	3.026	2.662	2.536	2.418	3.053	2.772	2.886	10.31
44) T 1,2,4-Trichlorobe	2.184	2.118	2.169	1.753	1.641	1.608	1.575	2.139	1.978	1.907	13.66
45) T Naphthalene	6.857	6.355	6.704	5.143	5.107	5.172	5.297	8.126	7.328	6.232	17.82
46) T Hexachlorobutadie	1.486	1.423	1.403	1.217	1.055	1.013	0.975	1.227	1.145	1.216	15.40

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_03\31\  
 Data File : 03311203.D  
 Acq On : 31 Mar 2012 4:54 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00185  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Mar 31 06:24:54 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane (IS1)	1.000	1.000	0.0	88	0.00
2 T	Dichlorodifluoromethane (CF	3.313	2.947	11.0	79	0.00
3 T	Chloromethane	0.949	0.874	7.9	85	0.00
4 T	Vinyl Chloride	2.615	2.401	8.2	81	0.00
5 T	Bromomethane	1.407	1.368	2.8	90	0.00
6 T	Chloroethane	1.340	1.238	7.6	84	0.00
7 T	Acetone	1.208	1.227	-1.6	89	0.00
8 T	Trichlorofluoromethane	2.761	2.452	11.2	80	0.00
9 T	1,1-Dichloroethene	1.376	1.319	4.1	88	0.00
10 T	Methylene Chloride	1.633	1.465	10.3	84	0.00
11 T	Trichlorotrifluoroethane	1.220	1.189	2.5	90	0.00
12 T	trans-1,2-Dichloroethene	1.588	1.446	8.9	85	0.00
13 T	1,1-Dichloroethane	2.980	2.772	7.0	83	0.00
14 T	Methyl tert-Butyl Ether	4.613	4.200	9.0	85	0.00
15 T	cis-1,2-Dichloroethene	1.591	1.507	5.3	87	0.00
16 T	Chloroform	2.636	2.615	0.8	89	0.00
17 S	1,2-Dichloroethane-d4 (SS1)	1.983	1.968	0.8	87	0.00
18 T	1,2-Dichloroethane	2.293	1.996	13.0	79	0.00
19 T	1,1,1-Trichloroethane	2.291	2.088	8.9	84	0.00
20 T	Benzene	6.222	6.142	1.3	89	0.00
21 T	Carbon Tetrachloride	1.790	1.570	12.3	82	0.00
22 I	1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	97	0.00
23 T	1,2-Dichloropropane	0.366	0.311	15.0	85	0.00
24 T	Bromodichloromethane	0.439	0.367	16.4	85	0.00
25 T	Trichloroethene	0.343	0.289	15.7	87	0.00
26 T	1,4-Dioxane	0.263	0.228	13.3	91	0.00
27 T	cis-1,3-Dichloropropene	0.521	0.456	12.5	89	0.00
28 T	trans-1,3-Dichloropropene	0.444	0.393	11.5	91	0.00
29 T	1,1,2-Trichloroethane	0.276	0.238	13.8	88	0.00
30 S	Toluene-d8 (SS2)	1.056	1.189	-12.6	110	0.00
31 T	Toluene	1.374	1.195	13.0	91	0.00
32 T	1,2-Dibromoethane	0.340	0.302	11.2	90	0.00
33 T	Tetrachloroethene	0.359	0.320	10.9	91	0.00
34 I	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	94	0.00
35 T	Chlorobenzene	3.844	3.609	6.1	92	0.00
36 T	Ethylbenzene	6.608	6.023	8.9	90	0.00
37 T	m,p-Xylene	5.125	4.811	6.1	92	0.00

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_03\31\  
 Data File : 03311203.D  
 Acq On : 31 Mar 2012 4:54 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00185  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Mar 31 06:24:54 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
38 T	o-Xylene	5.490	5.201	5.3	94	0.00
39 T	1,1,2,2-Tetrachloroethane	2.561	2.404	6.1	93	0.00
40 S	Bromofluorobenzene (SS3)	1.675	1.862	-11.2	104	0.00
41 T	1,3-Dichlorobenzene	3.015	3.021	-0.2	100	0.00
42 T	1,4-Dichlorobenzene	3.034	2.972	2.0	100	0.00
43 T	1,2-Dichlorobenzene	2.886	2.904	-0.6	103	0.00
44 T	1,2,4-Trichlorobenzene	1.907	2.056	-7.8	118	0.00
45 T	Naphthalene	6.232	7.293	-17.0	134	0.00
46 T	Hexachlorobutadiene	1.216	1.214	0.2	108	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*Handwritten signature*  
 KR 3/31/12

## LABORATORY REPORT

April 12, 2012

Christopher Cote  
Montana DEQ  
1100 N. Last Chance Gulch  
Helena, MT 59601

### RE: Background Indoor Air Study

Dear Christopher:

Enclosed are the results of the samples submitted to our laboratory on March 29, 2012. For your reference, these analyses have been assigned our service request number P1201247.

All analyses were performed according to our laboratory's NELAP and DoD-ELAP-approved quality assurance program. The test results meet requirements of the current NELAP and DoD-ELAP standards, where applicable, and except as noted in the laboratory case narrative provided. For a specific list of NELAP and DoD-ELAP-accredited analytes, refer to the certifications section at [www.caslab.com](http://www.caslab.com). Results are intended to be considered in their entirety and apply only to the samples analyzed and reported herein.

Columbia Analytical Services, Inc. is certified by the California Department of Health Services, NELAP Laboratory Certificate No. 02115CA; Arizona Department of Health Services, Certificate No. AZ0694; Florida Department of Health, NELAP Certification E871020; New Jersey Department of Environmental Protection, NELAP Laboratory Certification ID #CA009; New York State Department of Health, NELAP NY Lab ID No: 11221; Oregon Environmental Laboratory Accreditation Program, NELAP ID: CA20007; The American Industrial Hygiene Association, Laboratory #101661; United States Department of Defense Environmental Laboratory Accreditation Program (DoD-ELAP), Certificate No. L11-203; Pennsylvania Registration No. 68-03307; TX Commission of Environmental Quality, NELAP ID T104704413-11-2; Minnesota Department of Health, NELAP Certificate No. 362188; Washington State Department of Ecology, ELAP Lab ID: C946, State of Utah Department of Health, NELAP Certificate No. CA015272011-1; Los Angeles Department of Building and Safety, Approval No: TA00001. Each of the certifications listed above have an explicit Scope of Accreditation that applies to specific matrices/methods/analytes; therefore, please contact me for information corresponding to a particular certification.

If you have any questions, please call me at (805) 526-7161.

Respectfully submitted,

**Columbia Analytical Services, Inc.**

Kate Aguilera  
Project Manager



Client: Montana DEQ  
Project: Background Indoor Air Study

CAS Project No: P1201247

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## CASE NARRATIVE

The samples were received intact under chain of custody on March 29, 2012 and were stored in accordance with the analytical method requirements. Please refer to the sample acceptance check form for additional information. The results reported herein are applicable only to the condition of the samples at the time of sample receipt.

### Air-Phase Petroleum Hydrocarbons (APH) Analysis

The samples were analyzed for total aliphatic and aromatic gasoline range hydrocarbons by gas chromatography/mass spectrometry according to the Method for the Determination of Air-Phase Petroleum Hydrocarbons (APH), Massachusetts Department of Environmental Protection, Revision 1, December, 2009.

### Volatile Organic Compound Analysis

The samples were also analyzed in Scan and SIM mode for selected volatile organic compounds in accordance with EPA Method TO-15 from the Compendium of Methods for the Determination of Toxic Organic Compounds in Ambient Air, Second Edition (EPA/625/R-96/010b), January, 1999. The analytical system was comprised of a gas chromatograph / mass spectrometer (GC/MS) interfaced to a whole-air preconcentrator.

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*The results of analyses are given in the attached laboratory report. All results are intended to be considered in their entirety, and Columbia Analytical Services, Inc. (CAS) is not responsible for utilization of less than the complete report.*

*Use of Columbia Analytical Services, Inc. (CAS) Name. Client shall not use CAS's name or trademark in any marketing or reporting materials, press releases or in any other manner ("Materials") whatsoever and shall not attribute to CAS any test result, tolerance or specification derived from CAS's data ("Attribution") without CAS's prior written consent, which may be withheld by CAS for any reason in its sole discretion. To request CAS's consent, Client shall provide copies of the proposed Materials or Attribution and describe in writing Client's proposed use of such Materials or Attribution. If CAS has not provided written approval of the Materials or Attribution within ten (10) days of receipt from Client, Client's request to use CAS's name or trademark in any Materials or Attribution shall be deemed denied. CAS may, in its discretion, reasonably charge Client for its time in reviewing Materials or Attribution requests. Client acknowledges and agrees that the unauthorized use of CAS's name or trademark may cause CAS to incur irreparable harm for which the recovery of money damages will be inadequate. Accordingly, Client acknowledges and agrees that a violation shall justify preliminary injunctive relief. For questions contact the laboratory.*

DETAIL SUMMARY REPORT

Client: Montana DEQ  
 Project ID: Background Indoor Air Study

Service Request: P1201247

Date Received: 3/29/2012  
 Time Received: 09:55

Client Sample ID	Lab Code	Matrix	Date Collected	Time Collected	Container ID	Pi1 (psig)	Pf1 (psig)	MA APH 1.0 - MA VOC PH Can	TO-15 - VOC Cans	TO-15 - VOC SIM
14-1F-032712	P1201247-001	Air	3/28/2012	12:00	AC01282	-5.66	3.56	X	X	X
25-1F-032312	P1201247-002	Air	3/24/2012	09:50	AC00949	-4.74	3.69	X	X	X
27-1F-032312	P1201247-003	Air	3/24/2012	11:02	AC01400	-5.15	3.57	X	X	X
24-1F-032112	P1201247-004	Air	3/22/2012	18:30	AS00183	-2.08	3.68	X	X	X

**Air - Chain of Custody Record & Analytical Service Request**

**Requested Turnaround Time in Business Days (Surcharges) please circle**  
 1 Day (100%) 2 Day (75%) 3 Day (50%) 4 Day (35%) 5 Day (25%) 10 Day-Standard

CAS Project No. **P1201247**

**Company Name & Address (Reporting Information)**  
 Monterey Department of Environmental Quality  
 PO Box 200901  
 1100 North Lawrence Ave  
 Helena, MT 59620-0901

**Project Name**  
 Background Air Study

**CAS Contact:**  
**Analysis Method**  
 TO-15  
 TO-15 SIM  
 MA-APH

**Project Manager**  
 Chris Cole

**PO # / Billing Information**  
 Reference #5: Background Indoor Air Study  
 Bill to Chris Cole of MDEQ

**Comments**  
 e.g. Actual Preservative or specific instructions

**Phone**  
 406-841-5078

**Fax**  
 406-841-5050

**Sampler (Print & Sign)**  
 Bob Roll Oh BBA

**Email Address for Result Reporting**  
 cote2@mt.gov; Markhbb@cedm.com

**Client Sample ID**

**Canister ID (Bar code # - AC, SC, etc.)**  
**Flow Controller ID (Bar code # - FC #)**  
**Canister Start Pressure "Hg**  
**Canister End Pressure "Hg(psig)**  
**Sample Volume**

Client Sample ID	Laboratory ID Number	Date Collected	Time Collected	Canister ID (Bar code # - AC, SC, etc.)	Flow Controller ID (Bar code # - FC #)	Canister Start Pressure "Hg	Canister End Pressure "Hg(psig)	Sample Volume	Analysis Method	Comments
14-1F-032712	0-518	3/25/12	1200	FA00067	-27.7"	-10"			TO-15	Analyze
25-1F-032312	0-414	3/24/12	950	FA00049	-27.5"	-9.2"			TO-15 SIM	Project Specific
27-1F-032312	0-517	3/24/12	11:02	FA00046	-28.5"	-10.5"			TO-15 SIM	Project Specific
24-1F-032112	0-210	3/22/12	18:30	FA00083	-27"	-3.5"			TO-15 SIM	List of Compounds to Project Specific Reporting Limits

**Report Tier Levels - please select**

Tier I - Results (Default if not specified) \_\_\_\_\_  
 Tier II (Results + QC Summaries) \_\_\_\_\_  
 Tier III (Results + QC & Calibration Summaries) **X**  
 Tier IV (Data Validation Package) 10% Surcharge \_\_\_\_\_

**Relinquished by: (Signature)** *Bob Roll* **Date:** 3/25/12 **Time:** 1400  
**Received by: (Signature)** *Markhbb* **Date:** 3/26/12 **Time:** 0935  
**Relinquished by: (Signature)** \_\_\_\_\_ **Date:** \_\_\_\_\_ **Time:** \_\_\_\_\_  
**Received by: (Signature)** \_\_\_\_\_ **Date:** \_\_\_\_\_ **Time:** \_\_\_\_\_

**Project Requirements (MRLs, QAPP)**  
 Cooler / Blank \_\_\_\_\_ °C  
 Temperature \_\_\_\_\_ °C

**Sample Acceptance Check Form**

Client: Montana DEQ Work order: P1201247

Project: Background Indoor Air Study

Sample(s) received on: 3/29/12 Date opened: 3/29/12 by: MZAMORA

**Note:** This form is used for all samples received by CAS. The use of this form for custody seals is strictly meant to indicate presence/absence and not as an indication of compliance or nonconformity. Thermal preservation and pH will only be evaluated either at the request of the client and/or as required by the method/SOP.

- |  | <u>Yes</u>                          | <u>No</u>                           | <u>N/A</u>                          |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| 1 Were <b>sample containers</b> properly marked with client sample ID?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 2 Container(s) <b>supplied by CAS</b> ?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 3 Did <b>sample containers</b> arrive in good condition?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 4 Were <b>chain-of-custody</b> papers used and filled out?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 5 Did <b>sample container labels</b> and/or tags agree with custody papers?                                      | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 6 Was <b>sample volume</b> received adequate for analysis?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 7 Are samples within specified holding times?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 8 Was proper <b>temperature</b> (thermal preservation) of cooler at receipt adhered to?                          | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 9 Was a <b>trip blank</b> received?  | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| 10 Were <b>custody seals</b> on outside of cooler/Box?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| Location of seal(s)? _____ Sealing Lid?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were signature and date included?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were seals intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were custody seals on outside of sample container?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| Location of seal(s)? _____ Sealing Lid?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were signature and date included?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were seals intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 11 Do containers have appropriate <b>preservation</b> , according to method/SOP or Client specified information? | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Is there a client indication that the submitted samples are <b>pH</b> preserved?                                 | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were <b>VOA vials</b> checked for presence/absence of air bubbles?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Does the client/method/SOP require that the analyst check the sample pH and <u>if necessary</u> alter it?        | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 12 <b>Tubes:</b> Are the tubes capped and intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Do they contain moisture?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 13 <b>Badges:</b> Are the badges properly capped and intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Are dual bed badges separated and individually capped and intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |

Lab Sample ID	Container Description	Required pH *	Received pH	Adjusted pH	VOA Headspace (Presence/Absence)	Receipt / Preservation Comments
P1201247-001.01	6.0 L Ambient Can					
P1201247-002.01	6.0 L Ambient Can					
P1201247-003.01	6.0 L Ambient Can					
P1201247-004.01	6.0 L Silonite Can					

Explain any discrepancies: (include lab sample ID numbers): \_\_\_\_\_

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 14-1F-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P1201247-001

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01282

Date Collected: 3/28/12  
 Date Received: 3/29/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.66      Final Pressure (psig): 3.56

Canister Dilution Factor: 2.02

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	84	40	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	39	20	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	10	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ

**Client Sample ID:** 25-1F-032312

**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247

CAS Sample ID: P1201247-002

Test Code: Massachusetts APH, Revision 1, December 2009

Date Collected: 3/24/12

Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8

Date Received: 3/29/12

Analyst: Elsa Moctezuma

Date Analyzed: 4/7/12

Sampling Media: 6.0 L Summa Canister

Volume(s) Analyzed: 1.00 Liter(s)

Test Notes:

Container ID: AC00949

Initial Pressure (psig): -4.74      Final Pressure (psig): 3.69

Canister Dilution Factor: 1.85

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	170	37	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	28	19	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	9.3	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 27-1F-032312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P1201247-003

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01400

Date Collected: 3/24/12  
 Date Received: 3/29/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 0.50 Liter(s)

Initial Pressure (psig): -5.15      Final Pressure (psig): 3.57

Canister Dilution Factor: 1.91

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	97	76	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	850	38	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	19	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 24-1F-032112  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P1201247-004

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00183

Date Collected: 3/22/12  
 Date Received: 3/29/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.08      Final Pressure (psig): 3.68

Canister Dilution Factor: 1.46

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	41	29	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	26	15	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	7.3	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P120407-MB

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/07/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	20	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	10	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	5.0	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
CAS Sample ID: P120407-LCS

Test Code: Massachusetts APH, Revision 1, December 2009  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/07/12  
Volume(s) Analyzed: NA Liter(s)

Compound	Spike Amount ng	Result ng	% Recovery	CAS	Data Qualifier
				Acceptance Limits	
C5 - C8 Aliphatic Hydrocarbons	24.5	25.1	102	70-130	
C9 - C12 Aliphatic Hydrocarbons	25.0	25.2	101	70-130	
C9 - C10 Aromatic Hydrocarbons	48.5	46.3	95	70-130	

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 14-1F-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
CAS Sample ID: P1201247-001

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01282

Date Collected: 3/28/12  
Date Received: 3/29/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.66 Final Pressure (psig): 3.56

Canister Dilution Factor: 2.02

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	1.7	1.0	1.0	0.59	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.1	1.0	0.42	0.20	
74-87-3	Chloromethane	ND	1.0	ND	0.49	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	1.0	ND	0.14	
106-99-0	1,3-Butadiene	ND	1.0	ND	0.46	
74-83-9	Bromomethane	ND	1.0	ND	0.26	
75-00-3	Chloroethane	ND	1.0	ND	0.38	
64-17-5	Ethanol	610	10	320	5.4	
75-05-8	Acetonitrile	ND	1.0	ND	0.60	
107-02-8	Acrolein	ND	4.0	ND	1.8	
67-64-1	Acetone	53	10	22	4.3	
75-69-4	Trichlorofluoromethane	1.1	1.0	0.20	0.18	
67-63-0	2-Propanol (Isopropyl Alcohol)	4.6	2.0	1.9	0.82	
107-13-1	Acrylonitrile	ND	1.0	ND	0.47	
75-35-4	1,1-Dichloroethene	ND	1.0	ND	0.25	
75-09-2	Methylene Chloride	ND	1.0	ND	0.29	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	1.0	ND	0.32	
76-13-1	Trichlorotrifluoroethane	ND	1.0	ND	0.13	
75-15-0	Carbon Disulfide	ND	10	ND	3.2	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	ND	0.25	
75-34-3	1,1-Dichloroethane	ND	1.0	ND	0.25	
1634-04-4	Methyl tert-Butyl Ether	ND	1.0	ND	0.28	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 14-1F-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
CAS Sample ID: P1201247-001

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01282

Date Collected: 3/28/12  
Date Received: 3/29/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.66 Final Pressure (psig): 3.56

Canister Dilution Factor: 2.02

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	10	ND	2.9	
78-93-3	2-Butanone (MEK)	ND	10	ND	3.4	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	ND	0.25	
141-78-6	Ethyl Acetate	7.6	2.0	2.1	0.56	
110-54-3	n-Hexane	1.3	1.0	0.36	0.29	
67-66-3	Chloroform	ND	1.0	ND	0.21	
109-99-9	Tetrahydrofuran (THF)	ND	1.0	ND	0.34	
71-55-6	1,1,1-Trichloroethane	ND	1.0	ND	0.19	
56-23-5	Carbon Tetrachloride	ND	1.0	ND	0.16	
110-82-7	Cyclohexane	ND	2.0	ND	0.59	
78-87-5	1,2-Dichloropropane	ND	1.0	ND	0.22	
75-27-4	Bromodichloromethane	ND	1.0	ND	0.15	
123-91-1	1,4-Dioxane	ND	1.0	ND	0.28	
80-62-6	Methyl Methacrylate	ND	2.0	ND	0.49	
142-82-5	n-Heptane	ND	1.0	ND	0.25	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	ND	0.22	
108-10-1	4-Methyl-2-pentanone	ND	1.0	ND	0.25	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	ND	0.22	
108-88-3	Toluene	5.0	1.0	1.3	0.27	
591-78-6	2-Hexanone	ND	1.0	ND	0.25	
124-48-1	Dibromochloromethane	ND	1.0	ND	0.12	
106-93-4	1,2-Dibromoethane	ND	1.0	ND	0.13	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 14-1F-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P1201247-001

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01282

**Date Collected:** 3/28/12  
**Date Received:** 3/29/12  
**Date Analyzed:** 4/7/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -5.66 Final Pressure (psig): 3.56

Canister Dilution Factor: 2.02

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	2.5	1.0	0.52	0.21	
111-65-9	n-Octane	ND	1.0	ND	0.22	
108-90-7	Chlorobenzene	ND	1.0	ND	0.22	
179601-23-1	m,p-Xylenes	2.7	2.0	0.63	0.47	
75-25-2	Bromoform	ND	1.0	ND	0.098	
100-42-5	Styrene	ND	1.0	ND	0.24	
95-47-6	o-Xylene	ND	1.0	ND	0.23	
111-84-2	n-Nonane	ND	1.0	ND	0.19	
98-82-8	Cumene	ND	1.0	ND	0.21	
80-56-8	alpha-Pinene	11	1.0	2.0	0.18	
103-65-1	n-Propylbenzene	ND	1.0	ND	0.21	
622-96-8	4-Ethyltoluene	ND	1.0	ND	0.21	
108-67-8	1,3,5-Trimethylbenzene	ND	1.0	ND	0.21	
95-63-6	1,2,4-Trimethylbenzene	1.1	1.0	0.23	0.21	
100-44-7	Benzyl Chloride	ND	1.0	ND	0.20	
541-73-1	1,3-Dichlorobenzene	ND	1.0	ND	0.17	
106-46-7	1,4-Dichlorobenzene	ND	1.0	ND	0.17	
95-50-1	1,2-Dichlorobenzene	ND	1.0	ND	0.17	
5989-27-5	d-Limonene	13	1.0	2.4	0.18	
96-12-8	1,2-Dibromo-3-chloropropane	ND	1.0	ND	0.10	
120-82-1	1,2,4-Trichlorobenzene	ND	1.0	ND	0.14	
87-68-3	Hexachlorobutadiene	ND	1.0	ND	0.095	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 25-1F-032312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
CAS Sample ID: P1201247-002

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00949

Date Collected: 3/24/12  
Date Received: 3/29/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.74 Final Pressure (psig): 3.69

Canister Dilution Factor: 1.85

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	8.5	0.93	4.9	0.54	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.1	0.93	0.43	0.19	
74-87-3	Chloromethane	ND	0.93	ND	0.45	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.93	ND	0.13	
106-99-0	1,3-Butadiene	ND	0.93	ND	0.42	
74-83-9	Bromomethane	ND	0.93	ND	0.24	
75-00-3	Chloroethane	ND	0.93	ND	0.35	
64-17-5	Ethanol	270	9.3	140	4.9	
75-05-8	Acetonitrile	ND	0.93	ND	0.55	
107-02-8	Acrolein	ND	3.7	ND	1.6	
67-64-1	Acetone	24	9.3	10	3.9	
75-69-4	Trichlorofluoromethane	1.3	0.93	0.23	0.16	
67-63-0	2-Propanol (Isopropyl Alcohol)	42	1.9	17	0.75	
107-13-1	Acrylonitrile	ND	0.93	ND	0.43	
75-35-4	1,1-Dichloroethene	ND	0.93	ND	0.23	
75-09-2	Methylene Chloride	ND	0.93	ND	0.27	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.93	ND	0.30	
76-13-1	Trichlorotrifluoroethane	ND	0.93	ND	0.12	
75-15-0	Carbon Disulfide	ND	9.3	ND	3.0	
156-60-5	trans-1,2-Dichloroethene	ND	0.93	ND	0.23	
75-34-3	1,1-Dichloroethane	ND	0.93	ND	0.23	
1634-04-4	Methyl tert-Butyl Ether	ND	0.93	ND	0.26	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 25-1F-032312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P1201247-002

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC00949

**Date Collected:** 3/24/12  
**Date Received:** 3/29/12  
**Date Analyzed:** 4/7/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -4.74      Final Pressure (psig): 3.69

Canister Dilution Factor: 1.85

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	9.3	ND	2.6	
78-93-3	2-Butanone (MEK)	ND	9.3	ND	3.1	
156-59-2	cis-1,2-Dichloroethene	ND	0.93	ND	0.23	
141-78-6	Ethyl Acetate	<b>3.2</b>	1.9	<b>0.89</b>	0.51	
110-54-3	n-Hexane	ND	0.93	ND	0.26	
67-66-3	Chloroform	ND	0.93	ND	0.19	
109-99-9	Tetrahydrofuran (THF)	ND	0.93	ND	0.31	
71-55-6	1,1,1-Trichloroethane	ND	0.93	ND	0.17	
56-23-5	Carbon Tetrachloride	ND	0.93	ND	0.15	
110-82-7	Cyclohexane	ND	1.9	ND	0.54	
78-87-5	1,2-Dichloropropane	ND	0.93	ND	0.20	
75-27-4	Bromodichloromethane	ND	0.93	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.93	ND	0.26	
80-62-6	Methyl Methacrylate	ND	1.9	ND	0.45	
142-82-5	n-Heptane	ND	0.93	ND	0.23	
10061-01-5	cis-1,3-Dichloropropene	ND	0.93	ND	0.20	
108-10-1	4-Methyl-2-pentanone	ND	0.93	ND	0.23	
10061-02-6	trans-1,3-Dichloropropene	ND	0.93	ND	0.20	
108-88-3	Toluene	<b>4.4</b>	0.93	<b>1.2</b>	0.25	
591-78-6	2-Hexanone	ND	0.93	ND	0.23	
124-48-1	Dibromochloromethane	ND	0.93	ND	0.11	
106-93-4	1,2-Dibromoethane	ND	0.93	ND	0.12	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 25-1F-032312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
CAS Sample ID: P1201247-002

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00949

Date Collected: 3/24/12  
Date Received: 3/29/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.74 Final Pressure (psig): 3.69

Canister Dilution Factor: 1.85

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.93	ND	0.19	
111-65-9	n-Octane	ND	0.93	ND	0.20	
108-90-7	Chlorobenzene	ND	0.93	ND	0.20	
179601-23-1	m,p-Xylenes	ND	1.9	ND	0.43	
75-25-2	Bromoform	ND	0.93	ND	0.089	
100-42-5	Styrene	ND	0.93	ND	0.22	
95-47-6	o-Xylene	ND	0.93	ND	0.21	
111-84-2	n-Nonane	ND	0.93	ND	0.18	
98-82-8	Cumene	ND	0.93	ND	0.19	
80-56-8	alpha-Pinene	1.7	0.93	0.31	0.17	
103-65-1	n-Propylbenzene	ND	0.93	ND	0.19	
622-96-8	4-Ethyltoluene	ND	0.93	ND	0.19	
108-67-8	1,3,5-Trimethylbenzene	ND	0.93	ND	0.19	
95-63-6	1,2,4-Trimethylbenzene	ND	0.93	ND	0.19	
100-44-7	Benzyl Chloride	ND	0.93	ND	0.18	
541-73-1	1,3-Dichlorobenzene	ND	0.93	ND	0.15	
106-46-7	1,4-Dichlorobenzene	ND	0.93	ND	0.15	
95-50-1	1,2-Dichlorobenzene	ND	0.93	ND	0.15	
5989-27-5	d-Limonene	11	0.93	2.0	0.17	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.93	ND	0.096	
120-82-1	1,2,4-Trichlorobenzene	ND	0.93	ND	0.12	
87-68-3	Hexachlorobutadiene	ND	0.93	ND	0.087	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 27-1F-032312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P1201247-003

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01400

Date Collected: 3/24/12  
 Date Received: 3/29/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 0.50 Liter(s)

Initial Pressure (psig): -5.15 Final Pressure (psig): 3.57

Canister Dilution Factor: 1.91

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	19	1.9	11	1.1	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.2	1.9	0.45	0.39	
74-87-3	Chloromethane	ND	1.9	ND	0.93	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	1.9	ND	0.27	
106-99-0	1,3-Butadiene	ND	1.9	ND	0.86	
74-83-9	Bromomethane	ND	1.9	ND	0.49	
75-00-3	Chloroethane	ND	1.9	ND	0.72	
64-17-5	Ethanol	170	19	93	10	
75-05-8	Acetonitrile	ND	1.9	ND	1.1	
107-02-8	Acrolein	ND	7.6	ND	3.3	
67-64-1	Acetone	71	19	30	8.0	
75-69-4	Trichlorofluoromethane	ND	1.9	ND	0.34	
67-63-0	2-Propanol (Isopropyl Alcohol)	91	3.8	37	1.6	
107-13-1	Acrylonitrile	ND	1.9	ND	0.88	
75-35-4	1,1-Dichloroethene	ND	1.9	ND	0.48	
75-09-2	Methylene Chloride	13	1.9	3.9	0.55	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	1.9	ND	0.61	
76-13-1	Trichlorotrifluoroethane	ND	1.9	ND	0.25	
75-15-0	Carbon Disulfide	ND	19	ND	6.1	
156-60-5	trans-1,2-Dichloroethene	ND	1.9	ND	0.48	
75-34-3	1,1-Dichloroethane	ND	1.9	ND	0.47	
1634-04-4	Methyl tert-Butyl Ether	ND	1.9	ND	0.53	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 27-1F-032312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P1201247-003

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01400

**Date Collected:** 3/24/12  
**Date Received:** 3/29/12  
**Date Analyzed:** 4/7/12  
**Volume(s) Analyzed:** 0.50 Liter(s)

Initial Pressure (psig): -5.15      Final Pressure (psig): 3.57

Canister Dilution Factor: 1.91

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	19	ND	5.4	
78-93-3	2-Butanone (MEK)	ND	19	ND	6.5	
156-59-2	cis-1,2-Dichloroethene	ND	1.9	ND	0.48	
141-78-6	Ethyl Acetate	ND	3.8	ND	1.1	
110-54-3	n-Hexane	ND	1.9	ND	0.54	
67-66-3	Chloroform	ND	1.9	ND	0.39	
109-99-9	Tetrahydrofuran (THF)	ND	1.9	ND	0.65	
71-55-6	1,1,1-Trichloroethane	ND	1.9	ND	0.35	
56-23-5	Carbon Tetrachloride	ND	1.9	ND	0.30	
110-82-7	Cyclohexane	ND	3.8	ND	1.1	
78-87-5	1,2-Dichloropropane	ND	1.9	ND	0.41	
75-27-4	Bromodichloromethane	ND	1.9	ND	0.29	
123-91-1	1,4-Dioxane	ND	1.9	ND	0.53	
80-62-6	Methyl Methacrylate	ND	3.8	ND	0.93	
142-82-5	n-Heptane	<b>4.5</b>	1.9	<b>1.1</b>	0.47	
10061-01-5	cis-1,3-Dichloropropene	ND	1.9	ND	0.42	
108-10-1	4-Methyl-2-pentanone	ND	1.9	ND	0.47	
10061-02-6	trans-1,3-Dichloropropene	ND	1.9	ND	0.42	
108-88-3	Toluene	<b>29</b>	1.9	<b>7.8</b>	0.51	
591-78-6	2-Hexanone	ND	1.9	ND	0.47	
124-48-1	Dibromochloromethane	ND	1.9	ND	0.22	
106-93-4	1,2-Dibromoethane	ND	1.9	ND	0.25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 27-1F-032312  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201247  
**CAS Sample ID:** P1201247-003

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01400

**Date Collected:** 3/24/12  
**Date Received:** 3/29/12  
**Date Analyzed:** 4/7/12  
**Volume(s) Analyzed:** 0.50 Liter(s)

Initial Pressure (psig): -5.15      Final Pressure (psig): 3.57

Canister Dilution Factor: 1.91

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	1.9	ND	0.40	
111-65-9	n-Octane	ND	1.9	ND	0.41	
108-90-7	Chlorobenzene	ND	1.9	ND	0.41	
179601-23-1	m,p-Xylenes	12	3.8	2.7	0.88	
75-25-2	Bromoform	ND	1.9	ND	0.18	
100-42-5	Styrene	ND	1.9	ND	0.45	
95-47-6	o-Xylene	2.8	1.9	0.65	0.44	
111-84-2	n-Nonane	ND	1.9	ND	0.36	
98-82-8	Cumene	ND	1.9	ND	0.39	
80-56-8	alpha-Pinene	2.7	1.9	0.48	0.34	
103-65-1	n-Propylbenzene	ND	1.9	ND	0.39	
622-96-8	4-Ethyltoluene	ND	1.9	ND	0.39	
108-67-8	1,3,5-Trimethylbenzene	ND	1.9	ND	0.39	
95-63-6	1,2,4-Trimethylbenzene	ND	1.9	ND	0.39	
100-44-7	Benzyl Chloride	ND	1.9	ND	0.37	
541-73-1	1,3-Dichlorobenzene	ND	1.9	ND	0.32	
106-46-7	1,4-Dichlorobenzene	ND	1.9	ND	0.32	
95-50-1	1,2-Dichlorobenzene	ND	1.9	ND	0.32	
5989-27-5	d-Limonene	12	1.9	2.1	0.34	
96-12-8	1,2-Dibromo-3-chloropropane	ND	1.9	ND	0.20	
120-82-1	1,2,4-Trichlorobenzene	ND	1.9	ND	0.26	
87-68-3	Hexachlorobutadiene	ND	1.9	ND	0.18	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 24-1F-032112  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
CAS Sample ID: P1201247-004

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00183

Date Collected: 3/22/12  
Date Received: 3/29/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.08 Final Pressure (psig): 3.68

Canister Dilution Factor: 1.46

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	<b>0.94</b>	0.73	<b>0.54</b>	0.42	
75-71-8	Dichlorodifluoromethane (CFC 12)	<b>2.1</b>	0.73	<b>0.42</b>	0.15	
74-87-3	Chloromethane	ND	0.73	ND	0.35	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.73	ND	0.10	
106-99-0	1,3-Butadiene	ND	0.73	ND	0.33	
74-83-9	Bromomethane	ND	0.73	ND	0.19	
75-00-3	Chloroethane	ND	0.73	ND	0.28	
64-17-5	Ethanol	<b>640</b>	7.3	<b>340</b>	3.9	
75-05-8	Acetonitrile	ND	0.73	ND	0.43	
107-02-8	Acrolein	ND	2.9	ND	1.3	
67-64-1	Acetone	<b>16</b>	7.3	<b>6.9</b>	3.1	
75-69-4	Trichlorofluoromethane	<b>2.0</b>	0.73	<b>0.36</b>	0.13	
67-63-0	2-Propanol (Isopropyl Alcohol)	<b>3.1</b>	1.5	<b>1.2</b>	0.59	
107-13-1	Acrylonitrile	ND	0.73	ND	0.34	
75-35-4	1,1-Dichloroethene	ND	0.73	ND	0.18	
75-09-2	Methylene Chloride	ND	0.73	ND	0.21	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.73	ND	0.23	
76-13-1	Trichlorotrifluoroethane	ND	0.73	ND	0.095	
75-15-0	Carbon Disulfide	ND	7.3	ND	2.3	
156-60-5	trans-1,2-Dichloroethene	ND	0.73	ND	0.18	
75-34-3	1,1-Dichloroethane	ND	0.73	ND	0.18	
1634-04-4	Methyl tert-Butyl Ether	ND	0.73	ND	0.20	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 24-1F-032112  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
CAS Sample ID: P1201247-004

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00183

Date Collected: 3/22/12  
Date Received: 3/29/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.08 Final Pressure (psig): 3.68

Canister Dilution Factor: 1.46

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	7.3	ND	2.1	
78-93-3	2-Butanone (MEK)	ND	7.3	ND	2.5	
156-59-2	cis-1,2-Dichloroethene	ND	0.73	ND	0.18	
141-78-6	Ethyl Acetate	1.8	1.5	0.49	0.41	
110-54-3	n-Hexane	1.0	0.73	0.30	0.21	
67-66-3	Chloroform	1.5	0.73	0.30	0.15	
109-99-9	Tetrahydrofuran (THF)	ND	0.73	ND	0.25	
71-55-6	1,1,1-Trichloroethane	ND	0.73	ND	0.13	
56-23-5	Carbon Tetrachloride	ND	0.73	ND	0.12	
110-82-7	Cyclohexane	ND	1.5	ND	0.42	
78-87-5	1,2-Dichloropropane	ND	0.73	ND	0.16	
75-27-4	Bromodichloromethane	ND	0.73	ND	0.11	
123-91-1	1,4-Dioxane	ND	0.73	ND	0.20	
80-62-6	Methyl Methacrylate	ND	1.5	ND	0.36	
142-82-5	n-Heptane	ND	0.73	ND	0.18	
10061-01-5	cis-1,3-Dichloropropene	ND	0.73	ND	0.16	
108-10-1	4-Methyl-2-pentanone	3.3	0.73	0.80	0.18	
10061-02-6	trans-1,3-Dichloropropene	ND	0.73	ND	0.16	
108-88-3	Toluene	3.3	0.73	0.86	0.19	
591-78-6	2-Hexanone	ND	0.73	ND	0.18	
124-48-1	Dibromochloromethane	ND	0.73	ND	0.086	
106-93-4	1,2-Dibromoethane	ND	0.73	ND	0.095	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 24-1F-032112  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
CAS Sample ID: P1201247-004

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00183

Date Collected: 3/22/12  
Date Received: 3/29/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.08 Final Pressure (psig): 3.68

Canister Dilution Factor: 1.46

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.73	ND	0.15	
111-65-9	n-Octane	ND	0.73	ND	0.16	
108-90-7	Chlorobenzene	ND	0.73	ND	0.16	
179601-23-1	m,p-Xylenes	ND	1.5	ND	0.34	
75-25-2	Bromoform	ND	0.73	ND	0.071	
100-42-5	Styrene	ND	0.73	ND	0.17	
95-47-6	o-Xylene	ND	0.73	ND	0.17	
111-84-2	n-Nonane	ND	0.73	ND	0.14	
98-82-8	Cumene	ND	0.73	ND	0.15	
80-56-8	alpha-Pinene	<b>1.6</b>	0.73	<b>0.29</b>	0.13	
103-65-1	n-Propylbenzene	ND	0.73	ND	0.15	
622-96-8	4-Ethyltoluene	ND	0.73	ND	0.15	
108-67-8	1,3,5-Trimethylbenzene	ND	0.73	ND	0.15	
95-63-6	1,2,4-Trimethylbenzene	ND	0.73	ND	0.15	
100-44-7	Benzyl Chloride	ND	0.73	ND	0.14	
541-73-1	1,3-Dichlorobenzene	ND	0.73	ND	0.12	
106-46-7	1,4-Dichlorobenzene	ND	0.73	ND	0.12	
95-50-1	1,2-Dichlorobenzene	ND	0.73	ND	0.12	
5989-27-5	d-Limonene	<b>23</b>	0.73	<b>4.1</b>	0.13	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.73	ND	0.076	
120-82-1	1,2,4-Trichlorobenzene	ND	0.73	ND	0.098	
87-68-3	Hexachlorobutadiene	ND	0.73	ND	0.068	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P120407-MB

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 4/7/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	ND	0.50	ND	0.29	
75-71-8	Dichlorodifluoromethane (CFC 12)	ND	0.50	ND	0.10	
74-87-3	Chloromethane	ND	0.50	ND	0.24	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.50	ND	0.072	
106-99-0	1,3-Butadiene	ND	0.50	ND	0.23	
74-83-9	Bromomethane	ND	0.50	ND	0.13	
75-00-3	Chloroethane	ND	0.50	ND	0.19	
64-17-5	Ethanol	ND	5.0	ND	2.7	
75-05-8	Acetonitrile	ND	0.50	ND	0.30	
107-02-8	Acrolein	ND	2.0	ND	0.87	
67-64-1	Acetone	ND	5.0	ND	2.1	
75-69-4	Trichlorofluoromethane	ND	0.50	ND	0.089	
67-63-0	2-Propanol (Isopropyl Alcohol)	ND	1.0	ND	0.41	
107-13-1	Acrylonitrile	ND	0.50	ND	0.23	
75-35-4	1,1-Dichloroethene	ND	0.50	ND	0.13	
75-09-2	Methylene Chloride	ND	0.50	ND	0.14	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.50	ND	0.16	
76-13-1	Trichlorotrifluoroethane	ND	0.50	ND	0.065	
75-15-0	Carbon Disulfide	ND	5.0	ND	1.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.50	ND	0.13	
75-34-3	1,1-Dichloroethane	ND	0.50	ND	0.12	
1634-04-4	Methyl tert-Butyl Ether	ND	0.50	ND	0.14	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P120407-MB

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 4/7/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	5.0	ND	1.4	
78-93-3	2-Butanone (MEK)	ND	5.0	ND	1.7	
156-59-2	cis-1,2-Dichloroethene	ND	0.50	ND	0.13	
141-78-6	Ethyl Acetate	ND	1.0	ND	0.28	
110-54-3	n-Hexane	ND	0.50	ND	0.14	
67-66-3	Chloroform	ND	0.50	ND	0.10	
109-99-9	Tetrahydrofuran (THF)	ND	0.50	ND	0.17	
71-55-6	1,1,1-Trichloroethane	ND	0.50	ND	0.092	
56-23-5	Carbon Tetrachloride	ND	0.50	ND	0.080	
110-82-7	Cyclohexane	ND	1.0	ND	0.29	
78-87-5	1,2-Dichloropropane	ND	0.50	ND	0.11	
75-27-4	Bromodichloromethane	ND	0.50	ND	0.075	
123-91-1	1,4-Dioxane	ND	0.50	ND	0.14	
80-62-6	Methyl Methacrylate	ND	1.0	ND	0.24	
142-82-5	n-Heptane	ND	0.50	ND	0.12	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-10-1	4-Methyl-2-pentanone	ND	0.50	ND	0.12	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-88-3	Toluene	ND	0.50	ND	0.13	
591-78-6	2-Hexanone	ND	0.50	ND	0.12	
124-48-1	Dibromochloromethane	ND	0.50	ND	0.059	
106-93-4	1,2-Dibromoethane	ND	0.50	ND	0.065	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P120407-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.50	ND	0.11	
111-65-9	n-Octane	ND	0.50	ND	0.11	
108-90-7	Chlorobenzene	ND	0.50	ND	0.11	
179601-23-1	m,p-Xylenes	ND	1.0	ND	0.23	
75-25-2	Bromoform	ND	0.50	ND	0.048	
100-42-5	Styrene	ND	0.50	ND	0.12	
95-47-6	o-Xylene	ND	0.50	ND	0.12	
111-84-2	n-Nonane	ND	0.50	ND	0.095	
98-82-8	Cumene	ND	0.50	ND	0.10	
80-56-8	alpha-Pinene	ND	0.50	ND	0.090	
103-65-1	n-Propylbenzene	ND	0.50	ND	0.10	
622-96-8	4-Ethyltoluene	ND	0.50	ND	0.10	
108-67-8	1,3,5-Trimethylbenzene	ND	0.50	ND	0.10	
95-63-6	1,2,4-Trimethylbenzene	ND	0.50	ND	0.10	
100-44-7	Benzyl Chloride	ND	0.50	ND	0.097	
541-73-1	1,3-Dichlorobenzene	ND	0.50	ND	0.083	
106-46-7	1,4-Dichlorobenzene	ND	0.50	ND	0.083	
95-50-1	1,2-Dichlorobenzene	ND	0.50	ND	0.083	
5989-27-5	d-Limonene	ND	0.50	ND	0.090	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.50	ND	0.052	
120-82-1	1,2,4-Trichlorobenzene	ND	0.50	ND	0.067	
87-68-3	Hexachlorobutadiene	ND	0.50	ND	0.047	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister(s)  
 Test Notes:

Date(s) Collected: 3/22 - 3/28/12  
 Date(s) Received: 3/29/12  
 Date(s) Analyzed: 4/7/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		Percent Recovered	Percent Recovered	Percent Recovered		
Method Blank	P120407-MB	106	99	97	70-130	
Lab Control Sample	P120407-LCS	103	99	100	70-130	
14-1F-032712	P1201247-001	106	98	96	70-130	
25-1F-032312	P1201247-002	105	99	97	70-130	
25-1F-032312	P1201247-002DUP	106	98	97	70-130	
27-1F-032312	P1201247-003	107	99	97	70-130	
24-1F-032112	P1201247-004	108	102	93	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.

Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P120407-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 4/07/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
115-07-1	Propene	198	215	109	59-137	
75-71-8	Dichlorodifluoromethane (CFC 12)	196	170	87	63-115	
74-87-3	Chloromethane	190	173	91	59-124	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	200	170	85	65-113	
106-99-0	1,3-Butadiene	204	205	100	60-138	
74-83-9	Bromomethane	194	174	90	69-129	
75-00-3	Chloroethane	196	174	89	60-120	
64-17-5	Ethanol	928	861	93	58-121	
75-05-8	Acetonitrile	194	194	100	64-129	
107-02-8	Acrolein	198	183	92	54-127	
67-64-1	Acetone	1,010	922	91	59-114	
75-69-4	Trichlorofluoromethane	202	177	88	66-108	
67-63-0	2-Propanol (Isopropyl Alcohol)	382	299	78	50-113	
107-13-1	Acrylonitrile	198	199	101	72-135	
75-35-4	1,1-Dichloroethene	212	198	93	70-117	
75-09-2	Methylene Chloride	206	183	89	61-108	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	208	217	104	70-131	
76-13-1	Trichlorotrifluoroethane	206	191	93	70-113	
75-15-0	Carbon Disulfide	208	186	89	65-112	
156-60-5	trans-1,2-Dichloroethene	196	188	96	71-119	
75-34-3	1,1-Dichloroethane	200	192	96	71-116	
1634-04-4	Methyl tert-Butyl Ether	198	193	97	67-116	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
CAS Sample ID: P120407-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/07/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
108-05-4	Vinyl Acetate	952	992	104	59-142	
78-93-3	2-Butanone (MEK)	206	205	100	68-125	
156-59-2	cis-1,2-Dichloroethene	206	195	95	69-119	
141-78-6	Ethyl Acetate	398	391	98	63-130	
110-54-3	n-Hexane	198	180	91	57-120	
67-66-3	Chloroform	214	192	90	69-111	
109-99-9	Tetrahydrofuran (THF)	202	191	95	57-123	
71-55-6	1,1,1-Trichloroethane	198	186	94	73-119	
56-23-5	Carbon Tetrachloride	202	194	96	74-129	
110-82-7	Cyclohexane	390	355	91	70-113	
78-87-5	1,2-Dichloropropane	198	183	92	69-118	
75-27-4	Bromodichloromethane	198	188	95	75-124	
123-91-1	1,4-Dioxane	200	189	95	71-123	
80-62-6	Methyl Methacrylate	400	390	98	72-127	
142-82-5	n-Heptane	196	188	96	68-120	
10061-01-5	cis-1,3-Dichloropropene	188	183	97	71-130	
108-10-1	4-Methyl-2-pentanone	204	204	100	69-130	
10061-02-6	trans-1,3-Dichloropropene	210	212	101	76-133	
108-88-3	Toluene	202	179	89	67-111	
591-78-6	2-Hexanone	222	213	96	70-123	
124-48-1	Dibromochloromethane	206	197	96	75-129	
106-93-4	1,2-Dibromoethane	200	180	90	73-122	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
CAS Sample ID: P120407-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/07/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
123-86-4	n-Butyl Acetate	220	211	96	68-132	
111-65-9	n-Octane	200	195	98	68-116	
108-90-7	Chlorobenzene	202	177	88	69-113	
179601-23-1	m,p-Xylenes	392	360	92	70-116	
75-25-2	Bromoform	208	188	90	69-127	
100-42-5	Styrene	200	188	94	71-125	
95-47-6	o-Xylene	194	176	91	70-116	
111-84-2	n-Nonane	196	184	94	68-116	
98-82-8	Cumene	190	174	92	70-116	
80-56-8	alpha-Pinene	186	174	94	71-119	
103-65-1	n-Propylbenzene	192	172	90	71-119	
622-96-8	4-Ethyltoluene	198	174	88	71-119	
108-67-8	1,3,5-Trimethylbenzene	200	181	91	71-121	
95-63-6	1,2,4-Trimethylbenzene	194	180	93	73-127	
100-44-7	Benzyl Chloride	200	206	103	65-137	
541-73-1	1,3-Dichlorobenzene	200	164	82	68-123	
106-46-7	1,4-Dichlorobenzene	206	161	78	65-120	
95-50-1	1,2-Dichlorobenzene	198	163	82	67-121	
5989-27-5	d-Limonene	200	189	95	67-130	
96-12-8	1,2-Dibromo-3-chloropropane	196	181	92	72-133	
120-82-1	1,2,4-Trichlorobenzene	196	163	83	62-133	
87-68-3	Hexachlorobutadiene	202	166	82	60-128	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

## LABORATORY DUPLICATE SUMMARY RESULTS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 25-1F-032312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P1201247-002DUP

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC00949

**Date Collected:** 3/24/12  
**Date Received:** 3/29/12  
**Date Analyzed:** 4/7/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -4.74

Final Pressure (psig): 3.69

Canister Dilution Factor: 1.85

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
Propene	8.48	4.93	7.56	4.39	8.02	11	25	
Dichlorodifluoromethane (CFC 12)	2.11	0.427	2.07	0.419	2.09	2	25	
Chloromethane	ND	ND	ND	ND	-	-	25	
1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	ND	ND	ND	-	-	25	
1,3-Butadiene	ND	ND	ND	ND	-	-	25	
Bromomethane	ND	ND	ND	ND	-	-	25	
Chloroethane	ND	ND	ND	ND	-	-	25	
Ethanol	272	144	265	141	268.5	3	25	
Acetonitrile	ND	ND	ND	ND	-	-	25	
Acrolein	ND	ND	ND	ND	-	-	25	
Acetone	24.3	10.2	23.7	9.97	24	3	25	
Trichlorofluoromethane	1.30	0.231	1.24	0.220	1.27	5	25	
2-Propanol (Isopropyl Alcohol)	42.2	17.2	42.9	17.5	42.55	2	25	
Acrylonitrile	ND	ND	ND	ND	-	-	25	
1,1-Dichloroethene	ND	ND	ND	ND	-	-	25	
Methylene Chloride	ND	ND	ND	ND	-	-	25	
3-Chloro-1-propene (Allyl Chloride)	ND	ND	ND	ND	-	-	25	
Trichlorotrifluoroethane	ND	ND	ND	ND	-	-	25	
Carbon Disulfide	ND	ND	ND	ND	-	-	25	
trans-1,2-Dichloroethene	ND	ND	ND	ND	-	-	25	
1,1-Dichloroethane	ND	ND	ND	ND	-	-	25	
Methyl tert-Butyl Ether	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

## LABORATORY DUPLICATE SUMMARY RESULTS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 25-1F-032312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P1201247-002DUP

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC00949

**Date Collected:** 3/24/12  
**Date Received:** 3/29/12  
**Date Analyzed:** 4/7/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -4.74

Final Pressure (psig): 3.69

Canister Dilution Factor: 1.85

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
Vinyl Acetate	ND	ND	ND	ND	-	-	25	
2-Butanone (MEK)	ND	ND	ND	ND	-	-	25	
cis-1,2-Dichloroethene	ND	ND	ND	ND	-	-	25	
Ethyl Acetate	3.21	0.892	3.10	0.862	3.155	3	25	
n-Hexane	ND	ND	ND	ND	-	-	25	
Chloroform	ND	ND	ND	ND	-	-	25	
Tetrahydrofuran (THF)	ND	ND	ND	ND	-	-	25	
1,1,1-Trichloroethane	ND	ND	ND	ND	-	-	25	
Carbon Tetrachloride	ND	ND	ND	ND	-	-	25	
Cyclohexane	ND	ND	ND	ND	-	-	25	
1,2-Dichloropropane	ND	ND	ND	ND	-	-	25	
Bromodichloromethane	ND	ND	ND	ND	-	-	25	
1,4-Dioxane	ND	ND	ND	ND	-	-	25	
Methyl Methacrylate	ND	ND	ND	ND	-	-	25	
n-Heptane	ND	ND	ND	ND	-	-	25	
cis-1,3-Dichloropropene	ND	ND	ND	ND	-	-	25	
4-Methyl-2-pentanone	ND	ND	ND	ND	-	-	25	
trans-1,3-Dichloropropene	ND	ND	ND	ND	-	-	25	
Toluene	4.35	1.16	4.30	1.14	4.325	1	25	
2-Hexanone	ND	ND	ND	ND	-	-	25	
Dibromochloromethane	ND	ND	ND	ND	-	-	25	
1,2-Dibromoethane	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

LABORATORY DUPLICATE SUMMARY RESULTS

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 25-1F-032312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P1201247-002DUP

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00949

Date Collected: 3/24/12  
 Date Received: 3/29/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.74

Final Pressure (psig): 3.69

Canister Dilution Factor: 1.85

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
n-Butyl Acetate	ND	ND	ND	ND	-	-	25	
n-Octane	ND	ND	ND	ND	-	-	25	
Chlorobenzene	ND	ND	ND	ND	-	-	25	
m,p-Xylenes	ND	ND	ND	ND	-	-	25	
Bromoform	ND	ND	ND	ND	-	-	25	
Styrene	ND	ND	ND	ND	-	-	25	
o-Xylene	ND	ND	ND	ND	-	-	25	
n-Nonane	ND	ND	ND	ND	-	-	25	
Cumene	ND	ND	ND	ND	-	-	25	
alpha-Pinene	1.72	0.310	1.70	0.306	1.71	<b>1</b>	25	
n-Propylbenzene	ND	ND	ND	ND	-	-	25	
4-Ethyltoluene	ND	ND	ND	ND	-	-	25	
1,3,5-Trimethylbenzene	ND	ND	ND	ND	-	-	25	
1,2,4-Trimethylbenzene	ND	ND	ND	ND	-	-	25	
Benzyl Chloride	ND	ND	ND	ND	-	-	25	
1,3-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
1,4-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
1,2-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
d-Limonene	10.9	1.95	10.8	1.94	10.85	<b>0.9</b>	25	
1,2-Dibromo-3-chloropropane	ND	ND	ND	ND	-	-	25	
1,2,4-Trichlorobenzene	ND	ND	ND	ND	-	-	25	
Hexachlorobutadiene	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.



RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 14-1F-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P1201247-001

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01282

Date Collected: 3/28/12  
 Date Received: 3/29/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.66      Final Pressure (psig): 3.56

Canister Dilution Factor: 2.02

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.051	ND	0.020	
107-06-2	1,2-Dichloroethane	<b>0.15</b>	0.051	<b>0.038</b>	0.012	
71-43-2	Benzene	<b>1.1</b>	0.15	<b>0.35</b>	0.047	
79-01-6	Trichloroethene	<b>0.058</b>	0.051	<b>0.011</b>	0.0094	
79-00-5	1,1,2-Trichloroethane	ND	0.20	ND	0.037	
127-18-4	Tetrachloroethene	ND	0.051	ND	0.0075	
100-41-4	Ethylbenzene	<b>0.69</b>	0.20	<b>0.16</b>	0.047	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.051	ND	0.0074	
91-20-3	Naphthalene	<b>0.39</b>	0.20	<b>0.075</b>	0.039	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 25-1F-032312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P1201247-002

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00949

Date Collected: 3/24/12  
 Date Received: 3/29/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.74      Final Pressure (psig): 3.69

Canister Dilution Factor: 1.85

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.046	ND	0.018	
107-06-2	1,2-Dichloroethane	<b>0.089</b>	0.046	<b>0.022</b>	0.011	
71-43-2	Benzene	<b>0.48</b>	0.14	<b>0.15</b>	0.043	
79-01-6	Trichloroethene	ND	0.046	ND	0.0086	
79-00-5	1,1,2-Trichloroethane	ND	0.19	ND	0.034	
127-18-4	Tetrachloroethene	ND	0.046	ND	0.0068	
100-41-4	Ethylbenzene	<b>0.22</b>	0.19	<b>0.050</b>	0.043	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.046	ND	0.0067	
91-20-3	Naphthalene	<b>0.22</b>	0.19	<b>0.041</b>	0.035	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 27-1F-032312  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P1201247-003

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01400

Date Collected: 3/24/12  
 Date Received: 3/29/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 0.50 Liter(s)

Initial Pressure (psig): -5.15      Final Pressure (psig): 3.57

Canister Dilution Factor: 1.91

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.096	ND	0.037	
107-06-2	1,2-Dichloroethane	<b>0.17</b>	0.096	<b>0.042</b>	0.024	
71-43-2	Benzene	<b>1.1</b>	0.29	<b>0.36</b>	0.090	
79-01-6	Trichloroethene	ND	0.096	ND	0.018	
79-00-5	1,1,2-Trichloroethane	ND	0.38	ND	0.070	
127-18-4	Tetrachloroethene	<b>0.10</b>	0.096	<b>0.015</b>	0.014	
100-41-4	Ethylbenzene	<b>2.5</b>	0.38	<b>0.58</b>	0.088	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.096	ND	0.014	
91-20-3	Naphthalene	ND	0.38	ND	0.073	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 24-1F-032112  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
CAS Sample ID: P1201247-004

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00183

Date Collected: 3/22/12  
Date Received: 3/29/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.08      Final Pressure (psig): 3.68

Canister Dilution Factor: 1.46

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.037	ND	0.014	
107-06-2	1,2-Dichloroethane	<b>0.11</b>	0.037	<b>0.026</b>	0.0090	
71-43-2	Benzene	<b>0.35</b>	0.11	<b>0.11</b>	0.034	
79-01-6	Trichloroethene	<b>0.068</b>	0.037	<b>0.013</b>	0.0068	
79-00-5	1,1,2-Trichloroethane	ND	0.15	ND	0.027	
127-18-4	Tetrachloroethene	<b>0.097</b>	0.037	<b>0.014</b>	0.0054	
100-41-4	Ethylbenzene	<b>0.18</b>	0.15	<b>0.041</b>	0.034	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.037	ND	0.0053	
91-20-3	Naphthalene	ND	0.15	ND	0.028	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P120409-MB

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.025	ND	0.0098	
107-06-2	1,2-Dichloroethane	ND	0.025	ND	0.0062	
71-43-2	Benzene	ND	0.075	ND	0.023	
79-01-6	Trichloroethene	ND	0.025	ND	0.0047	
79-00-5	1,1,2-Trichloroethane	ND	0.10	ND	0.018	
127-18-4	Tetrachloroethene	ND	0.025	ND	0.0037	
100-41-4	Ethylbenzene	ND	0.10	ND	0.023	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.025	ND	0.0036	
91-20-3	Naphthalene	ND	0.10	ND	0.019	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister(s)  
 Test Notes:

Date(s) Collected: 3/22 - 3/28/12  
 Date(s) Received: 3/29/12  
 Date(s) Analyzed: 4/9/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		% Recovered	% Recovered	% Recovered		
Method Blank	P120409-MB	101	110	108	70-130	
Lab Control Sample	P120409-LCS	102	110	110	70-130	
14-1F-032712	P1201247-001	99	112	107	70-130	
14-1F-032712	P1201247-001DUP	101	114	106	70-130	
25-1F-032312	P1201247-002	100	112	108	70-130	
27-1F-032312	P1201247-003	100	110	107	70-130	
24-1F-032112	P1201247-004	99	110	106	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.

Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
 CAS Sample ID: P120409-LCS

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/09/12  
 Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
75-01-4	Vinyl Chloride	3.88	3.08	79	56-127	
107-06-2	1,2-Dichloroethane	4.00	3.13	78	51-140	
71-43-2	Benzene	3.96	3.51	89	56-125	
79-01-6	Trichloroethene	3.88	2.97	77	51-127	
79-00-5	1,1,2-Trichloroethane	3.92	2.89	74	49-137	
127-18-4	Tetrachloroethene	3.68	2.91	79	58-134	
100-41-4	Ethylbenzene	3.96	3.18	80	56-126	
79-34-5	1,1,2,2-Tetrachloroethane	3.84	3.04	79	53-148	
91-20-3	Naphthalene	3.44	3.08	90	25-149	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY DUPLICATE SUMMARY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 14-1F-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201247  
CAS Sample ID: P1201247-001DUP

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01282

Date Collected: 3/28/12  
Date Received: 3/29/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.66      Final Pressure (psig): 3.56

Canister Dilution Factor: 2.02

CAS #	Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
		µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
75-01-4	Vinyl Chloride	ND	ND	ND	ND	-	-	25	
107-06-2	1,2-Dichloroethane	0.154	0.0381	0.161	0.0398	0.1575	4	25	
71-43-2	Benzene	1.12	0.351	1.15	0.361	1.135	3	25	
79-01-6	Trichloroethene	0.0585	0.0109	0.0601	0.0112	0.0593	3	25	
79-00-5	1,1,2-Trichloroethane	ND	ND	ND	ND	-	-	25	
127-18-4	Tetrachloroethene	ND	ND	ND	ND	-	-	25	
100-41-4	Ethylbenzene	0.690	0.159	0.713	0.164	0.7015	3	25	
79-34-5	1,1,2,2-Tetrachloroethane	ND	ND	ND	ND	-	-	25	
91-20-3	Naphthalene	0.392	0.0749	0.397	0.0758	0.3945	1	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.



Massachusetts APH  
Hydrocarbon Ranges

ICAL Method: M8011412A.M

ICAL Date: 1/14/12

Instrument ID: MS08

	areas				masses							
	0.5	1	5	25	50	100	0.5	1	5	25	50	100
<b>C5-C8 Aliphatics</b>												
Isopentane	60478	112606	488243	2480105	4710733	9222808	0.520	1.04	5.20	26.0	52.0	104
n-Hexane	59953	110631	507126	2577865	4931102	9678199	0.490	0.98	4.90	24.5	49.0	98
Cyclohexane	63203	115230	548910	2874726	5524402	10890911	0.500	1.00	5.00	25.0	50.0	100
2,3-Dimethylpentane	68648	127222	585183	2998017	5742104	11319597	0.520	1.04	5.20	26.0	52.0	104
n-Heptane	66143	120286	558074	2884817	5509801	10873046	0.490	0.98	4.90	24.5	49.0	98
n-Octane	76169	137829	632891	3286017	6205035	12092514	0.505	1.01	5.05	25.3	50.5	101
<b>area sum:</b>	<b>394594</b>	<b>723804</b>	<b>3320427</b>	<b>17101547</b>	<b>32623177</b>	<b>64077075</b>	<b>3.025</b>	<b>6.050</b>	<b>30.25</b>	<b>151.30</b>	<b>302.50</b>	<b>605.00</b>

<b>C9-C12 Aliphatics</b>												
2,3-Dimethylheptane	78780	147724	682518	3501487	6691428	13148840	0.485	0.97	4.85	24.3	48.5	97
n-Nonane	77464	140680	662221	3417128	6555165	12940625	0.485	0.97	4.85	24.3	48.5	97
n-Decane	80593	147724	705259	3651028	7021424	13870530	0.490	0.98	4.90	24.5	49.0	98
n-Butylcyclohexane	88915	166782	776903	4026931	7749936	15415340	0.495	0.99	4.95	24.8	49.5	99
n-Undecane	80867	152557	719618	3772331	7263077	14385881	0.480	0.96	4.80	24.0	48.0	96
n-Dodecane	70045	143746	687255	3673524	7115012	14041583	0.500	1.00	5.00	25.0	50.0	100
<b>area sum:</b>	<b>476664</b>	<b>899213</b>	<b>4233774</b>	<b>22042429</b>	<b>42396042</b>	<b>83802799</b>	<b>2.935</b>	<b>5.870</b>	<b>29.35</b>	<b>146.90</b>	<b>293.50</b>	<b>587.00</b>

<b>C9-C10 Aromatics</b>												
Isopropylbenzene	10658	19169	87439	448287	849124	1695737	0.485	0.97	4.85	24.3	48.5	97
3-Ethyltoluene	11805	20564	97823	504510	977020	1973178	0.470	0.94	4.70	23.5	47.0	94
1,3,5-Trimethylbenzene	15759	29260	134073	700133	1353305	2726630	0.490	0.98	4.90	24.5	49.0	98
1,2,3-Trimethylbenzene	17195	31328	143473	757946	1493373	3063059	0.495	0.99	4.95	24.8	49.5	99
p-Isopropyltoluene	9490	18116	83401	440962	872799	1814314	0.460	0.92	4.60	23.0	46.0	92
<b>area sum:</b>	<b>64907</b>	<b>118437</b>	<b>546209</b>	<b>2851838</b>	<b>5545621</b>	<b>11272918</b>	<b>2.400</b>	<b>4.800</b>	<b>24.00</b>	<b>120.10</b>	<b>240.00</b>	<b>480.00</b>

Massachusetts APH  
Hydrocarbon Ranges

ICAL: M8011412A.M

ICAL Date: 1/14/12

Instrument ID: MS08

areas

<u>Internal Standards (TIC)</u>	0.5	1	5	25	50	100
Bromochloromethane (IS1)	1312793	1307309	1293586	1289283	1284194	1283817
1,4-Difluorobenzene (IS2)	2547079	2541159	2512498	2475255	2422119	2352159
Chlorobenzene-d5 (IS3)	2929189	2913374	2880258	2895080	2878314	2908396

Internal Standards (EIC)

Bromochloromethane (IS1)	249696	247899	245985	245585	244327	243561
1,4-Difluorobenzene (IS2)	1153657	1149947	1133233	1136223	1127203	1138730
Chlorobenzene-d5 (IS3)	1007702	1004638	991212	992925	982812	986950

Surrogates (TIC)

	0.5	1	5	25	50	100
1,2-Dichloroethane-d4	1162497	1154617	1105630	1126201	1103958	1095945
o-Toluene-d8	3508018	3490493	3455382	3466688	3437986	3457323
p-Bromofluorobenzene	2298220	2284797	2197239	2227396	2192159	2197999

C5-C8 Aliphatics

	<u>RRFs</u>						
0.5	1	5	25	50	100	<u>RRF<sub>avg</sub></u>	<u>%RSD</u>
2.8268	2.6009	2.4215	2.4870	2.3919	2.3252	2.509	7.24

C9-C12 Aliphatics

0.5	1	5	25	50	100	<u>RRF<sub>avg</sub></u>	<u>%RSD</u>
4.0291	3.8120	3.6383	3.7780	3.6744	3.6163	3.758	4.09

C9-C10 Aromatics

0.5	1	5	25	50	100	<u>RRF<sub>avg</sub></u>	<u>%RSD</u>
0.6709	0.6140	0.5740	0.5979	0.5878	0.5949	0.607	5.63

**Massachusetts APH**  
**Continuing Calibration Verification Check Sheet**

Data File Name: 04071202.D  
 Data File Path: J:\MS08\Data\2012\_04\06\  
 Operator: EM  
 Date Acquired: 4/7/12 3:06  
 Acq. Method File: TO15.M  
 Sample Name: 25ng TO-15/MAPH CCV STD  
 Misc Info: S25-04031203/S25-03161202  
 Instrument Name: MS08

Enter RRFs from current ICAL!

Internal Standards	RT	Area
7) 1,4-Difluorobenzene (IS2)	15.89	1082412
16) Chlorobenzene-d5 (IS3)	21.69	941687

C5-C8 Aliphatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
3) Isopentane	8.19	2388606	2.538	153.0	1.14	-30	30	Pass
4) n-Hexane	13.08	2507742						
9) Cyclohexane	15.81	2792389						
10) 2,3-Dimethylpentane	16.17	2922324	Spike	ICAL				
11) n-Heptane	17.35	2804136	Amt (ng)	RRF				
14) n-Octane	20.69	3208272	151.30	2.509				
		16623469						

C9-C12 Aliphatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
18) 2,3-Dimethylheptane	22.31	3401295	3.834	149.9	2.02	-30	30	Pass
19) n-Nonane	23.30	3317221						
25) n-Decane	25.26	3533579						
28) Butylcyclohexane	25.99	3853565	Spike	ICAL				
29) n-Undecane	26.76	3619770	Amt (ng)	RRF				
30) n-Dodecane	28.00	3488040	146.90	3.758				
		21213470						

C9-C10 Aromatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
22) Isopropylbenzene	23.78	439879	0.601	118.9	-1.03	-30	30	Pass
23) 1-Methyl-3-ethylbenzene	24.52	482296						
24) 1,3,5-Trimethylbenzene	24.66	665021						
26) p-Isopropyltoluene	25.68	416582	Spike	ICAL				
27) 1,2,3-Trimethylbenzene	25.68	714045	Amt (ng)	RRF				
		2717823	120.1	0.607				

*Em 4/7/12*

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Mon Dec 12 09:45:45 2011  
 Response Via : Initial Calibration

Calibration Files

0.1 =12091118.D 0.2 =12091119.D 0.5 =12091120.D 1.0 =12091121.D 5.0 =12091122.D 25 =12091123.D  
 50 =12091124.D 100 =12091125.D

Compound	0.1	0.2	0.5	1.0	5.0	25	50	100	Avg	%RSD
1) IR Bromochloromethane...				ISTD						
2) T Propene	2.487	1.965	1.952	1.692	2.194	2.303	2.185	2.036	2.102	11.62
3) T Dichlorodifluo...	3.545	2.805	2.940	2.748	2.432	2.361	2.278	2.199	2.664	16.73
4) T Chloromethane	3.290	2.811	2.875	2.512	2.176	2.286	1.977	1.664	2.449	21.62
5) T 1,2-Dichloro-1...	1.901	1.661	1.575	1.458	1.359	1.333	1.200	1.190	1.460	16.68
6) T Vinyl Chloride	2.906	2.590	2.631	2.374	2.145	2.129	1.950	1.939	2.333	15.07
7) T 1,3-Butadiene	2.145	1.788	1.790	1.718	1.581	1.689	1.621	1.616	1.743	10.33
8) T Bromomethane	1.687	1.328	1.349	1.212	1.094	1.142	1.111	1.092	1.252	16.21
9) T Chloroethane	1.671	1.259	1.333	1.193	1.101	1.091	1.046	1.044	1.217	17.29
10) T Ethanol	2.040	1.345	1.186	1.085	1.022	1.096	1.076	1.064	1.239	27.34
11) T Acetonitrile	4.365	2.973	2.723	2.546	2.480	2.562	2.482	2.468	2.825	22.85
12) T Acrolein	1.328	0.931	0.907	0.861	0.772	0.814	0.789	0.779	0.897	20.48
13) T Acetone	1.645	1.316	1.249	1.129	0.994	1.015	0.981	0.964	1.161	20.26
14) T Trichlorofluor...	2.776	2.385	2.496	2.272	2.097	2.135	2.042	2.011	2.277	11.59
15) T 2-Propanol (IS...		4.463	4.536	4.168	2.920	2.763	2.577	2.672	3.443	26.09
16) T Acrylonitrile	2.701	1.907	1.878	1.783	1.745	1.813	1.765	1.738	1.916	16.85
17) T 1,1-Dichloroet...	1.458	1.418	1.452	1.263	1.206	1.223	1.172	1.152	1.293	9.95
18) T 2-Methyl-2-Pro...	4.464	3.710	3.939	3.616	3.558	3.584			3.812	9.14
19) T Methylene Chlo...			1.958	1.578	1.294	1.273	1.220	1.201	1.421	20.88
20) T 3-Chloro-1-pro...	2.347	2.260	2.345	2.220	2.150	2.296	2.230	2.191	2.255	3.15
21) T Trichlorotrifl...	1.348	1.244	1.299	1.193	1.089	1.098	1.062	1.035	1.171	10.01
22) T Carbon Disulfide		6.210	5.678	5.057	4.654	4.699	4.488	4.424	5.030	13.40
23) T trans-1,2-Dich...	2.436	2.237	2.206	2.142	1.951	1.953	1.883	1.835	2.080	10.00
24) T 1,1-Dichloroet...	2.896	2.757	2.747	2.500	2.307	2.379	2.288	2.236	2.514	10.07
25) T Methyl tert-Bu...	5.128	4.611	4.650	4.307	4.119	4.185	4.042	3.984	4.378	8.93
26) T Vinyl Acetate	0.430	0.368	0.423	0.399	0.393	0.410	0.402	0.398	0.403	4.70
27) T 2-Butanone (MEK)	1.265	1.003	1.038	0.972	0.927	0.964	0.928	0.794	0.986	13.56
28) T cis-1,2-Dichlo...	2.442	2.056	2.101	1.938	1.826	1.836	1.764	1.720	1.960	12.05
29) T Diisopropyl Ether	1.446	1.223	1.323	1.262	1.171	1.188	1.153	1.148	1.239	8.26
30) T Ethyl Acetate	0.611	0.554	0.573	0.545	0.517	0.535	0.520	0.516	0.546	6.02
31) T n-Hexane	3.541	3.006	3.181	2.805	2.582	2.633	2.545	2.543	2.855	12.70

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M

Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

32) T	Chloroform	2.679	2.408	2.409	2.204	1.983	2.054	1.974	1.937	2.206	12.18
33) S	1,2-Dichloroet...	1.586	1.577	1.582	1.590	1.535	1.533	1.520	1.504	1.553	2.19
34) T	Tetrahydrofura...	1.274	0.932	1.100	0.938	0.910	0.879	0.830	0.821	0.961	15.96
35) T	Ethyl tert-But...	1.999	1.832	2.033	1.818	1.732	1.757	1.708	1.687	1.821	7.18
36) T	1,2-Dichloroet...	2.099	1.992	1.976	1.806	1.683	1.700	1.599	1.531	1.798	11.39
37) IR	1,4-Difluorobenzen...										
	ISTD										
38) T	1,1,1-Trichlor...	0.531	0.465	0.486	0.453	0.419	0.434	0.413	0.403	0.451	9.46
39) T	Isopropyl Acetate	0.227	0.205	0.220	0.206	0.194	0.201	0.192	0.188	0.204	6.71
40) T	1-Butanol		0.348	0.319	0.300	0.320	0.335	0.324	0.318	0.323	4.66
41) T	Benzene	1.633	1.385	1.359	1.227	1.136	1.158	1.101	1.070	1.258	15.11
42) T	Carbon Tetrach...	0.394	0.356	0.369	0.340	0.334	0.352	0.337	0.333	0.352	6.03
43) T	Cyclohexane	0.666	0.576	0.588	0.539	0.492	0.500	0.479	0.469	0.539	12.59
44) T	tert-Amyl Meth...	1.092	0.926	0.998	0.937	0.900	0.918	0.881	0.858	0.939	7.95
45) T	1,2-Dichloropr...	0.397	0.361	0.368	0.324	0.312	0.316	0.300	0.293	0.334	11.12
46) T	Bromodichlorom...	0.458	0.387	0.392	0.378	0.362	0.378	0.362	0.353	0.384	8.57
47) T	Trichloroethene	0.466	0.375	0.379	0.344	0.323	0.328	0.316	0.309	0.355	14.62
48) T	1,4-Dioxane	0.305	0.282	0.288	0.248	0.245	0.246	0.237	0.230	0.260	10.55
49) T	2,2,4-Trimethy...	1.964	1.629	1.686	1.546	1.447	1.461	1.387	1.337	1.557	12.98
50) T	Methyl Methacr...	0.156	0.139	0.150	0.135	0.136	0.142	0.136	0.135	0.141	5.53
51) T	n-Heptane	0.427	0.373	0.385	0.364	0.341	0.344	0.332	0.324	0.361	9.35
52) T	cis-1,3-Dichlo...	0.571	0.502	0.521	0.491	0.482	0.500	0.481	0.471	0.502	6.33
53) T	4-Methyl-2-pen...	0.359	0.295	0.308	0.303	0.297	0.303	0.289	0.284	0.305	7.68
54) T	trans-1,3-Dich...	0.476	0.413	0.443	0.421	0.429	0.453	0.437	0.427	0.437	4.59
55) T	1,1,2-Trichlor...	0.340	0.301	0.310	0.284	0.276	0.285	0.271	0.265	0.292	8.44
56) IR	Chlorobenzene-d5 (...)										
	ISTD										
57) S	Toluene-d8 (SS2)	2.305	2.306	2.305	2.309	2.306	2.300	2.310	2.304	2.306	0.14
58) T	Toluene	3.623	2.920	3.025	2.738	2.571	2.574	2.473	2.421	2.793	14.19
59) T	2-Hexanone	2.049	1.429	1.515	1.403	1.395	1.422	1.367	1.337	1.490	15.57
60) T	Dibromochlorom...	0.696	0.618	0.636	0.636	0.612	0.648	0.629	0.625	0.638	4.14
61) T	1,2-Dibromoethane	0.831	0.730	0.718	0.658	0.641	0.662	0.638	0.628	0.688	9.97
62) T	n-Butyl Acetate	2.338	1.850	1.857	1.721	1.704	1.765	1.737	1.737	1.839	11.40
63) T	n-Octane	0.785	0.652	0.714	0.643	0.602	0.611	0.580	0.572	0.645	11.28
64) T	Tetrachloroethene	1.090	0.904	0.904	0.826	0.767	0.779	0.756	0.755	0.848	13.63
65) T	Chlorobenzene	2.375	1.868	1.886	1.744	1.610	1.637	1.574	1.548	1.780	15.30
66) T	Ethylbenzene	3.846	3.171	3.188	2.956	2.803	2.848	2.743	2.680	3.030	12.51
67) T	m- & p-Xylenes	3.081	2.499	2.611	2.392	2.216	2.273	2.170	2.112	2.419	13.08
68) T	Bromoform	0.622	0.494	0.556	0.541	0.551	0.598	0.589	0.594	0.568	7.19
69) T	Styrene	2.478	1.895	1.937	1.800	1.735	1.803	1.726	1.675	1.881	13.64
70) T	o-Xylene	3.279	2.708	2.726	2.479	2.331	2.373	2.283	2.224	2.550	13.65

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M

Title	EPA TO-15 per SOP	VOA-TO15	(CASS TO-15/GC-MS)	1.518	1.470	1.681	10.69
71) T n-Nonane	2.005	1.753	1.847	1.675	1.581	1.597	1.518
72) T 1,1,2,2-Tetrac...	1.348	1.164	1.262	1.150	1.083	1.121	1.085
73) S Bromofluoroben...	0.774	0.771	0.779	0.781	0.775	0.778	0.782
74) T Cumene	4.522	3.495	3.524	3.169	2.962	3.002	2.884
75) T alpha-Pinene	1.782	1.636	1.734	1.601	1.537	1.578	1.518
76) T n-Propylbenzene	5.441	4.229	4.306	3.924	3.710	3.754	3.553
77) T 3-Ethyltoluene	4.091	3.373	3.461	3.038	2.964	2.981	2.893
78) T 4-Ethyltoluene	4.003	3.202	3.301	3.151	2.837	2.983	2.810
79) T 1,3,5-Trimethy...	3.324	2.662	2.723	2.537	2.378	2.426	2.344
80) T alpha-Methylst...	1.742	1.386	1.449	1.371	1.390	1.402	1.354
81) T 2-Ethyltoluene	4.143	3.362	3.430	3.169	2.975	3.048	2.919
82) T 1,2,4-Trimethy...	3.356	2.663	2.791	2.551	2.433	2.513	2.454
83) T n-Decane	1.969	1.702	1.844	1.705	1.582	1.626	1.554
84) T Benzyl Chloride	2.373	1.648	1.772	1.688	1.799	2.058	2.054
85) T 1,3-Dichlorobe...	2.498	1.658	1.649	1.455	1.377	1.409	1.374
86) T 1,4-Dichlorobe...	2.655	1.785	1.618	1.492	1.380	1.428	1.388
87) T sec-Butylbenzene	4.527	3.773	3.816	3.471	3.264	3.328	3.214
88) T 4-Isopropyltol...	4.426	3.674	3.683	3.416	3.254	3.356	3.253
89) T 1,2,3-Trimethy...	3.266	2.650	2.852	2.588	2.471	2.573	2.506
90) T 1,2-Dichlorobe...	2.173	1.589	1.545	1.435	1.332	1.387	1.370
91) T d-Limonene	1.124	0.998	1.106	1.069	1.064	1.072	1.040
92) T 1,2-Dibromo-3-...	0.563	0.488	0.501	0.486	0.494	0.525	0.532
93) T n-Undecane	1.922	1.637	1.755	1.671	1.608	1.630	1.608
94) T 1,2,4-Trichlor...		1.335	1.165	1.062	1.017	1.071	1.086
95) T Naphthalene		4.848	3.923	3.638	3.497	3.721	3.755
96) T n-Dodecane	1.866	1.599	1.669	1.586	1.545	1.575	1.553
97) T Hexachlorobuta...	0.841	0.749	0.730	0.673	0.635	0.664	0.663
98) T Cyclohexanone	1.172	0.979	1.021	0.969	0.952	0.962	0.931
99) T tert-Butylbenzene	3.486	2.799	2.880	2.658	2.449	2.533	2.453
100) T n-Butylbenzene	3.596	2.904	2.958	2.670	2.572	2.625	2.613

(#) = Out of Range

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_04\07\  
 Data File : 04071201.D  
 Acq On : 7 Apr 2012 2:23  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-04031203/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 07 07:01:39 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	IR Bromochloromethane (IS1)	1.000	1.000	0.0	82	-0.02
2	T Propene	2.102	2.305	-9.7	83	0.00
3	T Dichlorodifluoromethane (CF	2.664	2.312	13.2	81	0.00
4	T Chloromethane	2.449	2.172	11.3	78	0.00
5	T 1,2-Dichloro-1,1,2,2-tetra	1.460	1.212	17.0	75	0.00
6	T Vinyl Chloride	2.333	2.037	12.7	79	0.00
7	T 1,3-Butadiene	1.743	1.733	0.6	85	0.00
8	T Bromomethane	1.252	1.113	11.1	80	0.00
9	T Chloroethane	1.217	1.092	10.3	83	0.00
10	T Ethanol	1.239	1.182	4.6	89	-0.07
11	T Acetonitrile	2.825	2.653	6.1	85	-0.05
12	T Acrolein	0.897	0.829	7.6	84	-0.02
13	T Acetone	1.161	1.067	8.1	87	-0.05
14	T Trichlorofluoromethane	2.277	2.119	6.9	82	0.00
15	T 2-Propanol (Isopropanol)	3.443	2.761	19.8	82	-0.05
16	T Acrylonitrile	1.916	1.842	3.9	84	-0.03
17	T 1,1-Dichloroethene	1.293	1.185	8.4	80	0.00
18	T 2-Methyl-2-Propanol (tert-B	3.812	3.075	19.3	71	-0.03
19	T Methylene Chloride	1.421	1.250	12.0	81	-0.01
20	T 3-Chloro-1-propene (Allyl C	2.255	2.354	-4.4	85	-0.02
21	T Trichlorotrifluoroethane	1.171	1.055	9.9	79	0.00
22	T Carbon Disulfide	5.030	4.592	8.7	81	0.00
23	T trans-1,2-Dichloroethene	2.080	1.974	5.1	83	-0.01
24	T 1,1-Dichloroethane	2.514	2.417	3.9	84	-0.02
25	T Methyl tert-Butyl Ether	4.378	4.248	3.0	84	0.00
26	T Vinyl Acetate	0.403	0.414	-2.7	83	-0.03
27	T 2-Butanone (MEK)	0.986	0.977	0.9	84	-0.03
28	T cis-1,2-Dichloroethene	1.960	1.856	5.3	83	-0.01
29	T Diisopropyl Ether	1.239	1.196	3.5	83	-0.01
30	T Ethyl Acetate	0.546	0.542	0.7	84	-0.03
31	T n-Hexane	2.855	2.627	8.0	82	-0.01
32	T Chloroform	2.206	2.031	7.9	82	-0.03
33	S 1,2-Dichloroethane-d4 (SS1)	1.553	1.635	-5.3	88	-0.02
34	T Tetrahydrofuran (THF)	0.961	0.905	5.8	85	-0.01
35	T Ethyl tert-Butyl Ether	1.821	1.755	3.6	82	-0.01
36	T 1,2-Dichloroethane	1.798	1.708	5.0	83	-0.02
37	IR 1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	87	-0.01
38	T 1,1,1-Trichloroethane	0.451	0.412	8.6	82	-0.01

*Em* 4/7/12

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_04\07\  
 Data File : 04071201.D  
 Acq On : 7 Apr 2012 2:23  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-04031203/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 07 07:01:39 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
39 T	Isopropyl Acetate	0.204	0.194	4.9	83	-0.02
40 T	1-Butanol	0.323	0.356	-10.2	92	-0.05
41 T	Benzene	1.258	1.097	12.8	82	-0.02
42 T	Carbon Tetrachloride	0.352	0.326	7.4	80	-0.01
43 T	Cyclohexane	0.539	0.482	10.6	83	-0.02
44 T	tert-Amyl Methyl Ether	0.939	0.875	6.8	82	-0.02
45 T	1,2-Dichloropropane	0.334	0.303	9.3	83	-0.02
46 T	Bromodichloromethane	0.384	0.357	7.0	82	-0.01
47 T	Trichloroethene	0.355	0.302	14.9	80	-0.02
48 T	1,4-Dioxane	0.260	0.238	8.5	84	-0.02
49 T	2,2,4-Trimethylpentane (Iso	1.557	1.423	8.6	84	-0.01
50 T	Methyl Methacrylate	0.141	0.134	5.0	82	-0.03
51 T	n-Heptane	0.361	0.336	6.9	84	-0.02
52 T	cis-1,3-Dichloropropene	0.502	0.476	5.2	82	-0.01
53 T	4-Methyl-2-pentanone	0.305	0.298	2.3	85	-0.02
54 T	trans-1,3-Dichloropropene	0.437	0.429	1.8	82	-0.01
55 T	1,1,2-Trichloroethane	0.292	0.268	8.2	81	-0.02
56 IR	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	87	0.00
57 S	Toluene-d8 (SS2)	2.306	2.271	1.5	86	0.00
58 T	Toluene	2.793	2.427	13.1	82	-0.01
59 T	2-Hexanone	1.490	1.418	4.8	87	-0.02
60 T	Dibromochloromethane	0.638	0.594	6.9	80	-0.01
61 T	1,2-Dibromoethane	0.688	0.609	11.5	80	-0.01
62 T	n-Butyl Acetate	1.839	1.752	4.7	86	-0.02
63 T	n-Octane	0.645	0.620	3.9	88	-0.01
64 T	Tetrachloroethene	0.848	0.702	17.2	78	0.00
65 T	Chlorobenzene	1.780	1.527	14.2	81	0.00
66 T	Ethylbenzene	3.030	2.719	10.3	83	0.00
67 T	m- & p-Xylenes	2.419	2.179	9.9	83	-0.02
68 T	Bromoform	0.568	0.540	4.9	79	-0.01
69 T	Styrene	1.881	1.741	7.4	84	-0.01
70 T	o-Xylene	2.550	2.268	11.1	83	-0.01
71 T	n-Nonane	1.681	1.549	7.9	84	-0.01
72 T	1,1,2,2-Tetrachloroethane	1.160	1.053	9.2	82	-0.02
73 S	Bromofluorobenzene (SS3)	0.777	0.774	0.4	87	0.00
74 T	Cumene	3.299	2.975	9.8	86	-0.01
75 T	alpha-Pinene	1.611	1.486	7.8	82	-0.01
76 T	n-Propylbenzene	4.039	3.560	11.9	82	-0.01

*Em* 4/7/12



Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_04\07\  
 Data File : 04071201.D  
 Acq On : 7 Apr 2012 2:23  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-04031203/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 07 07:01:39 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
77 T	3-Ethyltoluene	3.206	2.844	11.3	83	-0.01
78 T	4-Ethyltoluene	3.133	2.708	13.6	79	-0.02
79 T	1,3,5-Trimethylbenzene	2.600	2.285	12.1	82	-0.01
80 T	alpha-Methylstyrene	1.436	1.322	7.9	82	-0.01
81 T	2-Ethyltoluene	3.254	2.841	12.7	81	-0.01
82 T	1,2,4-Trimethylbenzene	2.664	2.354	11.6	81	-0.02
83 T	n-Decane	1.699	1.553	8.6	83	-0.02
84 T	Benzyl Chloride	1.948	2.052	-5.3	87	-0.02
85 T	1,3-Dichlorobenzene	1.607	1.281	20.3	79	-0.01
86 T	1,4-Dichlorobenzene	1.649	1.304	20.9	79	-0.02
87 T	sec-Butylbenzene	3.571	3.140	12.1	82	-0.01
88 T	4-Isopropyltoluene (p-Cymen	3.554	3.145	11.5	82	-0.01
89 T	1,2,3-Trimethylbenzene	2.691	2.401	10.8	81	-0.01
90 T	1,2-Dichlorobenzene	1.531	1.260	17.7	79	-0.01
91 T	d-Limonene	1.067	1.048	1.8	85	-0.01
92 T	1,2-Dibromo-3-Chloropropane	0.517	0.480	7.2	80	-0.01
93 T	n-Undecane	1.678	1.570	6.4	84	-0.01
94 T	1,2,4-Trichlorobenzene	1.122	0.972	13.4	79	-0.01
95 T	Naphthalene	3.881	3.485	10.2	81	-0.01
96 T	n-Dodecane	1.613	1.578	2.2	87	0.00
97 T	Hexachlorobutadiene	0.704	0.592	15.9	78	0.00
98 T	Cyclohexanone	0.986	0.996	-1.0	90	-0.02
99 T	tert-Butylbenzene	2.723	2.389	12.3	82	-0.02
100 T	n-Butylbenzene	2.817	2.452	13.0	81	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*EM 4/7/12*

Response Factor Report MS07

Method : J:\Ms07\METHODS\X7021712B.M (RTE Integrator)  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
 1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	500	AVG	%RSD
1) I Bromochloromethan					ISTD							
2) T Dichlorodifluorom	3.784	3.527	3.467	3.822	3.279	2.986	2.875	3.155	2.920		3.313	10.83
3) T Chloromethane		1.470	0.956	1.190	0.904	0.778	0.569	0.897	0.827		0.949	28.81
4) T Vinyl Chloride	2.802	2.749	2.588	3.044	2.597	2.406	2.314	2.588	2.446		2.615	8.57
5) T Bromomethane	1.894	1.496	1.211	1.613	1.335	1.208	1.105	1.433	1.367		1.407	17.11
6) T Chloroethane	1.511	1.416	1.357	1.517	1.295	1.237	1.195	1.312	1.224		1.340	8.92
7) T Acetone				1.548	1.214	1.122	1.080	1.193	1.090		1.208	14.52
8) T Trichlorofluorome	3.051	2.887	2.831	3.145	2.692	2.565	2.447	2.691	2.536		2.761	8.59
9) T 1,1-Dichloroethen	1.591	1.440	1.366	1.547	1.322	1.229	1.214	1.385	1.295		1.376	9.52
10) T Methylene Chlorid		2.134	1.795	1.866	1.536	1.426	1.404	1.494	1.411		1.633	16.43
11) T Trichlorotrifluor	1.408	1.325	1.295	1.376	1.169	1.081	1.062	1.170	1.097		1.220	10.87
12) T trans-1,2-Dichlor	1.972	1.710	1.589	1.762	1.505	1.406	1.393	1.512	1.439		1.588	12.18
13) T 1,1-Dichloroethan	3.265	3.020	3.016	3.383	2.925	2.756	2.681	2.937	2.841		2.980	7.60
14) T Methyl tert-Butyl	5.168	4.740	4.762	5.045	4.335	4.137	4.166	4.690	4.471		4.613	7.90
15) T cis-1,2-Dichloroe	1.885	1.626	1.630	1.776	1.524	1.434	1.388	1.575	1.482		1.591	10.09
16) T Chloroform			3.041	3.257	2.586	2.390	2.280	2.524	2.374		2.636	14.03
17) S 1,2-Dichloroethan	2.000	2.002	2.032	2.018	1.989	1.973	1.960	1.938	1.932		1.983	1.75
18) T 1,2-Dichloroethan	2.582	2.460	2.375	2.604	2.222	2.090	1.988	2.222	2.094		2.293	9.73
19) T 1,1,1-Trichloroet	2.587	2.431	2.340	2.573	2.191	2.059	2.007	2.278	2.150		2.291	9.18
20) T Benzene			7.536	7.534	6.055	5.665	5.395	5.903	5.468		6.222	14.88
21) T Carbon Tetrachlor	2.280	1.842	1.769	1.950	1.676	1.579	1.549	1.771	1.692		1.790	12.41
22) I 1,4-Difluorobenze					ISTD							
23) T 1,2-Dichloropropa	0.417	0.397	0.373	0.409	0.354	0.331	0.319	0.358	0.336		0.366	9.68
24) T Bromodichlorometh	0.491	0.464	0.436	0.489	0.420	0.398	0.394	0.447	0.416		0.439	8.26
25) T Trichloroethene	0.418	0.379	0.349	0.390	0.324	0.305	0.287	0.326	0.307		0.343	12.96
26) T 1,4-Dioxane	0.324	0.290	0.270	0.278	0.242	0.231	0.227	0.258	0.243		0.263	11.99
27) T cis-1,3-Dichlorop	0.567	0.523	0.508	0.540	0.499	0.484	0.488	0.559	0.522		0.521	5.67
28) T trans-1,3-Dichlor	0.457	0.434	0.431	0.444	0.419	0.409	0.417	0.505	0.482		0.444	7.22

(#) = Out of Range ### Number of calibration levels exceeded format ###  
 X7021712B.M Mon Apr 09 08:16:11 2012

Response Factor Report MS07

Method : J:\Ms07\METHODS\X7021712B.M (RTE Integrator)  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
 1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	Avg	%RSD
29) T 1,1,2-Trichloroet	0.315	0.296	0.286	0.305	0.263	0.248	0.242	0.274	0.257	0.276	9.32
30) S Toluene-d8 (SS2)	1.053	1.071	1.054	1.047	1.051	1.052	1.071	1.060	1.049	1.056	0.84
31) T Toluene	1.782	1.541	1.420	1.485	1.269	1.198	1.177	1.307	1.186	1.374	14.76
32) T 1,2-Dibromoethane	0.379	0.362	0.345	0.362	0.323	0.307	0.304	0.351	0.331	0.340	7.65
33) T Tetrachloroethene	0.423	0.394	0.371	0.400	0.341	0.320	0.311	0.347	0.326	0.359	11.03
-----ISTD-----											
34) I Chlorobenzene-d5											
35) T Chlorobenzene	4.581	4.172	4.017	4.308	3.688	3.469	3.311	3.705	3.345	3.844	11.68
36) T Ethylbenzene	7.604	7.082	6.707	7.235	6.297	6.070	5.915	6.682	5.878	6.608	9.30
37) T m,p-Xylene	5.861	5.347	5.132	5.536	4.930	4.777	4.719	5.268	4.558	5.125	8.24
38) T o-Xylene	6.505	5.826	5.526	5.889	5.188	4.995	4.950	5.594	4.932	5.490	9.67
39) T 1,1,2,2-Tetrachlo	2.655	2.674	2.710	2.646	2.437	2.371	2.298	2.761	2.498	2.561	6.40
40) S Bromofluorobenzene	1.712	1.701	1.695	1.697	1.683	1.680	1.670	1.642	1.593	1.675	2.20
41) T 1,3-Dichlorobenze	3.304	3.261	3.222	3.204	2.835	2.719	2.608	3.138	2.846	3.015	8.71
42) T 1,4-Dichlorobenze	3.470	3.303	3.208	3.208	2.803	2.671	2.579	3.174	2.891	3.034	10.13
43) T 1,2-Dichlorobenze	3.271	3.123	3.112	3.026	2.662	2.536	2.418	3.053	2.772	2.886	10.31
44) T 1,2,4-Trichlorobe	2.184	2.118	2.169	1.753	1.641	1.608	1.575	2.139	1.978	1.907	13.66
45) T Naphthalene	6.857	6.355	6.704	5.143	5.107	5.172	5.297	8.126	7.328	6.232	17.82
46) T Hexachlorobutadie	1.486	1.423	1.403	1.217	1.055	1.013	0.975	1.227	1.145	1.216	15.40

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_04\09\  
 Data File : 04091203.D  
 Acq On : 9 Apr 2012 7:01 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00163  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Apr 09 08:15:00 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane (IS1)	1.000	1.000	0.0	82	0.00
2 T	Dichlorodifluoromethane (CF)	3.313	2.539	23.4	63	0.00
3 T	Chloromethane	0.949	0.738	22.2	67	0.00
4 T	Vinyl Chloride	2.615	2.065	21.0	65	0.00
5 T	Bromomethane	1.407	1.190	15.4	73	0.00
6 T	Chloroethane	1.340	1.065	20.5	67	0.00
7 T	Acetone	1.208	1.043	13.7	70	0.00
8 T	Trichlorofluoromethane	2.761	2.169	21.4	66	0.00
9 T	1,1-Dichloroethene	1.376	1.147	16.6	71	0.00
10 T	Methylene Chloride	1.633	1.261	22.8	67	0.00
11 T	Trichlorotrifluoroethane	1.220	1.046	14.3	73	0.00
12 T	trans-1,2-Dichloroethene	1.588	1.249	21.3	68	0.00
13 T	1,1-Dichloroethane	2.980	2.362	20.7	66	0.00
14 T	Methyl tert-Butyl Ether	4.613	3.609	21.8	68	0.00
15 T	cis-1,2-Dichloroethene	1.591	1.330	16.4	71	0.00
16 T	Chloroform	2.636	2.237	15.1	71	0.00
17 S	1,2-Dichloroethane-d4 (SS1)	1.983	1.917	3.3	79	0.00
18 T	1,2-Dichloroethane	2.293	1.706	25.6	63	0.00
19 T	1,1,1-Trichloroethane	2.291	1.879	18.0	70	0.00
20 T	Benzene	6.222	5.361	13.8	72	0.00
21 T	Carbon Tetrachloride	1.790	1.425	20.4	69	0.00
22 I	1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	88	0.00
23 T	1,2-Dichloropropane	0.366	0.268	26.8	67	0.00
24 T	Bromodichloromethane	0.439	0.322	26.7	68	0.00
25 T	Trichloroethene	0.343	0.255	25.7	70	0.00
26 T	1,4-Dioxane	0.263	0.198	24.7	73	0.00
27 T	cis-1,3-Dichloropropene	0.521	0.396	24.0	70	0.00
28 T	trans-1,3-Dichloropropene	0.444	0.343	22.7	72	0.00
29 T	1,1,2-Trichloroethane	0.276	0.207	25.0	70	0.00
30 S	Toluene-d8 (SS2)	1.056	1.179	-11.6	99	0.00
31 T	Toluene	1.374	1.051	23.5	73	0.00
32 T	1,2-Dibromoethane	0.340	0.264	22.4	72	0.00
33 T	Tetrachloroethene	0.359	0.287	20.1	75	0.00
34 I	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	85	0.00
35 T	Chlorobenzene	3.844	3.219	16.3	75	0.00
36 T	Ethylbenzene	6.608	5.416	18.0	74	0.00
37 T	m,p-Xylene	5.125	4.308	15.9	75	0.00

*Handwritten signature/initials*

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_04\09\  
 Data File : 04091203.D  
 Acq On : 9 Apr 2012 7:01 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00163  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Apr 09 08:15:00 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
38 T	o-Xylene	5.490	4.561	16.9	75	0.00
39 T	1,1,2,2-Tetrachloroethane	2.561	2.098	18.1	74	0.00
40 S	Bromofluorobenzene (SS3)	1.675	1.861	-11.1	95	0.00
41 T	1,3-Dichlorobenzene	3.015	2.704	10.3	82	0.00
42 T	1,4-Dichlorobenzene	3.034	2.684	11.5	82	0.00
43 T	1,2-Dichlorobenzene	2.886	2.603	9.8	84	0.00
44 T	1,2,4-Trichlorobenzene	1.907	1.808	5.2	94	0.00
45 T	Naphthalene	6.232	6.283	-0.8	105	0.00
46 T	Hexachlorobutadiene	1.216	1.083	10.9	88	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*KR 4/9/12*

## LABORATORY REPORT

April 18, 2012

Christopher Cote  
Montana DEQ  
1100 N. Last Chance Gulch  
Helena, MT 59601

### **RE: Background Indoor Air Study**

Dear Christopher:

Enclosed are the results of the samples submitted to our laboratory on April 4, 2012. For your reference, these analyses have been assigned our service request number P1201319.

All analyses were performed according to our laboratory's NELAP and DoD-ELAP-approved quality assurance program. The test results meet requirements of the current NELAP and DoD-ELAP standards, where applicable, and except as noted in the laboratory case narrative provided. For a specific list of NELAP and DoD-ELAP-accredited analytes, refer to the certifications section at [www.caslab.com](http://www.caslab.com). Results are intended to be considered in their entirety and apply only to the samples analyzed and reported herein.

Columbia Analytical Services, Inc. is certified by the California Department of Health Services, NELAP Laboratory Certificate No. 02115CA; Arizona Department of Health Services, Certificate No. AZ0694; Florida Department of Health, NELAP Certification E871020; New Jersey Department of Environmental Protection, NELAP Laboratory Certification ID #CA009; New York State Department of Health, NELAP NY Lab ID No: 11221; Oregon Environmental Laboratory Accreditation Program, NELAP ID: CA20007; The American Industrial Hygiene Association, Laboratory #101661; United States Department of Defense Environmental Laboratory Accreditation Program (DoD-ELAP), Certificate No. L11-203; Pennsylvania Registration No. 68-03307; TX Commission of Environmental Quality, NELAP ID T104704413-11-2; Minnesota Department of Health, NELAP Certificate No. 362188; Washington State Department of Ecology, ELAP Lab ID: C946, State of Utah Department of Health, NELAP Certificate No. CA015272011-1; Los Angeles Department of Building and Safety, Approval No: TA00001. Each of the certifications listed above have an explicit Scope of Accreditation that applies to specific matrices/methods/analytes; therefore, please contact me for information corresponding to a particular certification.

If you have any questions, please call me at (805) 526-7161.

Respectfully submitted,

**Columbia Analytical Services, Inc.**

Kate Aguilera  
Project Manager

Client: Montana DEQ  
Project: Background Indoor Air Study

CAS Project No: P1201319

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## CASE NARRATIVE

The samples were received intact under chain of custody on April 4, 2012 and were stored in accordance with the analytical method requirements. Please refer to the sample acceptance check form for additional information. The results reported herein are applicable only to the condition of the samples at the time of sample receipt.

### Air-Phase Petroleum Hydrocarbons (APH) Analysis

The samples were analyzed for total aliphatic and aromatic gasoline range hydrocarbons by gas chromatography/mass spectrometry according to the Method for the Determination of Air-Phase Petroleum Hydrocarbons (APH), Massachusetts Department of Environmental Protection, Revision 1, December, 2009.

### Volatile Organic Compound Analysis

The samples were also analyzed in scan and SIM mode for selected volatile organic compounds in accordance with EPA Method TO-15 from the Compendium of Methods for the Determination of Toxic Organic Compounds in Ambient Air, Second Edition (EPA/625/R-96/010b), January, 1999. The analytical system was comprised of a gas chromatograph / mass spectrometer (GC/MS) interfaced to a whole-air preconcentrator.

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*The results of analyses are given in the attached laboratory report. All results are intended to be considered in their entirety, and Columbia Analytical Services, Inc. (CAS) is not responsible for utilization of less than the complete report.*

*Use of Columbia Analytical Services, Inc. (CAS) Name. Client shall not use CAS's name or trademark in any marketing or reporting materials, press releases or in any other manner ("Materials") whatsoever and shall not attribute to CAS any test result, tolerance or specification derived from CAS's data ("Attribution") without CAS's prior written consent, which may be withheld by CAS for any reason in its sole discretion. To request CAS's consent, Client shall provide copies of the proposed Materials or Attribution and describe in writing Client's proposed use of such Materials or Attribution. If CAS has not provided written approval of the Materials or Attribution within ten (10) days of receipt from Client, Client's request to use CAS's name or trademark in any Materials or Attribution shall be deemed denied. CAS may, in its discretion, reasonably charge Client for its time in reviewing Materials or Attribution requests. Client acknowledges and agrees that the unauthorized use of CAS's name or trademark may cause CAS to incur irreparable harm for which the recovery of money damages will be inadequate. Accordingly, Client acknowledges and agrees that a violation shall justify preliminary injunctive relief. For questions contact the laboratory.*

DETAIL SUMMARY REPORT

Client: Montana DEQ  
 Project ID: Background Indoor Air Study

Service Request: P1201319

Date Received: 4/4/2012  
 Time Received: 09:50

Client Sample ID	Lab Code	Matrix	Date Collected	Time Collected	Container ID	Pi1 (psig)	Pf1 (psig)	2nd Pi (psig)	2nd Pf (psig)	MA APH 1.0 - MA VOC PH Can	TO-15 - VOC Cans	TO-15 - VOC SIM
50-2F-032512	P1201319-001	Air	3/25/2012	09:08	AC01225	-5.46	3.75	-3.36	2.06	X	X	X
49-1F-032712	P1201319-002	Air	3/27/2012	07:30	AC01223	-2.77	3.75			X	X	X
38-1F-032812	P1201319-003	Air	3/28/2012	07:20	AC00993	-3.55	3.63			X	X	X
39-1F-032812	P1201319-004	Air	3/28/2012	09:03	AC01628	-2.36	3.62			X	X	X
40-1F-032812	P1201319-005	Air	3/28/2012	10:11	AC00588	-4.22	3.74			X	X	X
55-1F-032812	P1201319-006	Air	3/28/2012	10:15	AS00131	-3.45	3.63			X	X	X
41-1F-032912	P1201319-007	Air	3/29/2012	11:13	AC01819	-3.37	3.65			X	X	X
48-1F-032912	P1201319-008	Air	3/29/2012	11:51	AC00142	-5.15	3.70			X	X	X
43-1F-032912	P1201319-009	Air	3/29/2012	14:05	AC00603	-4.54	3.74			X	X	X
45-1F-032912	P1201319-010	Air	3/29/2012	14:55	AC01626	-1.65	3.69			X	X	X
42-1F-032912	P1201319-011	Air	3/29/2012	16:09	AS00144	-2.74	3.58			X	X	X
44-1F-032912	P1201319-012	Air	3/29/2012	16:57	AC01183	-3.92	3.76			X	X	X
46-1F-033012	P1201319-013	Air	3/30/2012	18:00	AC00810	-4.91	3.67			X	X	X
47-1F-033112	P1201319-014	Air	3/31/2012	08:17	AC01122	-5.26	3.76			X	X	X



**Air - Chain of Custody Record & Analytical Service Request**

Requested Turnaround Time in Business Days (Surcharges) please circle  
 1 Day (100%) 2 Day (75%) 3 Day (50%) 4 Day (35%) 5 Day (25%) 10 Day-Standard

CAS Project No. **P1201319**

Company Name & Address (Reporting Information)  
 Montana Department of Environmental Quality  
 PO Box 200901  
 100 N. East Chance Gulch  
 Helena, MT 59620-0901

Project Name: **Background Indoor Air Study**  
 Project Number: **---**  
 P.O. # / Billing Information  
 Reference as: Background Indoor Air Study  
 Bill to: Chris Cote - MDEQ

CAS Contact: **Nicole Rannone**  
 Analysis Method: **MA-APH**

Comments  
 e.g. Actual Preservative or specific instructions

Project Manager: **Chris Cote**  
 Phone: (406) 841-5078 Fax: (406) 841-5050

Sampler (Print & Sign): **Chris Cote, Nick Sauer, Steve Dpp**

Client Sample ID	Laboratory ID Number	Date Collected	Time Collected	Canister ID (Bar code # - AC, SC, etc.)	Flow Controller ID (Bar code # - FC #)	Canister Start Pressure "Hg	Canister End Pressure "Hg/psig	Sample Volume	CAS Contact	CAS Project No.	Comments
50-2F-032512	0554	03-25-12	0908	AC01225	FCAD0151	-27.75"	-10.0	17.75	X	X	Analyze
49-1F-032712	0225	03-27-12	0730	AC01223	FCAD0200	-27.5	-6.0	21.5	X	X	Project
38-1F-032812	0336	03-28-12	0720	AC00993	FCAD0518	-26.0	-6.0	20.0	X	X	Specific
39-1F-032812	0245	03-28-12	0903	AC01628	FCAD0053	-27.0	-4.0	23.0	X	X	List of
40-1F-032812	0431	03-28-12	1011	AC00588	FCAD0093	-24.5	-4.5	20.0	X	X	Compounds
55-1F-032812	0352	03-28-12	1015	AS00131	FCAD0468	-27.5	-6.5	21.0	X	X	to Project
41-1F-032912	0348	03-29-12	1113	AC01819	FCAD0471	-25.0	-4.0	21.0	X	X	Specific
48-1F-032912	0524	03-29-12	11:51	AC00142	FCAD0487	-28.5	-10.0	18.5	X	X	Reporting
43-1F-032912	0446	03-29-12	1405	AC00603	FCAD0476	-27.0	-9.0	18.0	X	X	limits
45-1F-032912	0175	03-29-12	1455	AC01626	FCAD0192	-26.0	-1.0	25.0	X	X	
42-1F-032912	0282	03-29-12	1609	AS00144	FCAD0086	-29.0	-6.0	23.0	X	X	Digital Pressure
44-1F-032912	0400	03-29-12	1657	AC01183	FCAD0378	-28.5	-8.0	20.5	X	X	gauge included
46-1F-033012	0501	03-30-12	1800	AC00810	FCAD00383	-27.0	-9.0	18.0	X	X	in shipment
47-1F-033112	0597	03-31-12	0817	AC01122	FCAD00214	-28.5	-10.5	18.0	X	X	

Report Tier Levels - please select  
 Tier I - Results (Default if not specified)  Tier III (Results + QC & Calibration Summaries)  
 Tier II (Results + QC Summaries)  Tier IV (Data Validation Package) 10% Surcharge  Tier IV (Data Validation Package) 10% Surcharge

Relinquished by: (Signature) *Chris Cote* Date: 04-03-12 Time: 15:00  
 Received by: (Signature) *[Signature]* Date: 4/12/12 Time: 09:50  
 Cooler / Blank Temperature: \_\_\_\_\_ °C

**Sample Acceptance Check Form**

Client: Montana DEQ Work order: P1201319

Project: Background Indoor Air Study

Sample(s) received on: 4/4/12 Date opened: 4/4/12 by: MZAMORA

**Note:** This form is used for all samples received by CAS. The use of this form for custody seals is strictly meant to indicate presence/absence and not as an indication of compliance or nonconformity. Thermal preservation and pH will only be evaluated either at the request of the client and/or as required by the method/SOP.

- |    |   | Yes                                 | No                                  | N/A                                 |
|----|---|-------------------------------------|-------------------------------------|-------------------------------------|
| 1  | Were <b>sample containers</b> properly marked with client sample ID?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 2  | Container(s) <b>supplied by CAS</b> ?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 3  | Did <b>sample containers</b> arrive in good condition?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 4  | Were <b>chain-of-custody</b> papers used and filled out?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 5  | Did <b>sample container labels</b> and/or tags agree with custody papers?                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 6  | Was <b>sample volume</b> received adequate for analysis?  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 7  | Are samples within specified holding times?   | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 8  | Was proper <b>temperature</b> (thermal preservation) of cooler at receipt adhered to?                         | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 9  | Was a <b>trip blank</b> received?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| 10 | Were <b>custody seals</b> on outside of cooler/Box?   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
|    | Location of seal(s)? _____ Sealing Lid?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Were signature and date included?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Were seals intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Were custody seals on outside of sample container?  | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
|    | Location of seal(s)? _____ Sealing Lid?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Were signature and date included?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Were seals intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 11 | Do containers have appropriate <b>preservation</b> , according to method/SOP or Client specified information? | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Is there a client indication that the submitted samples are <b>pH</b> preserved?                              | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Were <b>VOA vials</b> checked for presence/absence of air bubbles?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Does the client/method/SOP require that the analyst check the sample pH and <u>if necessary</u> alter it?     | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 12 | <b>Tubes:</b> Are the tubes capped and intact?  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Do they contain moisture?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 13 | <b>Badges:</b> Are the badges properly capped and intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
|    | Are dual bed badges separated and individually capped and intact?   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |

Lab Sample ID	Container Description	Required pH *	Received pH	Adjusted pH	VOA Headspace (Presence/Absence)	Receipt / Preservation Comments
P1201319-001.01	6.0 L Ambient Can					
P1201319-002.01	6.0 L Ambient Can					
P1201319-003.01	6.0 L Ambient Can					
P1201319-004.01	6.0 L Ambient Can					
P1201319-005.01	6.0 L Ambient Can					
P1201319-006.01	6.0 L Silonite Can					
P1201319-007.01	6.0 L Ambient Can					
P1201319-008.01	6.0 L Ambient Can					

Explain any discrepancies: (include lab sample ID numbers): \_\_\_\_\_



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 50-2F-032512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-001

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01225

Date Collected: 3/25/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.46      Final Pressure (psig): 3.75

Canister Dilution Factor: 2.00

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	460	40	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	590	20	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	47	10	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 49-1F-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-002

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01223

Date Collected: 3/27/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.77      Final Pressure (psig): 3.75

Canister Dilution Factor: 1.55

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	120	31	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	200	16	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	8.6	7.8	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 38-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-003

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00993

Date Collected: 3/28/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.55      Final Pressure (psig): 3.63

Canister Dilution Factor: 1.64

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	50	33	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	100	16	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	8.2	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 39-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-004

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01628

Date Collected: 3/28/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.36      Final Pressure (psig): 3.62

Canister Dilution Factor: 1.48

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	110	30	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	29	15	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	7.4	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 40-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-005

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00588

Date Collected: 3/28/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.22      Final Pressure (psig): 3.74

Canister Dilution Factor: 1.76

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	35	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	<b>30</b>	18	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	8.8	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



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**Client:** Montana DEQ  
**Client Sample ID:** 55-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-006

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00131

Date Collected: 3/28/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.45      Final Pressure (psig): 3.63

Canister Dilution Factor: 1.63

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	33	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	19	16	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	8.2	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 41-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-007

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01819

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.37      Final Pressure (psig): 3.65

Canister Dilution Factor: 1.62

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	120	32	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	69	16	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	12	8.1	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 48-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-008

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00142

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.15      Final Pressure (psig): 3.70

Canister Dilution Factor: 1.93

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	170	39	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	81	19	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	22	9.7	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 43-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-009

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00603

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.54      Final Pressure (psig): 3.74

Canister Dilution Factor: 1.81

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	90	36	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	79	18	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	9.1	9.1	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 45-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-010

Test Code: Massachusetts APH, Revision 1, December 2009  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01626

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.65      Final Pressure (psig): 3.69

Canister Dilution Factor: 1.41

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	79	28	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	56	14	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	9.5	7.1	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 42-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-011

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00144

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.74      Final Pressure (psig): 3.58

Canister Dilution Factor: 1.53

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	31	31	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	26	15	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	7.7	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 44-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-012

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01183

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.92      Final Pressure (psig): 3.76

Canister Dilution Factor: 1.71

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	34	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	24	17	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	8.6	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 46-1F-033012  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-013

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00810

Date Collected: 3/30/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.91      Final Pressure (psig): 3.67

Canister Dilution Factor: 1.88

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	200	38	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	32	19	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	9.4	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 47-1F-033112  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-014

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01122

Date Collected: 3/31/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.26      Final Pressure (psig): 3.76

Canister Dilution Factor: 1.96

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	66	39	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	43	20	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	9.8	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P120407-MB

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/07/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	20	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	10	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	5.0	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P120409-MB

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/09/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	20	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	10	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	5.0	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P120407-LCS

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/07/12  
 Volume(s) Analyzed: NA Liter(s)

Compound	Spike Amount ng	Result ng	% Recovery	CAS	Data Qualifier
				Acceptance Limits	
C5 - C8 Aliphatic Hydrocarbons	24.5	19.7	80	70-130	
C9 - C12 Aliphatic Hydrocarbons	25.0	22.7	91	70-130	
C9 - C10 Aromatic Hydrocarbons	48.5	43.2	89	70-130	

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P120409-LCS

Test Code: Massachusetts APH, Revision 1, December 2009  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/09/12  
Volume(s) Analyzed: NA Liter(s)

Compound	Spike Amount ng	Result ng	% Recovery	CAS	Data Qualifier
				Acceptance Limits	
C5 - C8 Aliphatic Hydrocarbons	24.5	22.3	91	70-130	
C9 - C12 Aliphatic Hydrocarbons	25.0	26.0	104	70-130	
C9 - C10 Aromatic Hydrocarbons	48.5	48.8	101	70-130	

LABORATORY DUPLICATE SUMMARY RESULTS

Page 1 of 1

**Client:** Montana DEQ

**Client Sample ID:** 47-1F-033112

**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319

CAS Sample ID: P1201319-014DUP

Test Code: Massachusetts APH, Revision 1, December 2009

Date Collected: 3/31/12

Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Received: 4/4/12

Analyst: Lusine Hakobyan

Date Analyzed: 4/9/12

Sampling Media: 6.0 L Summa Canister

Volume(s) Analyzed: 1.00 Liter(s)

Test Notes:

Container ID: AC01122

Initial Pressure (psig): -5.26 Final Pressure (psig): 3.76

Canister Dilution Factor: 1.96

Compound	Sample Result	Duplicate Sample Result	Average	% RPD	RPD	Data
	µg/m <sup>3</sup>	µg/m <sup>3</sup>	µg/m <sup>3</sup>		Limit	Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	66.1	63.7	64.9	4	30	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	42.9	37.9	40.4	12	30	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	ND	-	-	30	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 50-2F-032512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-001

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01225

Date Collected: 3/25/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12 & 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.10 Liter(s)

Initial Pressure (psig): -5.46 Final Pressure (psig): 3.75

Canister Dilution Factor: 2.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	21	1.0	12	0.58	
75-71-8	Dichlorodifluoromethane (CFC 12)	1.8	1.0	0.36	0.20	
74-87-3	Chloromethane	ND	1.0	ND	0.48	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	1.0	ND	0.14	
106-99-0	1,3-Butadiene	ND	1.0	ND	0.45	
74-83-9	Bromomethane	ND	1.0	ND	0.26	
75-00-3	Chloroethane	ND	1.0	ND	0.38	
64-17-5	Ethanol	1,500	100	810	53	D
75-05-8	Acetonitrile	ND	1.0	ND	0.60	
107-02-8	Acrolein	ND	4.0	ND	1.7	
67-64-1	Acetone	66	10	28	4.2	
75-69-4	Trichlorofluoromethane	ND	1.0	ND	0.18	
67-63-0	2-Propanol (Isopropyl Alcohol)	98	2.0	40	0.81	
107-13-1	Acrylonitrile	ND	1.0	ND	0.46	
75-35-4	1,1-Dichloroethene	ND	1.0	ND	0.25	
75-09-2	Methylene Chloride	ND	1.0	ND	0.29	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	1.0	ND	0.32	
76-13-1	Trichlorotrifluoroethane	ND	1.0	ND	0.13	
75-15-0	Carbon Disulfide	ND	10	ND	3.2	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	ND	0.25	
75-34-3	1,1-Dichloroethane	ND	1.0	ND	0.25	
1634-04-4	Methyl tert-Butyl Ether	ND	1.0	ND	0.28	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 50-2F-032512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-001

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01225

Date Collected: 3/25/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12 & 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.10 Liter(s)

Initial Pressure (psig): -5.46 Final Pressure (psig): 3.75

Canister Dilution Factor: 2.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	10	ND	2.8	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	ND	0.25	
78-93-3	2-Butanone (MEK)	11	10	3.6	3.4	
141-78-6	Ethyl Acetate	4.6	2.0	1.3	0.56	
110-54-3	n-Hexane	19	1.0	5.5	0.28	
67-66-3	Chloroform	1.1	1.0	0.22	0.20	
109-99-9	Tetrahydrofuran (THF)	3.6	1.0	1.2	0.34	
71-55-6	1,1,1-Trichloroethane	ND	1.0	ND	0.18	
56-23-5	Carbon Tetrachloride	ND	1.0	ND	0.16	
110-82-7	Cyclohexane	3.9	2.0	1.1	0.58	
78-87-5	1,2-Dichloropropane	ND	1.0	ND	0.22	
75-27-4	Bromodichloromethane	ND	1.0	ND	0.15	
123-91-1	1,4-Dioxane	ND	1.0	ND	0.28	
80-62-6	Methyl Methacrylate	ND	2.0	ND	0.49	
142-82-5	n-Heptane	7.6	1.0	1.9	0.24	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	ND	0.22	
108-10-1	4-Methyl-2-pentanone	1.0	1.0	0.25	0.24	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	ND	0.22	
108-88-3	Toluene	51	1.0	14	0.27	
591-78-6	2-Hexanone	1.4	1.0	0.33	0.24	
124-48-1	Dibromochloromethane	ND	1.0	ND	0.12	
106-93-4	1,2-Dibromoethane	ND	1.0	ND	0.13	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 50-2F-032512  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201319  
**CAS Sample ID:** P1201319-001

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01225

**Date Collected:** 3/25/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/7/12 & 4/9/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -5.46 Final Pressure (psig): 3.75

Canister Dilution Factor: 2.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	1.3	1.0	0.28	0.21	
111-65-9	n-Octane	2.3	1.0	0.50	0.21	
108-90-7	Chlorobenzene	ND	1.0	ND	0.22	
179601-23-1	m,p-Xylenes	29	2.0	6.6	0.46	
75-25-2	Bromoform	ND	1.0	ND	0.097	
100-42-5	Styrene	ND	1.0	ND	0.23	
95-47-6	o-Xylene	8.8	1.0	2.0	0.23	
111-84-2	n-Nonane	39	1.0	7.4	0.19	
98-82-8	Cumene	ND	1.0	ND	0.20	
80-56-8	alpha-Pinene	30	1.0	5.4	0.18	
103-65-1	n-Propylbenzene	1.5	1.0	0.31	0.20	
622-96-8	4-Ethyltoluene	3.0	1.0	0.61	0.20	
108-67-8	1,3,5-Trimethylbenzene	2.8	1.0	0.58	0.20	
95-63-6	1,2,4-Trimethylbenzene	10	1.0	2.0	0.20	
100-44-7	Benzyl Chloride	ND	1.0	ND	0.19	
541-73-1	1,3-Dichlorobenzene	ND	1.0	ND	0.17	
106-46-7	1,4-Dichlorobenzene	ND	1.0	ND	0.17	
95-50-1	1,2-Dichlorobenzene	ND	1.0	ND	0.17	
5989-27-5	d-Limonene	11	1.0	2.0	0.18	
96-12-8	1,2-Dibromo-3-chloropropane	ND	1.0	ND	0.10	
120-82-1	1,2,4-Trichlorobenzene	ND	1.0	ND	0.13	
87-68-3	Hexachlorobutadiene	ND	1.0	ND	0.094	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 49-1F-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-002

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01223

Date Collected: 3/27/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.77 Final Pressure (psig): 3.75

Canister Dilution Factor: 1.55

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	6.2	0.78	3.6	0.45	
75-71-8	Dichlorodifluoromethane (CFC 12)	1.8	0.78	0.37	0.16	
74-87-3	Chloromethane	ND	0.78	ND	0.38	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.78	ND	0.11	
106-99-0	1,3-Butadiene	ND	0.78	ND	0.35	
74-83-9	Bromomethane	ND	0.78	ND	0.20	
75-00-3	Chloroethane	ND	0.78	ND	0.29	
64-17-5	Ethanol	580	7.8	310	4.1	
75-05-8	Acetonitrile	3.2	0.78	1.9	0.46	
107-02-8	Acrolein	ND	3.1	ND	1.4	
67-64-1	Acetone	52	7.8	22	3.3	
75-69-4	Trichlorofluoromethane	1.0	0.78	0.18	0.14	
67-63-0	2-Propanol (Isopropyl Alcohol)	32	1.6	13	0.63	
107-13-1	Acrylonitrile	ND	0.78	ND	0.36	
75-35-4	1,1-Dichloroethene	ND	0.78	ND	0.20	
75-09-2	Methylene Chloride	ND	0.78	ND	0.22	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.78	ND	0.25	
76-13-1	Trichlorotrifluoroethane	ND	0.78	ND	0.10	
75-15-0	Carbon Disulfide	ND	7.8	ND	2.5	
156-60-5	trans-1,2-Dichloroethene	ND	0.78	ND	0.20	
75-34-3	1,1-Dichloroethane	ND	0.78	ND	0.19	
1634-04-4	Methyl tert-Butyl Ether	ND	0.78	ND	0.22	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 49-1F-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-002

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01223

Date Collected: 3/27/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.77 Final Pressure (psig): 3.75

Canister Dilution Factor: 1.55

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	7.8	ND	2.2	
156-59-2	cis-1,2-Dichloroethene	ND	0.78	ND	0.20	
78-93-3	2-Butanone (MEK)	<b>11</b>	7.8	<b>3.7</b>	2.6	
141-78-6	Ethyl Acetate	<b>25</b>	1.6	<b>6.9</b>	0.43	
110-54-3	n-Hexane	<b>2.2</b>	0.78	<b>0.63</b>	0.22	
67-66-3	Chloroform	ND	0.78	ND	0.16	
109-99-9	Tetrahydrofuran (THF)	<b>4.8</b>	0.78	<b>1.6</b>	0.26	
71-55-6	1,1,1-Trichloroethane	ND	0.78	ND	0.14	
56-23-5	Carbon Tetrachloride	ND	0.78	ND	0.12	
110-82-7	Cyclohexane	ND	1.6	ND	0.45	
78-87-5	1,2-Dichloropropane	ND	0.78	ND	0.17	
75-27-4	Bromodichloromethane	ND	0.78	ND	0.12	
123-91-1	1,4-Dioxane	ND	0.78	ND	0.22	
80-62-6	Methyl Methacrylate	ND	1.6	ND	0.38	
142-82-5	n-Heptane	<b>1.1</b>	0.78	<b>0.27</b>	0.19	
10061-01-5	cis-1,3-Dichloropropene	ND	0.78	ND	0.17	
108-10-1	4-Methyl-2-pentanone	ND	0.78	ND	0.19	
10061-02-6	trans-1,3-Dichloropropene	ND	0.78	ND	0.17	
108-88-3	Toluene	<b>8.1</b>	0.78	<b>2.2</b>	0.21	
591-78-6	2-Hexanone	<b>0.78</b>	0.78	<b>0.19</b>	0.19	
124-48-1	Dibromochloromethane	ND	0.78	ND	0.091	
106-93-4	1,2-Dibromoethane	ND	0.78	ND	0.10	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 49-1F-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-002

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01223

Date Collected: 3/27/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.77 Final Pressure (psig): 3.75

Canister Dilution Factor: 1.55

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	5.4	0.78	1.1	0.16	
111-65-9	n-Octane	0.82	0.78	0.18	0.17	
108-90-7	Chlorobenzene	ND	0.78	ND	0.17	
179601-23-1	m,p-Xylenes	4.0	1.6	0.91	0.36	
75-25-2	Bromoform	ND	0.78	ND	0.075	
100-42-5	Styrene	1.3	0.78	0.31	0.18	
95-47-6	o-Xylene	1.2	0.78	0.27	0.18	
111-84-2	n-Nonane	2.5	0.78	0.48	0.15	
98-82-8	Cumene	ND	0.78	ND	0.16	
80-56-8	alpha-Pinene	5.1	0.78	0.92	0.14	
103-65-1	n-Propylbenzene	ND	0.78	ND	0.16	
622-96-8	4-Ethyltoluene	ND	0.78	ND	0.16	
108-67-8	1,3,5-Trimethylbenzene	ND	0.78	ND	0.16	
95-63-6	1,2,4-Trimethylbenzene	1.2	0.78	0.25	0.16	
100-44-7	Benzyl Chloride	ND	0.78	ND	0.15	
541-73-1	1,3-Dichlorobenzene	ND	0.78	ND	0.13	
106-46-7	1,4-Dichlorobenzene	ND	0.78	ND	0.13	
95-50-1	1,2-Dichlorobenzene	ND	0.78	ND	0.13	
5989-27-5	d-Limonene	16	0.78	2.8	0.14	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.78	ND	0.080	
120-82-1	1,2,4-Trichlorobenzene	ND	0.78	ND	0.10	
87-68-3	Hexachlorobutadiene	ND	0.78	ND	0.073	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 38-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-003

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC00993

**Date Collected:** 3/28/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/7/12 & 4/9/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -3.55      Final Pressure (psig): 3.63

Canister Dilution Factor: 1.64

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	3.2	0.82	1.9	0.48	
75-71-8	Dichlorodifluoromethane (CFC 12)	1.8	0.82	0.37	0.17	
74-87-3	Chloromethane	ND	0.82	ND	0.40	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.82	ND	0.12	
106-99-0	1,3-Butadiene	ND	0.82	ND	0.37	
74-83-9	Bromomethane	ND	0.82	ND	0.21	
75-00-3	Chloroethane	ND	0.82	ND	0.31	
64-17-5	Ethanol	1,100	82	580	44	D
75-05-8	Acetonitrile	ND	0.82	ND	0.49	
107-02-8	Acrolein	ND	3.3	ND	1.4	
67-64-1	Acetone	45	8.2	19	3.5	
75-69-4	Trichlorofluoromethane	1.3	0.82	0.23	0.15	
67-63-0	2-Propanol (Isopropyl Alcohol)	30	1.6	12	0.67	
107-13-1	Acrylonitrile	ND	0.82	ND	0.38	
75-35-4	1,1-Dichloroethene	ND	0.82	ND	0.21	
75-09-2	Methylene Chloride	ND	0.82	ND	0.24	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.82	ND	0.26	
76-13-1	Trichlorotrifluoroethane	ND	0.82	ND	0.11	
75-15-0	Carbon Disulfide	ND	8.2	ND	2.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.82	ND	0.21	
75-34-3	1,1-Dichloroethane	ND	0.82	ND	0.20	
1634-04-4	Methyl tert-Butyl Ether	ND	0.82	ND	0.23	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 38-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-003

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00993

Date Collected: 3/28/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12 & 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.10 Liter(s)

Initial Pressure (psig): -3.55 Final Pressure (psig): 3.63

Canister Dilution Factor: 1.64

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	8.2	ND	2.3	
156-59-2	cis-1,2-Dichloroethene	ND	0.82	ND	0.21	
78-93-3	2-Butanone (MEK)	16	8.2	5.3	2.8	
141-78-6	Ethyl Acetate	6.0	1.6	1.7	0.46	
110-54-3	n-Hexane	ND	0.82	ND	0.23	
67-66-3	Chloroform	2.1	0.82	0.44	0.17	
109-99-9	Tetrahydrofuran (THF)	4.5	0.82	1.5	0.28	
71-55-6	1,1,1-Trichloroethane	ND	0.82	ND	0.15	
56-23-5	Carbon Tetrachloride	ND	0.82	ND	0.13	
110-82-7	Cyclohexane	ND	1.6	ND	0.48	
78-87-5	1,2-Dichloropropane	ND	0.82	ND	0.18	
75-27-4	Bromodichloromethane	ND	0.82	ND	0.12	
123-91-1	1,4-Dioxane	ND	0.82	ND	0.23	
80-62-6	Methyl Methacrylate	ND	1.6	ND	0.40	
142-82-5	n-Heptane	ND	0.82	ND	0.20	
10061-01-5	cis-1,3-Dichloropropene	ND	0.82	ND	0.18	
108-10-1	4-Methyl-2-pentanone	2.3	0.82	0.56	0.20	
10061-02-6	trans-1,3-Dichloropropene	ND	0.82	ND	0.18	
108-88-3	Toluene	3.2	0.82	0.86	0.22	
591-78-6	2-Hexanone	ND	0.82	ND	0.20	
124-48-1	Dibromochloromethane	ND	0.82	ND	0.096	
106-93-4	1,2-Dibromoethane	ND	0.82	ND	0.11	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 38-1F-032812  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201319  
**CAS Sample ID:** P1201319-003

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC00993

**Date Collected:** 3/28/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/7/12 & 4/9/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -3.55 Final Pressure (psig): 3.63

Canister Dilution Factor: 1.64

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	1.3	0.82	0.27	0.17	
111-65-9	n-Octane	ND	0.82	ND	0.18	
108-90-7	Chlorobenzene	ND	0.82	ND	0.18	
179601-23-1	m,p-Xylenes	ND	1.6	ND	0.38	
75-25-2	Bromoform	ND	0.82	ND	0.079	
100-42-5	Styrene	ND	0.82	ND	0.19	
95-47-6	o-Xylene	ND	0.82	ND	0.19	
111-84-2	n-Nonane	ND	0.82	ND	0.16	
98-82-8	Cumene	ND	0.82	ND	0.17	
80-56-8	alpha-Pinene	5.2	0.82	0.93	0.15	
103-65-1	n-Propylbenzene	ND	0.82	ND	0.17	
622-96-8	4-Ethyltoluene	ND	0.82	ND	0.17	
108-67-8	1,3,5-Trimethylbenzene	ND	0.82	ND	0.17	
95-63-6	1,2,4-Trimethylbenzene	ND	0.82	ND	0.17	
100-44-7	Benzyl Chloride	ND	0.82	ND	0.16	
541-73-1	1,3-Dichlorobenzene	ND	0.82	ND	0.14	
106-46-7	1,4-Dichlorobenzene	27	0.82	4.5	0.14	
95-50-1	1,2-Dichlorobenzene	ND	0.82	ND	0.14	
5989-27-5	d-Limonene	13	0.82	2.4	0.15	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.82	ND	0.085	
120-82-1	1,2,4-Trichlorobenzene	ND	0.82	ND	0.11	
87-68-3	Hexachlorobutadiene	ND	0.82	ND	0.077	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 39-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-004

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01628

Date Collected: 3/28/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12 & 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.10 Liter(s)

Initial Pressure (psig): -2.36 Final Pressure (psig): 3.62

Canister Dilution Factor: 1.48

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	6.4	0.74	3.7	0.43	
75-71-8	Dichlorodifluoromethane (CFC 12)	1.8	0.74	0.36	0.15	
74-87-3	Chloromethane	ND	0.74	ND	0.36	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.74	ND	0.11	
106-99-0	1,3-Butadiene	ND	0.74	ND	0.33	
74-83-9	Bromomethane	ND	0.74	ND	0.19	
75-00-3	Chloroethane	ND	0.74	ND	0.28	
64-17-5	Ethanol	1,900	74	1,000	39	D
75-05-8	Acetonitrile	ND	0.74	ND	0.44	
107-02-8	Acrolein	ND	3.0	ND	1.3	
67-64-1	Acetone	60	7.4	25	3.1	
75-69-4	Trichlorofluoromethane	0.89	0.74	0.16	0.13	
67-63-0	2-Propanol (Isopropyl Alcohol)	46	1.5	19	0.60	
107-13-1	Acrylonitrile	ND	0.74	ND	0.34	
75-35-4	1,1-Dichloroethene	ND	0.74	ND	0.19	
75-09-2	Methylene Chloride	ND	0.74	ND	0.21	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.74	ND	0.24	
76-13-1	Trichlorotrifluoroethane	ND	0.74	ND	0.097	
75-15-0	Carbon Disulfide	ND	7.4	ND	2.4	
156-60-5	trans-1,2-Dichloroethene	ND	0.74	ND	0.19	
75-34-3	1,1-Dichloroethane	ND	0.74	ND	0.18	
1634-04-4	Methyl tert-Butyl Ether	ND	0.74	ND	0.21	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.



## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 39-1F-032812  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201319  
**CAS Sample ID:** P1201319-004

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01628

**Date Collected:** 3/28/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/7/12 & 4/9/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -2.36 Final Pressure (psig): 3.62

Canister Dilution Factor: 1.48

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	14	7.4	4.1	2.1	
156-59-2	cis-1,2-Dichloroethene	ND	0.74	ND	0.19	
78-93-3	2-Butanone (MEK)	ND	7.4	ND	2.5	
141-78-6	Ethyl Acetate	22	1.5	6.0	0.41	
110-54-3	n-Hexane	8.3	0.74	2.3	0.21	
67-66-3	Chloroform	ND	0.74	ND	0.15	
109-99-9	Tetrahydrofuran (THF)	ND	0.74	ND	0.25	
71-55-6	1,1,1-Trichloroethane	ND	0.74	ND	0.14	
56-23-5	Carbon Tetrachloride	ND	0.74	ND	0.12	
110-82-7	Cyclohexane	1.9	1.5	0.55	0.43	
78-87-5	1,2-Dichloropropane	ND	0.74	ND	0.16	
75-27-4	Bromodichloromethane	ND	0.74	ND	0.11	
123-91-1	1,4-Dioxane	ND	0.74	ND	0.21	
80-62-6	Methyl Methacrylate	ND	1.5	ND	0.36	
142-82-5	n-Heptane	3.4	0.74	0.84	0.18	
10061-01-5	cis-1,3-Dichloropropene	ND	0.74	ND	0.16	
108-10-1	4-Methyl-2-pentanone	ND	0.74	ND	0.18	
10061-02-6	trans-1,3-Dichloropropene	ND	0.74	ND	0.16	
108-88-3	Toluene	5.7	0.74	1.5	0.20	
591-78-6	2-Hexanone	ND	0.74	ND	0.18	
124-48-1	Dibromochloromethane	ND	0.74	ND	0.087	
106-93-4	1,2-Dibromoethane	ND	0.74	ND	0.096	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 39-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-004

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01628

Date Collected: 3/28/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12 & 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.10 Liter(s)

Initial Pressure (psig): -2.36 Final Pressure (psig): 3.62

Canister Dilution Factor: 1.48

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	2.4	0.74	0.51	0.16	
111-65-9	n-Octane	1.2	0.74	0.26	0.16	
108-90-7	Chlorobenzene	ND	0.74	ND	0.16	
179601-23-1	m,p-Xylenes	ND	1.5	ND	0.34	
75-25-2	Bromoform	ND	0.74	ND	0.072	
100-42-5	Styrene	ND	0.74	ND	0.17	
95-47-6	o-Xylene	ND	0.74	ND	0.17	
111-84-2	n-Nonane	1.4	0.74	0.27	0.14	
98-82-8	Cumene	ND	0.74	ND	0.15	
80-56-8	alpha-Pinene	0.83	0.74	0.15	0.13	
103-65-1	n-Propylbenzene	ND	0.74	ND	0.15	
622-96-8	4-Ethyltoluene	ND	0.74	ND	0.15	
108-67-8	1,3,5-Trimethylbenzene	ND	0.74	ND	0.15	
95-63-6	1,2,4-Trimethylbenzene	ND	0.74	ND	0.15	
100-44-7	Benzyl Chloride	ND	0.74	ND	0.14	
541-73-1	1,3-Dichlorobenzene	ND	0.74	ND	0.12	
106-46-7	1,4-Dichlorobenzene	ND	0.74	ND	0.12	
95-50-1	1,2-Dichlorobenzene	ND	0.74	ND	0.12	
5989-27-5	d-Limonene	13	0.74	2.4	0.13	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.74	ND	0.077	
120-82-1	1,2,4-Trichlorobenzene	ND	0.74	ND	0.10	
87-68-3	Hexachlorobutadiene	ND	0.74	ND	0.069	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 40-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-005

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00588

Date Collected: 3/28/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.22      Final Pressure (psig): 3.74

Canister Dilution Factor: 1.76

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	3.0	0.88	1.8	0.51	
75-71-8	Dichlorodifluoromethane (CFC 12)	1.9	0.88	0.38	0.18	
74-87-3	Chloromethane	ND	0.88	ND	0.43	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.88	ND	0.13	
106-99-0	1,3-Butadiene	ND	0.88	ND	0.40	
74-83-9	Bromomethane	ND	0.88	ND	0.23	
75-00-3	Chloroethane	ND	0.88	ND	0.33	
64-17-5	Ethanol	180	8.8	94	4.7	
75-05-8	Acetonitrile	ND	0.88	ND	0.52	
107-02-8	Acrolein	ND	3.5	ND	1.5	
67-64-1	Acetone	41	8.8	17	3.7	
75-69-4	Trichlorofluoromethane	1.9	0.88	0.33	0.16	
67-63-0	2-Propanol (Isopropyl Alcohol)	15	1.8	6.2	0.72	
107-13-1	Acrylonitrile	ND	0.88	ND	0.41	
75-35-4	1,1-Dichloroethene	ND	0.88	ND	0.22	
75-09-2	Methylene Chloride	0.90	0.88	0.26	0.25	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.88	ND	0.28	
76-13-1	Trichlorotrifluoroethane	ND	0.88	ND	0.11	
75-15-0	Carbon Disulfide	ND	8.8	ND	2.8	
156-60-5	trans-1,2-Dichloroethene	ND	0.88	ND	0.22	
75-34-3	1,1-Dichloroethane	ND	0.88	ND	0.22	
1634-04-4	Methyl tert-Butyl Ether	ND	0.88	ND	0.24	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 40-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-005

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00588

Date Collected: 3/28/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.22 Final Pressure (psig): 3.74

Canister Dilution Factor: 1.76

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	8.8	ND	2.5	
156-59-2	cis-1,2-Dichloroethene	ND	0.88	ND	0.22	
78-93-3	2-Butanone (MEK)	ND	8.8	ND	3.0	
141-78-6	Ethyl Acetate	ND	1.8	ND	0.49	
110-54-3	n-Hexane	ND	0.88	ND	0.25	
67-66-3	Chloroform	ND	0.88	ND	0.18	
109-99-9	Tetrahydrofuran (THF)	ND	0.88	ND	0.30	
71-55-6	1,1,1-Trichloroethane	ND	0.88	ND	0.16	
56-23-5	Carbon Tetrachloride	ND	0.88	ND	0.14	
110-82-7	Cyclohexane	ND	1.8	ND	0.51	
78-87-5	1,2-Dichloropropane	ND	0.88	ND	0.19	
75-27-4	Bromodichloromethane	ND	0.88	ND	0.13	
123-91-1	1,4-Dioxane	ND	0.88	ND	0.24	
80-62-6	Methyl Methacrylate	ND	1.8	ND	0.43	
142-82-5	n-Heptane	ND	0.88	ND	0.21	
10061-01-5	cis-1,3-Dichloropropene	ND	0.88	ND	0.19	
108-10-1	4-Methyl-2-pentanone	ND	0.88	ND	0.21	
10061-02-6	trans-1,3-Dichloropropene	ND	0.88	ND	0.19	
108-88-3	Toluene	<b>1.7</b>	0.88	<b>0.44</b>	0.23	
591-78-6	2-Hexanone	ND	0.88	ND	0.21	
124-48-1	Dibromochloromethane	ND	0.88	ND	0.10	
106-93-4	1,2-Dibromoethane	ND	0.88	ND	0.11	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 40-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-005

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00588

Date Collected: 3/28/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.22 Final Pressure (psig): 3.74

Canister Dilution Factor: 1.76

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.88	ND	0.19	
111-65-9	n-Octane	ND	0.88	ND	0.19	
108-90-7	Chlorobenzene	ND	0.88	ND	0.19	
179601-23-1	m,p-Xylenes	ND	1.8	ND	0.41	
75-25-2	Bromoform	ND	0.88	ND	0.085	
100-42-5	Styrene	ND	0.88	ND	0.21	
95-47-6	o-Xylene	ND	0.88	ND	0.20	
111-84-2	n-Nonane	ND	0.88	ND	0.17	
98-82-8	Cumene	ND	0.88	ND	0.18	
80-56-8	alpha-Pinene	23	0.88	4.1	0.16	
103-65-1	n-Propylbenzene	ND	0.88	ND	0.18	
622-96-8	4-Ethyltoluene	ND	0.88	ND	0.18	
108-67-8	1,3,5-Trimethylbenzene	ND	0.88	ND	0.18	
95-63-6	1,2,4-Trimethylbenzene	ND	0.88	ND	0.18	
100-44-7	Benzyl Chloride	ND	0.88	ND	0.17	
541-73-1	1,3-Dichlorobenzene	ND	0.88	ND	0.15	
106-46-7	1,4-Dichlorobenzene	ND	0.88	ND	0.15	
95-50-1	1,2-Dichlorobenzene	ND	0.88	ND	0.15	
5989-27-5	d-Limonene	7.3	0.88	1.3	0.16	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.88	ND	0.091	
120-82-1	1,2,4-Trichlorobenzene	ND	0.88	ND	0.12	
87-68-3	Hexachlorobutadiene	ND	0.88	ND	0.083	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 55-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-006

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00131

Date Collected: 3/28/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.45 Final Pressure (psig): 3.63

Canister Dilution Factor: 1.63

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	<b>3.8</b>	0.82	<b>2.2</b>	0.47	
75-71-8	Dichlorodifluoromethane (CFC 12)	<b>1.9</b>	0.82	<b>0.39</b>	0.16	
74-87-3	Chloromethane	ND	0.82	ND	0.39	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.82	ND	0.12	
106-99-0	1,3-Butadiene	ND	0.82	ND	0.37	
74-83-9	Bromomethane	ND	0.82	ND	0.21	
75-00-3	Chloroethane	ND	0.82	ND	0.31	
64-17-5	Ethanol	<b>180</b>	8.2	<b>98</b>	4.3	
75-05-8	Acetonitrile	ND	0.82	ND	0.49	
107-02-8	Acrolein	ND	3.3	ND	1.4	
67-64-1	Acetone	<b>38</b>	8.2	<b>16</b>	3.4	
75-69-4	Trichlorofluoromethane	<b>1.9</b>	0.82	<b>0.33</b>	0.15	
67-63-0	2-Propanol (Isopropyl Alcohol)	<b>17</b>	1.6	<b>7.0</b>	0.66	
107-13-1	Acrylonitrile	ND	0.82	ND	0.38	
75-35-4	1,1-Dichloroethene	ND	0.82	ND	0.21	
75-09-2	Methylene Chloride	<b>0.86</b>	0.82	<b>0.25</b>	0.23	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.82	ND	0.26	
76-13-1	Trichlorotrifluoroethane	<b>0.82</b>	0.82	<b>0.11</b>	0.11	
75-15-0	Carbon Disulfide	ND	8.2	ND	2.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.82	ND	0.21	
75-34-3	1,1-Dichloroethane	ND	0.82	ND	0.20	
1634-04-4	Methyl tert-Butyl Ether	ND	0.82	ND	0.23	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 55-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-006

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00131

Date Collected: 3/28/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.45 Final Pressure (psig): 3.63

Canister Dilution Factor: 1.63

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	8.2	ND	2.3	
156-59-2	cis-1,2-Dichloroethene	ND	0.82	ND	0.21	
78-93-3	2-Butanone (MEK)	ND	8.2	ND	2.8	
141-78-6	Ethyl Acetate	ND	1.6	ND	0.45	
110-54-3	n-Hexane	ND	0.82	ND	0.23	
67-66-3	Chloroform	ND	0.82	ND	0.17	
109-99-9	Tetrahydrofuran (THF)	ND	0.82	ND	0.28	
71-55-6	1,1,1-Trichloroethane	ND	0.82	ND	0.15	
56-23-5	Carbon Tetrachloride	ND	0.82	ND	0.13	
110-82-7	Cyclohexane	ND	1.6	ND	0.47	
78-87-5	1,2-Dichloropropane	ND	0.82	ND	0.18	
75-27-4	Bromodichloromethane	ND	0.82	ND	0.12	
123-91-1	1,4-Dioxane	ND	0.82	ND	0.23	
80-62-6	Methyl Methacrylate	ND	1.6	ND	0.40	
142-82-5	n-Heptane	ND	0.82	ND	0.20	
10061-01-5	cis-1,3-Dichloropropene	ND	0.82	ND	0.18	
108-10-1	4-Methyl-2-pentanone	ND	0.82	ND	0.20	
10061-02-6	trans-1,3-Dichloropropene	ND	0.82	ND	0.18	
108-88-3	Toluene	<b>1.6</b>	0.82	<b>0.43</b>	0.22	
591-78-6	2-Hexanone	ND	0.82	ND	0.20	
124-48-1	Dibromochloromethane	ND	0.82	ND	0.096	
106-93-4	1,2-Dibromoethane	ND	0.82	ND	0.11	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 55-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-006

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00131

Date Collected: 3/28/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.45 Final Pressure (psig): 3.63

Canister Dilution Factor: 1.63

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	0.83	0.82	0.17	0.17	
111-65-9	n-Octane	ND	0.82	ND	0.17	
108-90-7	Chlorobenzene	ND	0.82	ND	0.18	
179601-23-1	m,p-Xylenes	ND	1.6	ND	0.38	
75-25-2	Bromoform	ND	0.82	ND	0.079	
100-42-5	Styrene	ND	0.82	ND	0.19	
95-47-6	o-Xylene	ND	0.82	ND	0.19	
111-84-2	n-Nonane	ND	0.82	ND	0.16	
98-82-8	Cumene	ND	0.82	ND	0.17	
80-56-8	alpha-Pinene	25	0.82	4.6	0.15	
103-65-1	n-Propylbenzene	ND	0.82	ND	0.17	
622-96-8	4-Ethyltoluene	ND	0.82	ND	0.17	
108-67-8	1,3,5-Trimethylbenzene	ND	0.82	ND	0.17	
95-63-6	1,2,4-Trimethylbenzene	ND	0.82	ND	0.17	
100-44-7	Benzyl Chloride	ND	0.82	ND	0.16	
541-73-1	1,3-Dichlorobenzene	ND	0.82	ND	0.14	
106-46-7	1,4-Dichlorobenzene	ND	0.82	ND	0.14	
95-50-1	1,2-Dichlorobenzene	ND	0.82	ND	0.14	
5989-27-5	d-Limonene	9.4	0.82	1.7	0.15	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.82	ND	0.084	
120-82-1	1,2,4-Trichlorobenzene	ND	0.82	ND	0.11	
87-68-3	Hexachlorobutadiene	ND	0.82	ND	0.076	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 41-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-007

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01819

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12 & 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.10 Liter(s)

Initial Pressure (psig): -3.37 Final Pressure (psig): 3.65

Canister Dilution Factor: 1.62

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	<b>18</b>	0.81	<b>10</b>	0.47	
75-71-8	Dichlorodifluoromethane (CFC 12)	<b>1.8</b>	0.81	<b>0.37</b>	0.16	
74-87-3	Chloromethane	ND	0.81	ND	0.39	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.81	ND	0.12	
106-99-0	1,3-Butadiene	ND	0.81	ND	0.37	
74-83-9	Bromomethane	ND	0.81	ND	0.21	
75-00-3	Chloroethane	ND	0.81	ND	0.31	
64-17-5	Ethanol	<b>1,400</b>	81	<b>760</b>	43	<b>D</b>
75-05-8	Acetonitrile	ND	0.81	ND	0.48	
107-02-8	Acrolein	ND	3.2	ND	1.4	
67-64-1	Acetone	<b>52</b>	8.1	<b>22</b>	3.4	
75-69-4	Trichlorofluoromethane	<b>1.7</b>	0.81	<b>0.31</b>	0.14	
67-63-0	2-Propanol (Isopropyl Alcohol)	<b>120</b>	1.6	<b>49</b>	0.66	
107-13-1	Acrylonitrile	ND	0.81	ND	0.37	
75-35-4	1,1-Dichloroethene	ND	0.81	ND	0.20	
75-09-2	Methylene Chloride	<b>0.87</b>	0.81	<b>0.25</b>	0.23	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.81	ND	0.26	
76-13-1	Trichlorotrifluoroethane	ND	0.81	ND	0.11	
75-15-0	Carbon Disulfide	ND	8.1	ND	2.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.81	ND	0.20	
75-34-3	1,1-Dichloroethane	ND	0.81	ND	0.20	
1634-04-4	Methyl tert-Butyl Ether	ND	0.81	ND	0.22	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 41-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-007

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01819

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12 & 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.10 Liter(s)

Initial Pressure (psig): -3.37 Final Pressure (psig): 3.65

Canister Dilution Factor: 1.62

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	8.1	ND	2.3	
156-59-2	cis-1,2-Dichloroethene	ND	0.81	ND	0.20	
78-93-3	2-Butanone (MEK)	ND	8.1	ND	2.7	
141-78-6	Ethyl Acetate	5.7	1.6	1.6	0.45	
110-54-3	n-Hexane	1.7	0.81	0.47	0.23	
67-66-3	Chloroform	2.7	0.81	0.55	0.17	
109-99-9	Tetrahydrofuran (THF)	ND	0.81	ND	0.27	
71-55-6	1,1,1-Trichloroethane	ND	0.81	ND	0.15	
56-23-5	Carbon Tetrachloride	ND	0.81	ND	0.13	
110-82-7	Cyclohexane	ND	1.6	ND	0.47	
78-87-5	1,2-Dichloropropane	ND	0.81	ND	0.18	
75-27-4	Bromodichloromethane	ND	0.81	ND	0.12	
123-91-1	1,4-Dioxane	ND	0.81	ND	0.22	
80-62-6	Methyl Methacrylate	ND	1.6	ND	0.40	
142-82-5	n-Heptane	1.1	0.81	0.27	0.20	
10061-01-5	cis-1,3-Dichloropropene	ND	0.81	ND	0.18	
108-10-1	4-Methyl-2-pentanone	2.1	0.81	0.52	0.20	
10061-02-6	trans-1,3-Dichloropropene	ND	0.81	ND	0.18	
108-88-3	Toluene	12	0.81	3.1	0.22	
591-78-6	2-Hexanone	ND	0.81	ND	0.20	
124-48-1	Dibromochloromethane	ND	0.81	ND	0.095	
106-93-4	1,2-Dibromoethane	ND	0.81	ND	0.11	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 41-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-007

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01819

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12 & 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.10 Liter(s)

Initial Pressure (psig): -3.37 Final Pressure (psig): 3.65

Canister Dilution Factor: 1.62

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	0.85	0.81	0.18	0.17	
111-65-9	n-Octane	0.93	0.81	0.20	0.17	
108-90-7	Chlorobenzene	ND	0.81	ND	0.18	
179601-23-1	m,p-Xylenes	4.3	1.6	0.99	0.37	
75-25-2	Bromoform	ND	0.81	ND	0.078	
100-42-5	Styrene	ND	0.81	ND	0.19	
95-47-6	o-Xylene	1.6	0.81	0.38	0.19	
111-84-2	n-Nonane	1.9	0.81	0.36	0.15	
98-82-8	Cumene	ND	0.81	ND	0.16	
80-56-8	alpha-Pinene	2.9	0.81	0.52	0.15	
103-65-1	n-Propylbenzene	ND	0.81	ND	0.16	
622-96-8	4-Ethyltoluene	ND	0.81	ND	0.16	
108-67-8	1,3,5-Trimethylbenzene	ND	0.81	ND	0.16	
95-63-6	1,2,4-Trimethylbenzene	2.0	0.81	0.42	0.16	
100-44-7	Benzyl Chloride	ND	0.81	ND	0.16	
541-73-1	1,3-Dichlorobenzene	ND	0.81	ND	0.13	
106-46-7	1,4-Dichlorobenzene	ND	0.81	ND	0.13	
95-50-1	1,2-Dichlorobenzene	ND	0.81	ND	0.13	
5989-27-5	d-Limonene	9.8	0.81	1.8	0.15	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.81	ND	0.084	
120-82-1	1,2,4-Trichlorobenzene	ND	0.81	ND	0.11	
87-68-3	Hexachlorobutadiene	ND	0.81	ND	0.076	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 48-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-008

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC00142

**Date Collected:** 3/29/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/7/12 & 4/9/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -5.15      Final Pressure (psig): 3.70

Canister Dilution Factor: 1.93

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	6.2	0.97	3.6	0.56	
75-71-8	Dichlorodifluoromethane (CFC 12)	1.4	0.97	0.28	0.20	
74-87-3	Chloromethane	ND	0.97	ND	0.47	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.97	ND	0.14	
106-99-0	1,3-Butadiene	ND	0.97	ND	0.44	
74-83-9	Bromomethane	ND	0.97	ND	0.25	
75-00-3	Chloroethane	ND	0.97	ND	0.37	
64-17-5	Ethanol	3,200	97	1,700	51	D
75-05-8	Acetonitrile	ND	0.97	ND	0.58	
107-02-8	Acrolein	4.5	3.9	1.9	1.7	
67-64-1	Acetone	110	9.7	45	4.1	
75-69-4	Trichlorofluoromethane	1.5	0.97	0.27	0.17	
67-63-0	2-Propanol (Isopropyl Alcohol)	56	1.9	23	0.79	
107-13-1	Acrylonitrile	ND	0.97	ND	0.44	
75-35-4	1,1-Dichloroethene	ND	0.97	ND	0.24	
75-09-2	Methylene Chloride	ND	0.97	ND	0.28	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.97	ND	0.31	
76-13-1	Trichlorotrifluoroethane	ND	0.97	ND	0.13	
75-15-0	Carbon Disulfide	ND	9.7	ND	3.1	
156-60-5	trans-1,2-Dichloroethene	ND	0.97	ND	0.24	
75-34-3	1,1-Dichloroethane	ND	0.97	ND	0.24	
1634-04-4	Methyl tert-Butyl Ether	ND	0.97	ND	0.27	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 48-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-008

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00142

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12 & 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.10 Liter(s)

Initial Pressure (psig): -5.15 Final Pressure (psig): 3.70

Canister Dilution Factor: 1.93

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	9.7	ND	2.7	
156-59-2	cis-1,2-Dichloroethene	ND	0.97	ND	0.24	
78-93-3	2-Butanone (MEK)	ND	9.7	ND	3.3	
141-78-6	Ethyl Acetate	<b>13</b>	1.9	<b>3.6</b>	0.54	
110-54-3	n-Hexane	<b>6.7</b>	0.97	<b>1.9</b>	0.27	
67-66-3	Chloroform	ND	0.97	ND	0.20	
109-99-9	Tetrahydrofuran (THF)	ND	0.97	ND	0.33	
71-55-6	1,1,1-Trichloroethane	ND	0.97	ND	0.18	
56-23-5	Carbon Tetrachloride	ND	0.97	ND	0.15	
110-82-7	Cyclohexane	ND	1.9	ND	0.56	
78-87-5	1,2-Dichloropropane	ND	0.97	ND	0.21	
75-27-4	Bromodichloromethane	ND	0.97	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.97	ND	0.27	
80-62-6	Methyl Methacrylate	ND	1.9	ND	0.47	
142-82-5	n-Heptane	<b>2.6</b>	0.97	<b>0.63</b>	0.24	
10061-01-5	cis-1,3-Dichloropropene	ND	0.97	ND	0.21	
108-10-1	4-Methyl-2-pentanone	<b>2.2</b>	0.97	<b>0.54</b>	0.24	
10061-02-6	trans-1,3-Dichloropropene	ND	0.97	ND	0.21	
108-88-3	Toluene	<b>31</b>	0.97	<b>8.2</b>	0.26	
591-78-6	2-Hexanone	ND	0.97	ND	0.24	
124-48-1	Dibromochloromethane	ND	0.97	ND	0.11	
106-93-4	1,2-Dibromoethane	ND	0.97	ND	0.13	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 48-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-008

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00142

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12 & 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.10 Liter(s)

Initial Pressure (psig): -5.15 Final Pressure (psig): 3.70

Canister Dilution Factor: 1.93

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	1.3	0.97	0.28	0.20	
111-65-9	n-Octane	1.2	0.97	0.27	0.21	
108-90-7	Chlorobenzene	ND	0.97	ND	0.21	
179601-23-1	m,p-Xylenes	18	1.9	4.1	0.44	
75-25-2	Bromoform	ND	0.97	ND	0.093	
100-42-5	Styrene	ND	0.97	ND	0.23	
95-47-6	o-Xylene	6.0	0.97	1.4	0.22	
111-84-2	n-Nonane	ND	0.97	ND	0.18	
98-82-8	Cumene	ND	0.97	ND	0.20	
80-56-8	alpha-Pinene	2.0	0.97	0.35	0.17	
103-65-1	n-Propylbenzene	0.97	0.97	0.20	0.20	
622-96-8	4-Ethyltoluene	1.8	0.97	0.38	0.20	
108-67-8	1,3,5-Trimethylbenzene	1.5	0.97	0.31	0.20	
95-63-6	1,2,4-Trimethylbenzene	4.9	0.97	1.0	0.20	
100-44-7	Benzyl Chloride	ND	0.97	ND	0.19	
541-73-1	1,3-Dichlorobenzene	ND	0.97	ND	0.16	
106-46-7	1,4-Dichlorobenzene	ND	0.97	ND	0.16	
95-50-1	1,2-Dichlorobenzene	ND	0.97	ND	0.16	
5989-27-5	d-Limonene	23	0.97	4.2	0.17	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.97	ND	0.10	
120-82-1	1,2,4-Trichlorobenzene	ND	0.97	ND	0.13	
87-68-3	Hexachlorobutadiene	ND	0.97	ND	0.091	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 43-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-009

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00603

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.54 Final Pressure (psig): 3.74

Canister Dilution Factor: 1.81

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	6.3	0.91	3.7	0.53	
75-71-8	Dichlorodifluoromethane (CFC 12)	1.7	0.91	0.34	0.18	
74-87-3	Chloromethane	ND	0.91	ND	0.44	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.91	ND	0.13	
106-99-0	1,3-Butadiene	ND	0.91	ND	0.41	
74-83-9	Bromomethane	ND	0.91	ND	0.23	
75-00-3	Chloroethane	ND	0.91	ND	0.34	
64-17-5	Ethanol	380	9.1	200	4.8	
75-05-8	Acetonitrile	ND	0.91	ND	0.54	
107-02-8	Acrolein	ND	3.6	ND	1.6	
67-64-1	Acetone	94	9.1	39	3.8	
75-69-4	Trichlorofluoromethane	1.8	0.91	0.33	0.16	
67-63-0	2-Propanol (Isopropyl Alcohol)	20	1.8	8.2	0.74	
107-13-1	Acrylonitrile	ND	0.91	ND	0.42	
75-35-4	1,1-Dichloroethene	ND	0.91	ND	0.23	
75-09-2	Methylene Chloride	100	0.91	29	0.26	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.91	ND	0.29	
76-13-1	Trichlorotrifluoroethane	ND	0.91	ND	0.12	
75-15-0	Carbon Disulfide	ND	9.1	ND	2.9	
156-60-5	trans-1,2-Dichloroethene	ND	0.91	ND	0.23	
75-34-3	1,1-Dichloroethane	ND	0.91	ND	0.22	
1634-04-4	Methyl tert-Butyl Ether	ND	0.91	ND	0.25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 43-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-009

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00603

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.54 Final Pressure (psig): 3.74

Canister Dilution Factor: 1.81

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	9.1	ND	2.6	
156-59-2	cis-1,2-Dichloroethene	ND	0.91	ND	0.23	
78-93-3	2-Butanone (MEK)	<b>13</b>	9.1	<b>4.5</b>	3.1	
141-78-6	Ethyl Acetate	<b>2.3</b>	1.8	<b>0.63</b>	0.50	
110-54-3	n-Hexane	ND	0.91	ND	0.26	
67-66-3	Chloroform	ND	0.91	ND	0.19	
109-99-9	Tetrahydrofuran (THF)	<b>4.6</b>	0.91	<b>1.6</b>	0.31	
71-55-6	1,1,1-Trichloroethane	ND	0.91	ND	0.17	
56-23-5	Carbon Tetrachloride	ND	0.91	ND	0.14	
110-82-7	Cyclohexane	ND	1.8	ND	0.53	
78-87-5	1,2-Dichloropropane	ND	0.91	ND	0.20	
75-27-4	Bromodichloromethane	ND	0.91	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.91	ND	0.25	
80-62-6	Methyl Methacrylate	ND	1.8	ND	0.44	
142-82-5	n-Heptane	<b>2.2</b>	0.91	<b>0.54</b>	0.22	
10061-01-5	cis-1,3-Dichloropropene	ND	0.91	ND	0.20	
108-10-1	4-Methyl-2-pentanone	ND	0.91	ND	0.22	
10061-02-6	trans-1,3-Dichloropropene	ND	0.91	ND	0.20	
108-88-3	Toluene	<b>20</b>	0.91	<b>5.3</b>	0.24	
591-78-6	2-Hexanone	ND	0.91	ND	0.22	
124-48-1	Dibromochloromethane	ND	0.91	ND	0.11	
106-93-4	1,2-Dibromoethane	ND	0.91	ND	0.12	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 43-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-009

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00603

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.54 Final Pressure (psig): 3.74

Canister Dilution Factor: 1.81

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	2.1	0.91	0.44	0.19	
111-65-9	n-Octane	ND	0.91	ND	0.19	
108-90-7	Chlorobenzene	ND	0.91	ND	0.20	
179601-23-1	m,p-Xylenes	5.8	1.8	1.3	0.42	
75-25-2	Bromoform	ND	0.91	ND	0.088	
100-42-5	Styrene	ND	0.91	ND	0.21	
95-47-6	o-Xylene	1.7	0.91	0.39	0.21	
111-84-2	n-Nonane	2.6	0.91	0.49	0.17	
98-82-8	Cumene	ND	0.91	ND	0.18	
80-56-8	alpha-Pinene	4.2	0.91	0.76	0.16	
103-65-1	n-Propylbenzene	ND	0.91	ND	0.18	
622-96-8	4-Ethyltoluene	ND	0.91	ND	0.18	
108-67-8	1,3,5-Trimethylbenzene	ND	0.91	ND	0.18	
95-63-6	1,2,4-Trimethylbenzene	1.2	0.91	0.23	0.18	
100-44-7	Benzyl Chloride	ND	0.91	ND	0.17	
541-73-1	1,3-Dichlorobenzene	ND	0.91	ND	0.15	
106-46-7	1,4-Dichlorobenzene	ND	0.91	ND	0.15	
95-50-1	1,2-Dichlorobenzene	ND	0.91	ND	0.15	
5989-27-5	d-Limonene	12	0.91	2.1	0.16	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.91	ND	0.094	
120-82-1	1,2,4-Trichlorobenzene	ND	0.91	ND	0.12	
87-68-3	Hexachlorobutadiene	ND	0.91	ND	0.085	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 45-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-010

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01626

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.65 Final Pressure (psig): 3.69

Canister Dilution Factor: 1.41

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	3.4	0.71	2.0	0.41	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.1	0.71	0.42	0.14	
74-87-3	Chloromethane	0.78	0.71	0.38	0.34	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.71	ND	0.10	
106-99-0	1,3-Butadiene	ND	0.71	ND	0.32	
74-83-9	Bromomethane	ND	0.71	ND	0.18	
75-00-3	Chloroethane	ND	0.71	ND	0.27	
64-17-5	Ethanol	480	7.1	250	3.7	
75-05-8	Acetonitrile	ND	0.71	ND	0.42	
107-02-8	Acrolein	ND	2.8	ND	1.2	
67-64-1	Acetone	43	7.1	18	3.0	
75-69-4	Trichlorofluoromethane	1.1	0.71	0.19	0.13	
67-63-0	2-Propanol (Isopropyl Alcohol)	21	1.4	8.7	0.57	
107-13-1	Acrylonitrile	ND	0.71	ND	0.32	
75-35-4	1,1-Dichloroethene	ND	0.71	ND	0.18	
75-09-2	Methylene Chloride	ND	0.71	ND	0.20	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.71	ND	0.23	
76-13-1	Trichlorotrifluoroethane	ND	0.71	ND	0.092	
75-15-0	Carbon Disulfide	ND	7.1	ND	2.3	
156-60-5	trans-1,2-Dichloroethene	ND	0.71	ND	0.18	
75-34-3	1,1-Dichloroethane	ND	0.71	ND	0.17	
1634-04-4	Methyl tert-Butyl Ether	ND	0.71	ND	0.20	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 45-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-010

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01626

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.65 Final Pressure (psig): 3.69

Canister Dilution Factor: 1.41

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	7.1	ND	2.0	
156-59-2	cis-1,2-Dichloroethene	ND	0.71	ND	0.18	
78-93-3	2-Butanone (MEK)	ND	7.1	ND	2.4	
141-78-6	Ethyl Acetate	<b>11</b>	1.4	<b>3.0</b>	0.39	
110-54-3	n-Hexane	<b>0.92</b>	0.71	<b>0.26</b>	0.20	
67-66-3	Chloroform	ND	0.71	ND	0.14	
109-99-9	Tetrahydrofuran (THF)	<b>1.6</b>	0.71	<b>0.53</b>	0.24	
71-55-6	1,1,1-Trichloroethane	ND	0.71	ND	0.13	
56-23-5	Carbon Tetrachloride	ND	0.71	ND	0.11	
110-82-7	Cyclohexane	ND	1.4	ND	0.41	
78-87-5	1,2-Dichloropropane	ND	0.71	ND	0.15	
75-27-4	Bromodichloromethane	ND	0.71	ND	0.11	
123-91-1	1,4-Dioxane	ND	0.71	ND	0.20	
80-62-6	Methyl Methacrylate	ND	1.4	ND	0.34	
142-82-5	n-Heptane	<b>0.71</b>	0.71	<b>0.17</b>	0.17	
10061-01-5	cis-1,3-Dichloropropene	ND	0.71	ND	0.16	
108-10-1	4-Methyl-2-pentanone	ND	0.71	ND	0.17	
10061-02-6	trans-1,3-Dichloropropene	ND	0.71	ND	0.16	
108-88-3	Toluene	<b>9.3</b>	0.71	<b>2.5</b>	0.19	
591-78-6	2-Hexanone	<b>0.73</b>	0.71	<b>0.18</b>	0.17	
124-48-1	Dibromochloromethane	ND	0.71	ND	0.083	
106-93-4	1,2-Dibromoethane	ND	0.71	ND	0.092	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 45-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-010

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01626

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.65 Final Pressure (psig): 3.69

Canister Dilution Factor: 1.41

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	3.5	0.71	0.73	0.15	
111-65-9	n-Octane	ND	0.71	ND	0.15	
108-90-7	Chlorobenzene	ND	0.71	ND	0.15	
179601-23-1	m,p-Xylenes	3.8	1.4	0.88	0.32	
75-25-2	Bromoform	ND	0.71	ND	0.068	
100-42-5	Styrene	1.2	0.71	0.28	0.17	
95-47-6	o-Xylene	1.3	0.71	0.30	0.16	
111-84-2	n-Nonane	ND	0.71	ND	0.13	
98-82-8	Cumene	ND	0.71	ND	0.14	
80-56-8	alpha-Pinene	3.4	0.71	0.61	0.13	
103-65-1	n-Propylbenzene	ND	0.71	ND	0.14	
622-96-8	4-Ethyltoluene	ND	0.71	ND	0.14	
108-67-8	1,3,5-Trimethylbenzene	ND	0.71	ND	0.14	
95-63-6	1,2,4-Trimethylbenzene	1.8	0.71	0.37	0.14	
100-44-7	Benzyl Chloride	ND	0.71	ND	0.14	
541-73-1	1,3-Dichlorobenzene	ND	0.71	ND	0.12	
106-46-7	1,4-Dichlorobenzene	ND	0.71	ND	0.12	
95-50-1	1,2-Dichlorobenzene	ND	0.71	ND	0.12	
5989-27-5	d-Limonene	13	0.71	2.4	0.13	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.71	ND	0.073	
120-82-1	1,2,4-Trichlorobenzene	ND	0.71	ND	0.095	
87-68-3	Hexachlorobutadiene	ND	0.71	ND	0.066	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 42-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-011

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00144

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.050 Liter(s)

Initial Pressure (psig): -2.74 Final Pressure (psig): 3.58

Canister Dilution Factor: 1.53

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	ND	0.77	ND	0.44	
75-71-8	Dichlorodifluoromethane (CFC 12)	<b>1.6</b>	0.77	<b>0.32</b>	0.15	
74-87-3	Chloromethane	ND	0.77	ND	0.37	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.77	ND	0.11	
106-99-0	1,3-Butadiene	ND	0.77	ND	0.35	
74-83-9	Bromomethane	ND	0.77	ND	0.20	
75-00-3	Chloroethane	ND	0.77	ND	0.29	
64-17-5	Ethanol	<b>2,600</b>	150	<b>1,400</b>	81	<b>D</b>
75-05-8	Acetonitrile	ND	0.77	ND	0.46	
107-02-8	Acrolein	<b>4.8</b>	3.1	<b>2.1</b>	1.3	
67-64-1	Acetone	<b>44</b>	7.7	<b>19</b>	3.2	
75-69-4	Trichlorofluoromethane	<b>0.84</b>	0.77	<b>0.15</b>	0.14	
67-63-0	2-Propanol (Isopropyl Alcohol)	<b>46</b>	1.5	<b>19</b>	0.62	
107-13-1	Acrylonitrile	ND	0.77	ND	0.35	
75-35-4	1,1-Dichloroethene	ND	0.77	ND	0.19	
75-09-2	Methylene Chloride	<b>1.1</b>	0.77	<b>0.33</b>	0.22	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.77	ND	0.24	
76-13-1	Trichlorotrifluoroethane	ND	0.77	ND	0.10	
75-15-0	Carbon Disulfide	ND	7.7	ND	2.5	
156-60-5	trans-1,2-Dichloroethene	ND	0.77	ND	0.19	
75-34-3	1,1-Dichloroethane	ND	0.77	ND	0.19	
1634-04-4	Methyl tert-Butyl Ether	ND	0.77	ND	0.21	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 42-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-011

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00144

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.050 Liter(s)

Initial Pressure (psig): -2.74 Final Pressure (psig): 3.58

Canister Dilution Factor: 1.53

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	7.7	ND	2.2	
156-59-2	cis-1,2-Dichloroethene	ND	0.77	ND	0.19	
78-93-3	2-Butanone (MEK)	ND	7.7	ND	2.6	
141-78-6	Ethyl Acetate	<b>6.3</b>	1.5	<b>1.8</b>	0.42	
110-54-3	n-Hexane	ND	0.77	ND	0.22	
67-66-3	Chloroform	ND	0.77	ND	0.16	
109-99-9	Tetrahydrofuran (THF)	ND	0.77	ND	0.26	
71-55-6	1,1,1-Trichloroethane	ND	0.77	ND	0.14	
56-23-5	Carbon Tetrachloride	ND	0.77	ND	0.12	
110-82-7	Cyclohexane	ND	1.5	ND	0.44	
78-87-5	1,2-Dichloropropane	ND	0.77	ND	0.17	
75-27-4	Bromodichloromethane	ND	0.77	ND	0.11	
123-91-1	1,4-Dioxane	ND	0.77	ND	0.21	
80-62-6	Methyl Methacrylate	ND	1.5	ND	0.37	
142-82-5	n-Heptane	<b>1.1</b>	0.77	<b>0.28</b>	0.19	
10061-01-5	cis-1,3-Dichloropropene	ND	0.77	ND	0.17	
108-10-1	4-Methyl-2-pentanone	ND	0.77	ND	0.19	
10061-02-6	trans-1,3-Dichloropropene	ND	0.77	ND	0.17	
108-88-3	Toluene	<b>3.3</b>	0.77	<b>0.87</b>	0.20	
591-78-6	2-Hexanone	ND	0.77	ND	0.19	
124-48-1	Dibromochloromethane	ND	0.77	ND	0.090	
106-93-4	1,2-Dibromoethane	ND	0.77	ND	0.10	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 42-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-011

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00144

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)  
 0.050 Liter(s)

Initial Pressure (psig): -2.74 Final Pressure (psig): 3.58

Canister Dilution Factor: 1.53

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	0.83	0.77	0.17	0.16	
111-65-9	n-Octane	ND	0.77	ND	0.16	
108-90-7	Chlorobenzene	ND	0.77	ND	0.17	
179601-23-1	m,p-Xylenes	ND	1.5	ND	0.35	
75-25-2	Bromoform	ND	0.77	ND	0.074	
100-42-5	Styrene	ND	0.77	ND	0.18	
95-47-6	o-Xylene	ND	0.77	ND	0.18	
111-84-2	n-Nonane	ND	0.77	ND	0.15	
98-82-8	Cumene	ND	0.77	ND	0.16	
80-56-8	alpha-Pinene	4.9	0.77	0.88	0.14	
103-65-1	n-Propylbenzene	ND	0.77	ND	0.16	
622-96-8	4-Ethyltoluene	ND	0.77	ND	0.16	
108-67-8	1,3,5-Trimethylbenzene	ND	0.77	ND	0.16	
95-63-6	1,2,4-Trimethylbenzene	ND	0.77	ND	0.16	
100-44-7	Benzyl Chloride	ND	0.77	ND	0.15	
541-73-1	1,3-Dichlorobenzene	ND	0.77	ND	0.13	
106-46-7	1,4-Dichlorobenzene	ND	0.77	ND	0.13	
95-50-1	1,2-Dichlorobenzene	ND	0.77	ND	0.13	
5989-27-5	d-Limonene	24	0.77	4.3	0.14	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.77	ND	0.079	
120-82-1	1,2,4-Trichlorobenzene	ND	0.77	ND	0.10	
87-68-3	Hexachlorobutadiene	ND	0.77	ND	0.072	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 44-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-012

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01183

**Date Collected:** 3/29/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/9/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -3.92 Final Pressure (psig): 3.76

Canister Dilution Factor: 1.71

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	0.95	0.86	0.55	0.50	
75-71-8	Dichlorodifluoromethane (CFC 12)	1.8	0.86	0.36	0.17	
74-87-3	Chloromethane	ND	0.86	ND	0.41	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.86	ND	0.12	
106-99-0	1,3-Butadiene	ND	0.86	ND	0.39	
74-83-9	Bromomethane	ND	0.86	ND	0.22	
75-00-3	Chloroethane	ND	0.86	ND	0.32	
64-17-5	Ethanol	1,400	86	740	45	D
75-05-8	Acetonitrile	ND	0.86	ND	0.51	
107-02-8	Acrolein	ND	3.4	ND	1.5	
67-64-1	Acetone	19	8.6	8.1	3.6	
75-69-4	Trichlorofluoromethane	1.0	0.86	0.18	0.15	
67-63-0	2-Propanol (Isopropyl Alcohol)	4.0	1.7	1.6	0.70	
107-13-1	Acrylonitrile	ND	0.86	ND	0.39	
75-35-4	1,1-Dichloroethene	ND	0.86	ND	0.22	
75-09-2	Methylene Chloride	ND	0.86	ND	0.25	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.86	ND	0.27	
76-13-1	Trichlorotrifluoroethane	ND	0.86	ND	0.11	
75-15-0	Carbon Disulfide	ND	8.6	ND	2.7	
156-60-5	trans-1,2-Dichloroethene	ND	0.86	ND	0.22	
75-34-3	1,1-Dichloroethane	ND	0.86	ND	0.21	
1634-04-4	Methyl tert-Butyl Ether	ND	0.86	ND	0.24	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.



## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 44-1F-032912  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201319  
**CAS Sample ID:** P1201319-012

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01183

**Date Collected:** 3/29/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/9/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -3.92 Final Pressure (psig): 3.76

Canister Dilution Factor: 1.71

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	8.6	ND	2.4	
156-59-2	cis-1,2-Dichloroethene	ND	0.86	ND	0.22	
78-93-3	2-Butanone (MEK)	ND	8.6	ND	2.9	
141-78-6	Ethyl Acetate	4.5	1.7	1.2	0.47	
110-54-3	n-Hexane	ND	0.86	ND	0.24	
67-66-3	Chloroform	ND	0.86	ND	0.18	
109-99-9	Tetrahydrofuran (THF)	ND	0.86	ND	0.29	
71-55-6	1,1,1-Trichloroethane	ND	0.86	ND	0.16	
56-23-5	Carbon Tetrachloride	ND	0.86	ND	0.14	
110-82-7	Cyclohexane	ND	1.7	ND	0.50	
78-87-5	1,2-Dichloropropane	ND	0.86	ND	0.19	
75-27-4	Bromodichloromethane	ND	0.86	ND	0.13	
123-91-1	1,4-Dioxane	ND	0.86	ND	0.24	
80-62-6	Methyl Methacrylate	ND	1.7	ND	0.42	
142-82-5	n-Heptane	ND	0.86	ND	0.21	
10061-01-5	cis-1,3-Dichloropropene	ND	0.86	ND	0.19	
108-10-1	4-Methyl-2-pentanone	ND	0.86	ND	0.21	
10061-02-6	trans-1,3-Dichloropropene	ND	0.86	ND	0.19	
108-88-3	Toluene	3.3	0.86	0.87	0.23	
591-78-6	2-Hexanone	ND	0.86	ND	0.21	
124-48-1	Dibromochloromethane	ND	0.86	ND	0.10	
106-93-4	1,2-Dibromoethane	ND	0.86	ND	0.11	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 44-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-012

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01183

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.10 Liter(s)

Initial Pressure (psig): -3.92 Final Pressure (psig): 3.76

Canister Dilution Factor: 1.71

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.86	ND	0.18	
111-65-9	n-Octane	ND	0.86	ND	0.18	
108-90-7	Chlorobenzene	ND	0.86	ND	0.19	
179601-23-1	m,p-Xylenes	<b>2.6</b>	1.7	<b>0.60</b>	0.39	
75-25-2	Bromoform	ND	0.86	ND	0.083	
100-42-5	Styrene	ND	0.86	ND	0.20	
95-47-6	o-Xylene	ND	0.86	ND	0.20	
111-84-2	n-Nonane	ND	0.86	ND	0.16	
98-82-8	Cumene	ND	0.86	ND	0.17	
80-56-8	alpha-Pinene	<b>0.99</b>	0.86	<b>0.18</b>	0.15	
103-65-1	n-Propylbenzene	ND	0.86	ND	0.17	
622-96-8	4-Ethyltoluene	ND	0.86	ND	0.17	
108-67-8	1,3,5-Trimethylbenzene	ND	0.86	ND	0.17	
95-63-6	1,2,4-Trimethylbenzene	ND	0.86	ND	0.17	
100-44-7	Benzyl Chloride	ND	0.86	ND	0.17	
541-73-1	1,3-Dichlorobenzene	ND	0.86	ND	0.14	
106-46-7	1,4-Dichlorobenzene	ND	0.86	ND	0.14	
95-50-1	1,2-Dichlorobenzene	ND	0.86	ND	0.14	
5989-27-5	d-Limonene	<b>3.3</b>	0.86	<b>0.60</b>	0.15	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.86	ND	0.088	
120-82-1	1,2,4-Trichlorobenzene	ND	0.86	ND	0.12	
87-68-3	Hexachlorobutadiene	ND	0.86	ND	0.080	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 46-1F-033012  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-013

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00810

Date Collected: 3/30/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.10 Liter(s)

Initial Pressure (psig): -4.91 Final Pressure (psig): 3.67

Canister Dilution Factor: 1.88

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	7.4	0.94	4.3	0.55	
75-71-8	Dichlorodifluoromethane (CFC 12)	1.8	0.94	0.36	0.19	
74-87-3	Chloromethane	ND	0.94	ND	0.46	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.94	ND	0.13	
106-99-0	1,3-Butadiene	ND	0.94	ND	0.43	
74-83-9	Bromomethane	ND	0.94	ND	0.24	
75-00-3	Chloroethane	ND	0.94	ND	0.36	
64-17-5	Ethanol	3,700	94	2,000	50	D
75-05-8	Acetonitrile	ND	0.94	ND	0.56	
107-02-8	Acrolein	ND	3.8	ND	1.6	
67-64-1	Acetone	56	9.4	23	4.0	
75-69-4	Trichlorofluoromethane	1.1	0.94	0.19	0.17	
67-63-0	2-Propanol (Isopropyl Alcohol)	81	1.9	33	0.77	
107-13-1	Acrylonitrile	ND	0.94	ND	0.43	
75-35-4	1,1-Dichloroethene	ND	0.94	ND	0.24	
75-09-2	Methylene Chloride	1.3	0.94	0.37	0.27	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.94	ND	0.30	
76-13-1	Trichlorotrifluoroethane	ND	0.94	ND	0.12	
75-15-0	Carbon Disulfide	ND	9.4	ND	3.0	
156-60-5	trans-1,2-Dichloroethene	ND	0.94	ND	0.24	
75-34-3	1,1-Dichloroethane	ND	0.94	ND	0.23	
1634-04-4	Methyl tert-Butyl Ether	ND	0.94	ND	0.26	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 46-1F-033012  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-013

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC00810

**Date Collected:** 3/30/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/9/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -4.91 Final Pressure (psig): 3.67

Canister Dilution Factor: 1.88

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	9.4	ND	2.7	
156-59-2	cis-1,2-Dichloroethene	ND	0.94	ND	0.24	
78-93-3	2-Butanone (MEK)	ND	9.4	ND	3.2	
141-78-6	Ethyl Acetate	13	1.9	3.7	0.52	
110-54-3	n-Hexane	2.2	0.94	0.64	0.27	
67-66-3	Chloroform	1.4	0.94	0.29	0.19	
109-99-9	Tetrahydrofuran (THF)	ND	0.94	ND	0.32	
71-55-6	1,1,1-Trichloroethane	ND	0.94	ND	0.17	
56-23-5	Carbon Tetrachloride	ND	0.94	ND	0.15	
110-82-7	Cyclohexane	ND	1.9	ND	0.55	
78-87-5	1,2-Dichloropropane	ND	0.94	ND	0.20	
75-27-4	Bromodichloromethane	ND	0.94	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.94	ND	0.26	
80-62-6	Methyl Methacrylate	ND	1.9	ND	0.46	
142-82-5	n-Heptane	1.7	0.94	0.42	0.23	
10061-01-5	cis-1,3-Dichloropropene	ND	0.94	ND	0.21	
108-10-1	4-Methyl-2-pentanone	ND	0.94	ND	0.23	
10061-02-6	trans-1,3-Dichloropropene	ND	0.94	ND	0.21	
108-88-3	Toluene	130	0.94	34	0.25	
591-78-6	2-Hexanone	ND	0.94	ND	0.23	
124-48-1	Dibromochloromethane	ND	0.94	ND	0.11	
106-93-4	1,2-Dibromoethane	ND	0.94	ND	0.12	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 46-1F-033012  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-013

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00810

Date Collected: 3/30/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -4.91 Final Pressure (psig): 3.67

Canister Dilution Factor: 1.88

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	3.6	0.94	0.77	0.20	
111-65-9	n-Octane	12	0.94	2.7	0.20	
108-90-7	Chlorobenzene	ND	0.94	ND	0.20	
179601-23-1	m,p-Xylenes	2.0	1.9	0.47	0.43	
75-25-2	Bromoform	ND	0.94	ND	0.091	
100-42-5	Styrene	ND	0.94	ND	0.22	
95-47-6	o-Xylene	ND	0.94	ND	0.22	
111-84-2	n-Nonane	4.9	0.94	0.93	0.18	
98-82-8	Cumene	ND	0.94	ND	0.19	
80-56-8	alpha-Pinene	13	0.94	2.3	0.17	
103-65-1	n-Propylbenzene	ND	0.94	ND	0.19	
622-96-8	4-Ethyltoluene	ND	0.94	ND	0.19	
108-67-8	1,3,5-Trimethylbenzene	ND	0.94	ND	0.19	
95-63-6	1,2,4-Trimethylbenzene	ND	0.94	ND	0.19	
100-44-7	Benzyl Chloride	ND	0.94	ND	0.18	
541-73-1	1,3-Dichlorobenzene	ND	0.94	ND	0.16	
106-46-7	1,4-Dichlorobenzene	ND	0.94	ND	0.16	
95-50-1	1,2-Dichlorobenzene	ND	0.94	ND	0.16	
5989-27-5	d-Limonene	15	0.94	2.7	0.17	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.94	ND	0.097	
120-82-1	1,2,4-Trichlorobenzene	ND	0.94	ND	0.13	
87-68-3	Hexachlorobutadiene	ND	0.94	ND	0.088	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 47-1F-033112  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201319  
**CAS Sample ID:** P1201319-014

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01122

**Date Collected:** 3/31/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/9/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -5.26 Final Pressure (psig): 3.76

Canister Dilution Factor: 1.96

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	2.0	0.98	1.2	0.57	
75-71-8	Dichlorodifluoromethane (CFC 12)	1.7	0.98	0.35	0.20	
74-87-3	Chloromethane	ND	0.98	ND	0.47	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.98	ND	0.14	
106-99-0	1,3-Butadiene	ND	0.98	ND	0.44	
74-83-9	Bromomethane	ND	0.98	ND	0.25	
75-00-3	Chloroethane	ND	0.98	ND	0.37	
64-17-5	Ethanol	710	9.8	380	5.2	
75-05-8	Acetonitrile	ND	0.98	ND	0.58	
107-02-8	Acrolein	ND	3.9	ND	1.7	
67-64-1	Acetone	33	9.8	14	4.1	
75-69-4	Trichlorofluoromethane	ND	0.98	ND	0.17	
67-63-0	2-Propanol (Isopropyl Alcohol)	13	2.0	5.4	0.80	
107-13-1	Acrylonitrile	ND	0.98	ND	0.45	
75-35-4	1,1-Dichloroethene	ND	0.98	ND	0.25	
75-09-2	Methylene Chloride	ND	0.98	ND	0.28	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.98	ND	0.31	
76-13-1	Trichlorotrifluoroethane	ND	0.98	ND	0.13	
75-15-0	Carbon Disulfide	ND	9.8	ND	3.1	
156-60-5	trans-1,2-Dichloroethene	ND	0.98	ND	0.25	
75-34-3	1,1-Dichloroethane	ND	0.98	ND	0.24	
1634-04-4	Methyl tert-Butyl Ether	ND	0.98	ND	0.27	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 47-1F-033112  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-014

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01122

Date Collected: 3/31/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.26 Final Pressure (psig): 3.76

Canister Dilution Factor: 1.96

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	9.8	ND	2.8	
156-59-2	cis-1,2-Dichloroethene	ND	0.98	ND	0.25	
78-93-3	2-Butanone (MEK)	ND	9.8	ND	3.3	
141-78-6	Ethyl Acetate	3.9	2.0	1.1	0.54	
110-54-3	n-Hexane	1.5	0.98	0.44	0.28	
67-66-3	Chloroform	ND	0.98	ND	0.20	
109-99-9	Tetrahydrofuran (THF)	2.6	0.98	0.87	0.33	
71-55-6	1,1,1-Trichloroethane	ND	0.98	ND	0.18	
56-23-5	Carbon Tetrachloride	ND	0.98	ND	0.16	
110-82-7	Cyclohexane	ND	2.0	ND	0.57	
78-87-5	1,2-Dichloropropane	ND	0.98	ND	0.21	
75-27-4	Bromodichloromethane	ND	0.98	ND	0.15	
123-91-1	1,4-Dioxane	ND	0.98	ND	0.27	
80-62-6	Methyl Methacrylate	ND	2.0	ND	0.48	
142-82-5	n-Heptane	ND	0.98	ND	0.24	
10061-01-5	cis-1,3-Dichloropropene	ND	0.98	ND	0.22	
108-10-1	4-Methyl-2-pentanone	ND	0.98	ND	0.24	
10061-02-6	trans-1,3-Dichloropropene	ND	0.98	ND	0.22	
108-88-3	Toluene	5.6	0.98	1.5	0.26	
591-78-6	2-Hexanone	ND	0.98	ND	0.24	
124-48-1	Dibromochloromethane	ND	0.98	ND	0.12	
106-93-4	1,2-Dibromoethane	ND	0.98	ND	0.13	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 47-1F-033112  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-014

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01122

Date Collected: 3/31/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.26 Final Pressure (psig): 3.76

Canister Dilution Factor: 1.96

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	1.2	0.98	0.26	0.21	
111-65-9	n-Octane	ND	0.98	ND	0.21	
108-90-7	Chlorobenzene	ND	0.98	ND	0.21	
179601-23-1	m,p-Xylenes	2.0	2.0	0.46	0.45	
75-25-2	Bromoform	ND	0.98	ND	0.095	
100-42-5	Styrene	1.2	0.98	0.28	0.23	
95-47-6	o-Xylene	ND	0.98	ND	0.23	
111-84-2	n-Nonane	ND	0.98	ND	0.19	
98-82-8	Cumene	ND	0.98	ND	0.20	
80-56-8	alpha-Pinene	5.0	0.98	0.90	0.18	
103-65-1	n-Propylbenzene	ND	0.98	ND	0.20	
622-96-8	4-Ethyltoluene	ND	0.98	ND	0.20	
108-67-8	1,3,5-Trimethylbenzene	ND	0.98	ND	0.20	
95-63-6	1,2,4-Trimethylbenzene	ND	0.98	ND	0.20	
100-44-7	Benzyl Chloride	ND	0.98	ND	0.19	
541-73-1	1,3-Dichlorobenzene	ND	0.98	ND	0.16	
106-46-7	1,4-Dichlorobenzene	ND	0.98	ND	0.16	
95-50-1	1,2-Dichlorobenzene	ND	0.98	ND	0.16	
5989-27-5	d-Limonene	11	0.98	2.0	0.18	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.98	ND	0.10	
120-82-1	1,2,4-Trichlorobenzene	ND	0.98	ND	0.13	
87-68-3	Hexachlorobutadiene	ND	0.98	ND	0.092	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P120407-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	ND	0.50	ND	0.29	
75-71-8	Dichlorodifluoromethane (CFC 12)	ND	0.50	ND	0.10	
74-87-3	Chloromethane	ND	0.50	ND	0.24	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.50	ND	0.072	
106-99-0	1,3-Butadiene	ND	0.50	ND	0.23	
74-83-9	Bromomethane	ND	0.50	ND	0.13	
75-00-3	Chloroethane	ND	0.50	ND	0.19	
64-17-5	Ethanol	ND	5.0	ND	2.7	
75-05-8	Acetonitrile	ND	0.50	ND	0.30	
107-02-8	Acrolein	ND	2.0	ND	0.87	
67-64-1	Acetone	ND	5.0	ND	2.1	
75-69-4	Trichlorofluoromethane	ND	0.50	ND	0.089	
67-63-0	2-Propanol (Isopropyl Alcohol)	ND	1.0	ND	0.41	
107-13-1	Acrylonitrile	ND	0.50	ND	0.23	
75-35-4	1,1-Dichloroethene	ND	0.50	ND	0.13	
75-09-2	Methylene Chloride	ND	0.50	ND	0.14	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.50	ND	0.16	
76-13-1	Trichlorotrifluoroethane	ND	0.50	ND	0.065	
75-15-0	Carbon Disulfide	ND	5.0	ND	1.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.50	ND	0.13	
75-34-3	1,1-Dichloroethane	ND	0.50	ND	0.12	
1634-04-4	Methyl tert-Butyl Ether	ND	0.50	ND	0.14	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P120407-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	5.0	ND	1.4	
156-59-2	cis-1,2-Dichloroethene	ND	0.50	ND	0.13	
78-93-3	2-Butanone (MEK)	ND	5.0	ND	1.7	
141-78-6	Ethyl Acetate	ND	1.0	ND	0.28	
110-54-3	n-Hexane	ND	0.50	ND	0.14	
67-66-3	Chloroform	ND	0.50	ND	0.10	
109-99-9	Tetrahydrofuran (THF)	ND	0.50	ND	0.17	
71-55-6	1,1,1-Trichloroethane	ND	0.50	ND	0.092	
56-23-5	Carbon Tetrachloride	ND	0.50	ND	0.080	
110-82-7	Cyclohexane	ND	1.0	ND	0.29	
78-87-5	1,2-Dichloropropane	ND	0.50	ND	0.11	
75-27-4	Bromodichloromethane	ND	0.50	ND	0.075	
123-91-1	1,4-Dioxane	ND	0.50	ND	0.14	
80-62-6	Methyl Methacrylate	ND	1.0	ND	0.24	
142-82-5	n-Heptane	ND	0.50	ND	0.12	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-10-1	4-Methyl-2-pentanone	ND	0.50	ND	0.12	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-88-3	Toluene	ND	0.50	ND	0.13	
591-78-6	2-Hexanone	ND	0.50	ND	0.12	
124-48-1	Dibromochloromethane	ND	0.50	ND	0.059	
106-93-4	1,2-Dibromoethane	ND	0.50	ND	0.065	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P120407-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.50	ND	0.11	
111-65-9	n-Octane	ND	0.50	ND	0.11	
108-90-7	Chlorobenzene	ND	0.50	ND	0.11	
179601-23-1	m,p-Xylenes	ND	1.0	ND	0.23	
75-25-2	Bromoform	ND	0.50	ND	0.048	
100-42-5	Styrene	ND	0.50	ND	0.12	
95-47-6	o-Xylene	ND	0.50	ND	0.12	
111-84-2	n-Nonane	ND	0.50	ND	0.095	
98-82-8	Cumene	ND	0.50	ND	0.10	
80-56-8	alpha-Pinene	ND	0.50	ND	0.090	
103-65-1	n-Propylbenzene	ND	0.50	ND	0.10	
622-96-8	4-Ethyltoluene	ND	0.50	ND	0.10	
108-67-8	1,3,5-Trimethylbenzene	ND	0.50	ND	0.10	
95-63-6	1,2,4-Trimethylbenzene	ND	0.50	ND	0.10	
100-44-7	Benzyl Chloride	ND	0.50	ND	0.097	
541-73-1	1,3-Dichlorobenzene	ND	0.50	ND	0.083	
106-46-7	1,4-Dichlorobenzene	ND	0.50	ND	0.083	
95-50-1	1,2-Dichlorobenzene	ND	0.50	ND	0.083	
5989-27-5	d-Limonene	ND	0.50	ND	0.090	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.50	ND	0.052	
120-82-1	1,2,4-Trichlorobenzene	ND	0.50	ND	0.067	
87-68-3	Hexachlorobutadiene	ND	0.50	ND	0.047	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P120409-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result		MRL		Data Qualifier
		$\mu\text{g}/\text{m}^3$	$\mu\text{g}/\text{m}^3$	ppbV	ppbV	
115-07-1	Propene	ND	0.50	ND	0.29	
75-71-8	Dichlorodifluoromethane (CFC 12)	ND	0.50	ND	0.10	
74-87-3	Chloromethane	ND	0.50	ND	0.24	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.50	ND	0.072	
106-99-0	1,3-Butadiene	ND	0.50	ND	0.23	
74-83-9	Bromomethane	ND	0.50	ND	0.13	
75-00-3	Chloroethane	ND	0.50	ND	0.19	
64-17-5	Ethanol	ND	5.0	ND	2.7	
75-05-8	Acetonitrile	ND	0.50	ND	0.30	
107-02-8	Acrolein	ND	2.0	ND	0.87	
67-64-1	Acetone	ND	5.0	ND	2.1	
75-69-4	Trichlorofluoromethane	ND	0.50	ND	0.089	
67-63-0	2-Propanol (Isopropyl Alcohol)	ND	1.0	ND	0.41	
107-13-1	Acrylonitrile	ND	0.50	ND	0.23	
75-35-4	1,1-Dichloroethene	ND	0.50	ND	0.13	
75-09-2	Methylene Chloride	ND	0.50	ND	0.14	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.50	ND	0.16	
76-13-1	Trichlorotrifluoroethane	ND	0.50	ND	0.065	
75-15-0	Carbon Disulfide	ND	5.0	ND	1.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.50	ND	0.13	
75-34-3	1,1-Dichloroethane	ND	0.50	ND	0.12	
1634-04-4	Methyl tert-Butyl Ether	ND	0.50	ND	0.14	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P120409-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	5.0	ND	1.4	
156-59-2	cis-1,2-Dichloroethene	ND	0.50	ND	0.13	
78-93-3	2-Butanone (MEK)	ND	5.0	ND	1.7	
141-78-6	Ethyl Acetate	ND	1.0	ND	0.28	
110-54-3	n-Hexane	ND	0.50	ND	0.14	
67-66-3	Chloroform	ND	0.50	ND	0.10	
109-99-9	Tetrahydrofuran (THF)	ND	0.50	ND	0.17	
71-55-6	1,1,1-Trichloroethane	ND	0.50	ND	0.092	
56-23-5	Carbon Tetrachloride	ND	0.50	ND	0.080	
110-82-7	Cyclohexane	ND	1.0	ND	0.29	
78-87-5	1,2-Dichloropropane	ND	0.50	ND	0.11	
75-27-4	Bromodichloromethane	ND	0.50	ND	0.075	
123-91-1	1,4-Dioxane	ND	0.50	ND	0.14	
80-62-6	Methyl Methacrylate	ND	1.0	ND	0.24	
142-82-5	n-Heptane	ND	0.50	ND	0.12	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-10-1	4-Methyl-2-pentanone	ND	0.50	ND	0.12	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-88-3	Toluene	ND	0.50	ND	0.13	
591-78-6	2-Hexanone	ND	0.50	ND	0.12	
124-48-1	Dibromochloromethane	ND	0.50	ND	0.059	
106-93-4	1,2-Dibromoethane	ND	0.50	ND	0.065	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P120409-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.50	ND	0.11	
111-65-9	n-Octane	ND	0.50	ND	0.11	
108-90-7	Chlorobenzene	ND	0.50	ND	0.11	
179601-23-1	m,p-Xylenes	ND	1.0	ND	0.23	
75-25-2	Bromoform	ND	0.50	ND	0.048	
100-42-5	Styrene	ND	0.50	ND	0.12	
95-47-6	o-Xylene	ND	0.50	ND	0.12	
111-84-2	n-Nonane	ND	0.50	ND	0.095	
98-82-8	Cumene	ND	0.50	ND	0.10	
80-56-8	alpha-Pinene	ND	0.50	ND	0.090	
103-65-1	n-Propylbenzene	ND	0.50	ND	0.10	
622-96-8	4-Ethyltoluene	ND	0.50	ND	0.10	
108-67-8	1,3,5-Trimethylbenzene	ND	0.50	ND	0.10	
95-63-6	1,2,4-Trimethylbenzene	ND	0.50	ND	0.10	
100-44-7	Benzyl Chloride	ND	0.50	ND	0.097	
541-73-1	1,3-Dichlorobenzene	ND	0.50	ND	0.083	
106-46-7	1,4-Dichlorobenzene	ND	0.50	ND	0.083	
95-50-1	1,2-Dichlorobenzene	ND	0.50	ND	0.083	
5989-27-5	d-Limonene	ND	0.50	ND	0.090	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.50	ND	0.052	
120-82-1	1,2,4-Trichlorobenzene	ND	0.50	ND	0.067	
87-68-3	Hexachlorobutadiene	ND	0.50	ND	0.047	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister(s)  
**Test Notes:**

**Date(s) Collected:** 3/25 - 3/31/12  
**Date(s) Received:** 4/4/12  
**Date(s) Analyzed:** 4/7 - 4/9/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		Percent Recovered	Percent Recovered	Percent Recovered		
Method Blank	P120407-MB	91	96	102	70-130	
Method Blank	P120409-MB	93	99	101	70-130	
Lab Control Sample	P120407-LCS	90	97	103	70-130	
Lab Control Sample	P120409-LCS	91	99	102	70-130	
50-2F-032512	P1201319-001	92	98	99	70-130	
49-1F-032712	P1201319-002	92	98	100	70-130	
38-1F-032812	P1201319-003	93	97	100	70-130	
39-1F-032812	P1201319-004	92	98	102	70-130	
40-1F-032812	P1201319-005	93	97	100	70-130	
55-1F-032812	P1201319-006	93	98	99	70-130	
41-1F-032912	P1201319-007	93	99	99	70-130	
48-1F-032912	P1201319-008	93	96	99	70-130	
43-1F-032912	P1201319-009	92	98	100	70-130	
45-1F-032912	P1201319-010	93	99	100	70-130	
42-1F-032912	P1201319-011	90	99	103	70-130	
44-1F-032912	P1201319-012	90	99	102	70-130	
46-1F-033012	P1201319-013	90	101	103	70-130	
47-1F-033112	P1201319-014	90	98	101	70-130	
47-1F-033112	P1201319-014DUP	90	98	102	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.

Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P120407-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
**Analyst:** Lusine Hakobyan  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 4/07/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
115-07-1	Propene	198	205	104	59-137	
75-71-8	Dichlorodifluoromethane (CFC 12)	196	145	74	63-115	
74-87-3	Chloromethane	190	175	92	59-124	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	200	155	78	65-113	
106-99-0	1,3-Butadiene	204	193	95	60-138	
74-83-9	Bromomethane	194	158	81	69-129	
75-00-3	Chloroethane	196	177	90	60-120	
64-17-5	Ethanol	928	866	93	58-121	
75-05-8	Acetonitrile	194	185	95	64-129	
107-02-8	Acrolein	198	169	85	54-127	
67-64-1	Acetone	1,010	859	85	59-114	
75-69-4	Trichlorofluoromethane	202	143	71	66-108	
67-63-0	2-Propanol (Isopropyl Alcohol)	382	282	74	50-113	
107-13-1	Acrylonitrile	198	207	105	72-135	
75-35-4	1,1-Dichloroethene	212	182	86	70-117	
75-09-2	Methylene Chloride	206	170	83	61-108	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	208	194	93	70-131	
76-13-1	Trichlorotrifluoroethane	206	172	83	70-113	
75-15-0	Carbon Disulfide	208	175	84	65-112	
156-60-5	trans-1,2-Dichloroethene	196	175	89	71-119	
75-34-3	1,1-Dichloroethane	200	173	87	71-116	
1634-04-4	Methyl tert-Butyl Ether	198	148	75	67-116	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.



LABORATORY CONTROL SAMPLE SUMMARY

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P120407-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/07/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
108-05-4	Vinyl Acetate	952	805	85	59-142	
156-59-2	cis-1,2-Dichloroethene	206	175	85	69-119	
78-93-3	2-Butanone (MEK)	206	188	91	68-125	
141-78-6	Ethyl Acetate	398	337	85	63-130	
110-54-3	n-Hexane	198	151	76	57-120	
67-66-3	Chloroform	214	164	77	69-111	
109-99-9	Tetrahydrofuran (THF)	202	175	87	57-123	
71-55-6	1,1,1-Trichloroethane	198	155	78	73-119	
56-23-5	Carbon Tetrachloride	202	174	86	74-129	
110-82-7	Cyclohexane	390	314	81	70-113	
78-87-5	1,2-Dichloropropane	198	173	87	69-118	
75-27-4	Bromodichloromethane	198	161	81	75-124	
123-91-1	1,4-Dioxane	200	172	86	71-123	
80-62-6	Methyl Methacrylate	400	353	88	72-127	
142-82-5	n-Heptane	196	165	84	68-120	
10061-01-5	cis-1,3-Dichloropropene	188	168	89	71-130	
108-10-1	4-Methyl-2-pentanone	204	187	92	69-130	
10061-02-6	trans-1,3-Dichloropropene	210	194	92	76-133	
108-88-3	Toluene	202	157	78	67-111	
591-78-6	2-Hexanone	222	167	75	70-123	
124-48-1	Dibromochloromethane	206	171	83	75-129	
106-93-4	1,2-Dibromoethane	200	166	83	73-122	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P120407-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/07/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
123-86-4	n-Butyl Acetate	220	177	80	68-132	
111-65-9	n-Octane	200	161	81	68-116	
108-90-7	Chlorobenzene	202	161	80	69-113	
179601-23-1	m,p-Xylenes	392	304	78	70-116	
75-25-2	Bromoform	208	169	81	69-127	
100-42-5	Styrene	200	166	83	71-125	
95-47-6	o-Xylene	194	152	78	70-116	
111-84-2	n-Nonane	196	150	77	68-116	
98-82-8	Cumene	190	149	78	70-116	
80-56-8	alpha-Pinene	186	147	79	71-119	
103-65-1	n-Propylbenzene	192	150	78	71-119	
622-96-8	4-Ethyltoluene	198	163	82	71-119	
108-67-8	1,3,5-Trimethylbenzene	200	158	79	71-121	
95-63-6	1,2,4-Trimethylbenzene	194	151	78	73-127	
100-44-7	Benzyl Chloride	200	166	83	65-137	
541-73-1	1,3-Dichlorobenzene	200	164	82	68-123	
106-46-7	1,4-Dichlorobenzene	206	162	79	65-120	
95-50-1	1,2-Dichlorobenzene	198	157	79	67-121	
5989-27-5	d-Limonene	200	151	76	67-130	
96-12-8	1,2-Dibromo-3-chloropropane	196	173	88	72-133	
120-82-1	1,2,4-Trichlorobenzene	196	178	91	62-133	
87-68-3	Hexachlorobutadiene	202	178	88	60-128	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P120409-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/09/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
115-07-1	Propene	198	226	114	59-137	
75-71-8	Dichlorodifluoromethane (CFC 12)	196	162	83	63-115	
74-87-3	Chloromethane	190	202	106	59-124	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	200	173	87	65-113	
106-99-0	1,3-Butadiene	204	224	110	60-138	
74-83-9	Bromomethane	194	178	92	69-129	
75-00-3	Chloroethane	196	203	104	60-120	
64-17-5	Ethanol	928	1020	110	58-121	
75-05-8	Acetonitrile	194	217	112	64-129	
107-02-8	Acrolein	198	195	98	54-127	
67-64-1	Acetone	1,010	989	98	59-114	
75-69-4	Trichlorofluoromethane	202	160	79	66-108	
67-63-0	2-Propanol (Isopropyl Alcohol)	382	347	91	50-113	
107-13-1	Acrylonitrile	198	237	120	72-135	
75-35-4	1,1-Dichloroethene	212	202	95	70-117	
75-09-2	Methylene Chloride	206	189	92	61-108	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	208	227	109	70-131	
76-13-1	Trichlorotrifluoroethane	206	192	93	70-113	
75-15-0	Carbon Disulfide	208	195	94	65-112	
156-60-5	trans-1,2-Dichloroethene	196	198	101	71-119	
75-34-3	1,1-Dichloroethane	200	196	98	71-116	
1634-04-4	Methyl tert-Butyl Ether	198	168	85	67-116	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result. Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P120409-LCS

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/09/12  
 Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
108-05-4	Vinyl Acetate	952	908	95	59-142	
156-59-2	cis-1,2-Dichloroethene	206	197	96	69-119	
78-93-3	2-Butanone (MEK)	206	210	102	68-125	
141-78-6	Ethyl Acetate	398	383	96	63-130	
110-54-3	n-Hexane	198	173	87	57-120	
67-66-3	Chloroform	214	182	85	69-111	
109-99-9	Tetrahydrofuran (THF)	202	198	98	57-123	
71-55-6	1,1,1-Trichloroethane	198	171	86	73-119	
56-23-5	Carbon Tetrachloride	202	192	95	74-129	
110-82-7	Cyclohexane	390	350	90	70-113	
78-87-5	1,2-Dichloropropane	198	193	97	69-118	
75-27-4	Bromodichloromethane	198	178	90	75-124	
123-91-1	1,4-Dioxane	200	190	95	71-123	
80-62-6	Methyl Methacrylate	400	395	99	72-127	
142-82-5	n-Heptane	196	185	94	68-120	
10061-01-5	cis-1,3-Dichloropropene	188	189	101	71-130	
108-10-1	4-Methyl-2-pentanone	204	210	103	69-130	
10061-02-6	trans-1,3-Dichloropropene	210	217	103	76-133	
108-88-3	Toluene	202	178	88	67-111	
591-78-6	2-Hexanone	222	193	87	70-123	
124-48-1	Dibromochloromethane	206	193	94	75-129	
106-93-4	1,2-Dibromoethane	200	185	93	73-122	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result. Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P120409-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/09/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
123-86-4	n-Butyl Acetate	220	205	93	68-132	
111-65-9	n-Octane	200	187	94	68-116	
108-90-7	Chlorobenzene	202	181	90	69-113	
179601-23-1	m,p-Xylenes	392	346	88	70-116	
75-25-2	Bromoform	208	193	93	69-127	
100-42-5	Styrene	200	188	94	71-125	
95-47-6	o-Xylene	194	173	89	70-116	
111-84-2	n-Nonane	196	175	89	68-116	
98-82-8	Cumene	190	170	89	70-116	
80-56-8	alpha-Pinene	186	166	89	71-119	
103-65-1	n-Propylbenzene	192	171	89	71-119	
622-96-8	4-Ethyltoluene	198	173	87	71-119	
108-67-8	1,3,5-Trimethylbenzene	200	181	91	71-121	
95-63-6	1,2,4-Trimethylbenzene	194	174	90	73-127	
100-44-7	Benzyl Chloride	200	190	95	65-137	
541-73-1	1,3-Dichlorobenzene	200	185	93	68-123	
106-46-7	1,4-Dichlorobenzene	206	183	89	65-120	
95-50-1	1,2-Dichlorobenzene	198	179	90	67-121	
5989-27-5	d-Limonene	200	171	86	67-130	
96-12-8	1,2-Dibromo-3-chloropropane	196	197	101	72-133	
120-82-1	1,2,4-Trichlorobenzene	196	202	103	62-133	
87-68-3	Hexachlorobutadiene	202	204	101	60-128	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY DUPLICATE SUMMARY RESULTS

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**Client:** Montana DEQ  
**Client Sample ID:** 47-1F-033112  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-014DUP

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
Analyst: Lusine Hakobyan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01122

Date Collected: 3/31/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.26

Final Pressure (psig): 3.76

Canister Dilution Factor: 1.96

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
Propene	2.00	1.16	1.87	1.09	1.935	7	25	
Dichlorodifluoromethane (CFC 12)	1.74	0.352	1.75	0.353	1.745	0.6	25	
Chloromethane	ND	ND	ND	ND	-	-	25	
1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	ND	ND	ND	-	-	25	
1,3-Butadiene	ND	ND	ND	ND	-	-	25	
Bromomethane	ND	ND	ND	ND	-	-	25	
Chloroethane	ND	ND	ND	ND	-	-	25	
Ethanol	713	378	705	375	709	1	25	
Acetonitrile	ND	ND	ND	ND	-	-	25	
Acrolein	ND	ND	ND	ND	-	-	25	
Acetone	32.7	13.8	32.4	13.6	32.55	0.9	25	
Trichlorofluoromethane	ND	ND	ND	ND	-	-	25	
2-Propanol (Isopropyl Alcohol)	13.2	5.38	13.1	5.35	13.15	0.8	25	
Acrylonitrile	ND	ND	ND	ND	-	-	25	
1,1-Dichloroethene	ND	ND	ND	ND	-	-	25	
Methylene Chloride	ND	ND	ND	ND	-	-	25	
3-Chloro-1-propene (Allyl Chloride)	ND	ND	ND	ND	-	-	25	
Trichlorotrifluoroethane	ND	ND	ND	ND	-	-	25	
Carbon Disulfide	ND	ND	ND	ND	-	-	25	
trans-1,2-Dichloroethene	ND	ND	ND	ND	-	-	25	
1,1-Dichloroethane	ND	ND	ND	ND	-	-	25	
Methyl tert-Butyl Ether	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

LABORATORY DUPLICATE SUMMARY RESULTS

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**Client:** Montana DEQ  
**Client Sample ID:** 47-1F-033112  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-014DUP

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01122

Date Collected: 3/31/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.26

Final Pressure (psig): 3.76

Canister Dilution Factor: 1.96

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
Vinyl Acetate	ND	ND	ND	ND	-	-	25	
cis-1,2-Dichloroethene	ND	ND	ND	ND	-	-	25	
2-Butanone (MEK)	ND	ND	ND	ND	-	-	25	
Ethyl Acetate	3.88	1.08	3.82	1.06	3.85	2	25	
n-Hexane	1.55	0.439	1.55	0.440	1.55	0	25	
Chloroform	ND	ND	ND	ND	-	-	25	
Tetrahydrofuran (THF)	2.55	0.866	2.50	0.848	2.525	2	25	
1,1,1-Trichloroethane	ND	ND	ND	ND	-	-	25	
Carbon Tetrachloride	ND	ND	ND	ND	-	-	25	
Cyclohexane	ND	ND	ND	ND	-	-	25	
1,2-Dichloropropane	ND	ND	ND	ND	-	-	25	
Bromodichloromethane	ND	ND	ND	ND	-	-	25	
1,4-Dioxane	ND	ND	ND	ND	-	-	25	
Methyl Methacrylate	ND	ND	ND	ND	-	-	25	
n-Heptane	ND	ND	ND	ND	-	-	25	
cis-1,3-Dichloropropene	ND	ND	ND	ND	-	-	25	
4-Methyl-2-pentanone	ND	ND	ND	ND	-	-	25	
trans-1,3-Dichloropropene	ND	ND	ND	ND	-	-	25	
Toluene	5.57	1.48	5.47	1.45	5.52	2	25	
2-Hexanone	ND	ND	ND	ND	-	-	25	
Dibromochloromethane	ND	ND	ND	ND	-	-	25	
1,2-Dibromoethane	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

LABORATORY DUPLICATE SUMMARY RESULTS

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 47-1F-033112  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-014DUP

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16  
 Analyst: Lusine Hakobyan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01122

Date Collected: 3/31/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.26

Final Pressure (psig): 3.76

Canister Dilution Factor: 1.96

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
n-Butyl Acetate	1.24	0.262	1.20	0.252	1.22	<b>3</b>	25	
n-Octane	ND	ND	ND	ND	-	-	25	
Chlorobenzene	ND	ND	ND	ND	-	-	25	
m,p-Xylenes	1.98	0.455	1.96	0.452	1.97	<b>1</b>	25	
Bromoform	ND	ND	ND	ND	-	-	25	
Styrene	1.20	0.282	1.18	0.278	1.19	<b>2</b>	25	
o-Xylene	ND	ND	ND	ND	-	-	25	
n-Nonane	ND	ND	ND	ND	-	-	25	
Cumene	ND	ND	ND	ND	-	-	25	
alpha-Pinene	5.00	0.897	4.95	0.888	4.975	<b>1</b>	25	
n-Propylbenzene	ND	ND	ND	ND	-	-	25	
4-Ethyltoluene	ND	ND	ND	ND	-	-	25	
1,3,5-Trimethylbenzene	ND	ND	ND	ND	-	-	25	
1,2,4-Trimethylbenzene	ND	ND	ND	ND	-	-	25	
Benzyl Chloride	ND	ND	ND	ND	-	-	25	
1,3-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
1,4-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
1,2-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
d-Limonene	11.0	1.98	11.0	1.97	11	<b>0</b>	25	
1,2-Dibromo-3-chloropropane	ND	ND	ND	ND	-	-	25	
1,2,4-Trichlorobenzene	ND	ND	ND	ND	-	-	25	
Hexachlorobutadiene	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 50-2F-032512  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-001

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01225

Date Collected: 3/25/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.46      Final Pressure (psig): 3.75  
 Initial Pressure 2 (psig): -3.36      Final Pressure 2 (psig): 2.06

Canister Dilution Factor: 2.95

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.074	ND	0.029	
107-06-2	1,2-Dichloroethane	<b>0.12</b>	0.074	<b>0.030</b>	0.018	
71-43-2	Benzene	<b>15</b>	0.22	<b>4.6</b>	0.069	
79-01-6	Trichloroethene	ND	0.074	ND	0.014	
79-00-5	1,1,2-Trichloroethane	ND	0.30	ND	0.054	
127-18-4	Tetrachloroethene	<b>2.0</b>	0.074	<b>0.30</b>	0.011	
100-41-4	Ethylbenzene	<b>6.1</b>	0.30	<b>1.4</b>	0.068	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.074	ND	0.011	
91-20-3	Naphthalene	<b>1.9</b>	0.30	<b>0.37</b>	0.056	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 49-1F-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-002

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01223

Date Collected: 3/27/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.77      Final Pressure (psig): 3.75

Canister Dilution Factor: 1.55

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.039	ND	0.015	
107-06-2	1,2-Dichloroethane	<b>0.36</b>	0.039	<b>0.090</b>	0.0096	
71-43-2	Benzene	<b>3.1</b>	0.12	<b>0.96</b>	0.036	
79-01-6	Trichloroethene	<b>0.050</b>	0.039	<b>0.0093</b>	0.0072	
79-00-5	1,1,2-Trichloroethane	ND	0.16	ND	0.028	
127-18-4	Tetrachloroethene	<b>0.097</b>	0.039	<b>0.014</b>	0.0057	
100-41-4	Ethylbenzene	<b>1.3</b>	0.16	<b>0.30</b>	0.036	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.039	ND	0.0056	
91-20-3	Naphthalene	<b>0.35</b>	0.16	<b>0.066</b>	0.030	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 38-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-003

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00993

Date Collected: 3/28/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9 - 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -3.55      Final Pressure (psig): 3.63

Canister Dilution Factor: 1.64

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.041	ND	0.016	
107-06-2	1,2-Dichloroethane	<b>0.068</b>	0.041	<b>0.017</b>	0.010	
71-43-2	Benzene	<b>0.38</b>	0.12	<b>0.12</b>	0.039	
79-01-6	Trichloroethene	<b>0.042</b>	0.041	<b>0.0078</b>	0.0076	
79-00-5	1,1,2-Trichloroethane	ND	0.16	ND	0.030	
127-18-4	Tetrachloroethene	<b>0.047</b>	0.041	<b>0.0069</b>	0.0060	
100-41-4	Ethylbenzene	<b>0.31</b>	0.16	<b>0.072</b>	0.038	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.041	ND	0.0060	
91-20-3	Naphthalene	<b>130</b>	1.6	<b>24</b>	0.31	<b>D</b>

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 39-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-004

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01628

Date Collected: 3/28/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.36      Final Pressure (psig): 3.62

Canister Dilution Factor: 1.48

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.037	ND	0.014	
107-06-2	1,2-Dichloroethane	<b>0.067</b>	0.037	<b>0.017</b>	0.0091	
71-43-2	Benzene	<b>0.48</b>	0.11	<b>0.15</b>	0.035	
79-01-6	Trichloroethene	ND	0.037	ND	0.0069	
79-00-5	1,1,2-Trichloroethane	ND	0.15	ND	0.027	
127-18-4	Tetrachloroethene	<b>0.082</b>	0.037	<b>0.012</b>	0.0055	
100-41-4	Ethylbenzene	<b>0.31</b>	0.15	<b>0.071</b>	0.034	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.037	ND	0.0054	
91-20-3	Naphthalene	<b>0.26</b>	0.15	<b>0.049</b>	0.028	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 40-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-005

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00588

Date Collected: 3/28/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.22      Final Pressure (psig): 3.74

Canister Dilution Factor: 1.76

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.044	ND	0.017	
107-06-2	1,2-Dichloroethane	<b>0.36</b>	0.044	<b>0.090</b>	0.011	
71-43-2	Benzene	<b>0.73</b>	0.13	<b>0.23</b>	0.041	
79-01-6	Trichloroethene	<b>0.54</b>	0.044	<b>0.10</b>	0.0082	
79-00-5	1,1,2-Trichloroethane	ND	0.18	ND	0.032	
127-18-4	Tetrachloroethene	ND	0.044	ND	0.0065	
100-41-4	Ethylbenzene	ND	0.18	ND	0.041	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.044	ND	0.0064	
91-20-3	Naphthalene	<b>0.42</b>	0.18	<b>0.080</b>	0.034	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 55-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-006

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00131

Date Collected: 3/28/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.45      Final Pressure (psig): 3.63

Canister Dilution Factor: 1.63

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.041	ND	0.016	
107-06-2	1,2-Dichloroethane	<b>0.24</b>	0.041	<b>0.060</b>	0.010	
71-43-2	Benzene	<b>0.70</b>	0.12	<b>0.22</b>	0.038	
79-01-6	Trichloroethene	<b>0.54</b>	0.041	<b>0.10</b>	0.0076	
79-00-5	1,1,2-Trichloroethane	ND	0.16	ND	0.030	
127-18-4	Tetrachloroethene	ND	0.041	ND	0.0060	
100-41-4	Ethylbenzene	ND	0.16	ND	0.038	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.041	ND	0.0059	
91-20-3	Naphthalene	<b>0.30</b>	0.16	<b>0.058</b>	0.031	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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**Client:** Montana DEQ  
**Client Sample ID:** 41-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-007

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01819

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.37      Final Pressure (psig): 3.65

Canister Dilution Factor: 1.62

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.041	ND	0.016	
107-06-2	1,2-Dichloroethane	<b>0.11</b>	0.041	<b>0.027</b>	0.010	
71-43-2	Benzene	<b>1.9</b>	0.12	<b>0.59</b>	0.038	
79-01-6	Trichloroethene	ND	0.041	ND	0.0075	
79-00-5	1,1,2-Trichloroethane	ND	0.16	ND	0.030	
127-18-4	Tetrachloroethene	<b>0.22</b>	0.041	<b>0.032</b>	0.0060	
100-41-4	Ethylbenzene	<b>1.0</b>	0.16	<b>0.24</b>	0.037	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.041	ND	0.0059	
91-20-3	Naphthalene	<b>0.51</b>	0.16	<b>0.098</b>	0.031	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 48-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-008

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00142

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.15      Final Pressure (psig): 3.70

Canister Dilution Factor: 1.93

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.048	ND	0.019	
107-06-2	1,2-Dichloroethane	<b>0.62</b>	0.048	<b>0.15</b>	0.012	
71-43-2	Benzene	<b>9.8</b>	0.14	<b>3.1</b>	0.045	
79-01-6	Trichloroethene	<b>0.057</b>	0.048	<b>0.011</b>	0.0090	
79-00-5	1,1,2-Trichloroethane	ND	0.19	ND	0.035	
127-18-4	Tetrachloroethene	<b>3.2</b>	0.048	<b>0.47</b>	0.0071	
100-41-4	Ethylbenzene	<b>4.5</b>	0.19	<b>1.0</b>	0.044	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.048	ND	0.0070	
91-20-3	Naphthalene	<b>0.50</b>	0.19	<b>0.096</b>	0.037	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 43-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-009

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00603

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.54      Final Pressure (psig): 3.74

Canister Dilution Factor: 1.81

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.045	ND	0.018	
107-06-2	1,2-Dichloroethane	<b>0.12</b>	0.045	<b>0.029</b>	0.011	
71-43-2	Benzene	<b>0.51</b>	0.14	<b>0.16</b>	0.043	
79-01-6	Trichloroethene	ND	0.045	ND	0.0084	
79-00-5	1,1,2-Trichloroethane	ND	0.18	ND	0.033	
127-18-4	Tetrachloroethene	<b>0.69</b>	0.045	<b>0.10</b>	0.0067	
100-41-4	Ethylbenzene	<b>1.5</b>	0.18	<b>0.34</b>	0.042	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.045	ND	0.0066	
91-20-3	Naphthalene	<b>0.18</b>	0.18	<b>0.035</b>	0.035	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 45-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-010

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01626

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.65      Final Pressure (psig): 3.69

Canister Dilution Factor: 1.41

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.035	ND	0.014	
107-06-2	1,2-Dichloroethane	<b>0.13</b>	0.035	<b>0.032</b>	0.0087	
71-43-2	Benzene	<b>1.3</b>	0.11	<b>0.42</b>	0.033	
79-01-6	Trichloroethene	ND	0.035	ND	0.0066	
79-00-5	1,1,2-Trichloroethane	ND	0.14	ND	0.026	
127-18-4	Tetrachloroethene	<b>0.24</b>	0.035	<b>0.035</b>	0.0052	
100-41-4	Ethylbenzene	<b>0.89</b>	0.14	<b>0.20</b>	0.032	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.035	ND	0.0051	
91-20-3	Naphthalene	<b>0.28</b>	0.14	<b>0.054</b>	0.027	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 42-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-011

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00144

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.74      Final Pressure (psig): 3.58

Canister Dilution Factor: 1.53

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.038	ND	0.015	
107-06-2	1,2-Dichloroethane	<b>0.19</b>	0.038	<b>0.046</b>	0.0095	
71-43-2	Benzene	<b>0.46</b>	0.11	<b>0.14</b>	0.036	
79-01-6	Trichloroethene	ND	0.038	ND	0.0071	
79-00-5	1,1,2-Trichloroethane	ND	0.15	ND	0.028	
127-18-4	Tetrachloroethene	<b>2.9</b>	0.038	<b>0.43</b>	0.0056	
100-41-4	Ethylbenzene	<b>0.39</b>	0.15	<b>0.091</b>	0.035	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.038	ND	0.0056	
91-20-3	Naphthalene	<b>0.64</b>	0.15	<b>0.12</b>	0.029	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 44-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-012

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01183

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.92      Final Pressure (psig): 3.76

Canister Dilution Factor: 1.71

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.043	ND	0.017	
107-06-2	1,2-Dichloroethane	<b>0.50</b>	0.043	<b>0.12</b>	0.011	
71-43-2	Benzene	<b>0.83</b>	0.13	<b>0.26</b>	0.040	
79-01-6	Trichloroethene	ND	0.043	ND	0.0080	
79-00-5	1,1,2-Trichloroethane	ND	0.17	ND	0.031	
127-18-4	Tetrachloroethene	<b>0.064</b>	0.043	<b>0.0095</b>	0.0063	
100-41-4	Ethylbenzene	<b>0.79</b>	0.17	<b>0.18</b>	0.039	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.043	ND	0.0062	
91-20-3	Naphthalene	ND	0.17	ND	0.033	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 46-1F-033012  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-013

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00810

Date Collected: 3/30/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.91      Final Pressure (psig): 3.67

Canister Dilution Factor: 1.88

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.047	ND	0.018	
107-06-2	1,2-Dichloroethane	<b>0.091</b>	0.047	<b>0.023</b>	0.012	
71-43-2	Benzene	<b>1.4</b>	0.14	<b>0.43</b>	0.044	
79-01-6	Trichloroethene	ND	0.047	ND	0.0087	
79-00-5	1,1,2-Trichloroethane	ND	0.19	ND	0.034	
127-18-4	Tetrachloroethene	<b>0.50</b>	0.047	<b>0.074</b>	0.0069	
100-41-4	Ethylbenzene	<b>0.52</b>	0.19	<b>0.12</b>	0.043	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.047	ND	0.0068	
91-20-3	Naphthalene	<b>0.40</b>	0.19	<b>0.077</b>	0.036	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 47-1F-033112  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P1201319-014

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01122

Date Collected: 3/31/12  
Date Received: 4/4/12  
Date Analyzed: 4/10/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.26      Final Pressure (psig): 3.76

Canister Dilution Factor: 1.96

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.049	ND	0.019	
107-06-2	1,2-Dichloroethane	<b>2.0</b>	0.049	<b>0.49</b>	0.012	
71-43-2	Benzene	<b>1.4</b>	0.15	<b>0.43</b>	0.046	
79-01-6	Trichloroethene	ND	0.049	ND	0.0091	
79-00-5	1,1,2-Trichloroethane	ND	0.20	ND	0.036	
127-18-4	Tetrachloroethene	<b>0.13</b>	0.049	<b>0.020</b>	0.0072	
100-41-4	Ethylbenzene	<b>0.61</b>	0.20	<b>0.14</b>	0.045	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.049	ND	0.0071	
91-20-3	Naphthalene	<b>0.22</b>	0.20	<b>0.042</b>	0.037	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P120409-MB

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.025	ND	0.0098	
107-06-2	1,2-Dichloroethane	ND	0.025	ND	0.0062	
71-43-2	Benzene	ND	0.075	ND	0.023	
79-01-6	Trichloroethene	ND	0.025	ND	0.0047	
79-00-5	1,1,2-Trichloroethane	ND	0.10	ND	0.018	
127-18-4	Tetrachloroethene	ND	0.025	ND	0.0037	
100-41-4	Ethylbenzene	ND	0.10	ND	0.023	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.025	ND	0.0036	
91-20-3	Naphthalene	ND	0.10	ND	0.019	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P120410-MB

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result		MRL		Data Qualifier
		$\mu\text{g}/\text{m}^3$	$\mu\text{g}/\text{m}^3$	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.025	ND	0.0098	
107-06-2	1,2-Dichloroethane	ND	0.025	ND	0.0062	
71-43-2	Benzene	ND	0.075	ND	0.023	
79-01-6	Trichloroethene	ND	0.025	ND	0.0047	
79-00-5	1,1,2-Trichloroethane	ND	0.10	ND	0.018	
127-18-4	Tetrachloroethene	ND	0.025	ND	0.0037	
100-41-4	Ethylbenzene	ND	0.10	ND	0.023	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.025	ND	0.0036	
91-20-3	Naphthalene	ND	0.10	ND	0.019	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister(s)  
 Test Notes:

Date(s) Collected: 3/25 - 3/31/12  
 Date(s) Received: 4/4/12  
 Date(s) Analyzed: 4/9 - 4/10/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		% Recovered	% Recovered	% Recovered		
Method Blank	P120409-MB	101	110	108	70-130	
Method Blank	P120410-MB	100	109	107	70-130	
Lab Control Sample	P120409-LCS	102	110	110	70-130	
Lab Control Sample	P120410-LCS	96	108	109	70-130	
50-2F-032512	P1201319-001	99	107	102	70-130	
49-1F-032712	P1201319-002	96	110	104	70-130	
38-1F-032812	P1201319-003	97	112	109	70-130	
39-1F-032812	P1201319-004	99	112	102	70-130	
40-1F-032812	P1201319-005	97	108	110	70-130	
55-1F-032812	P1201319-006	97	109	110	70-130	
41-1F-032912	P1201319-007	96	110	105	70-130	
48-1F-032912	P1201319-008	96	107	109	70-130	
43-1F-032912	P1201319-009	97	110	106	70-130	
45-1F-032912	P1201319-010	97	110	107	70-130	
42-1F-032912	P1201319-011	96	111	108	70-130	
44-1F-032912	P1201319-012	98	108	108	70-130	
44-1F-032912	P1201319-012DUP	97	110	108	70-130	
46-1F-033012	P1201319-013	98	111	99	70-130	
47-1F-033112	P1201319-014	99	109	108	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.

Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.

LABORATORY CONTROL SAMPLE SUMMARY

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**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P120409-LCS

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/09/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
75-01-4	Vinyl Chloride	3.88	<b>3.08</b>	<b>79</b>	56-127	
107-06-2	1,2-Dichloroethane	4.00	<b>3.13</b>	<b>78</b>	51-140	
71-43-2	Benzene	3.96	<b>3.51</b>	<b>89</b>	56-125	
79-01-6	Trichloroethene	3.88	<b>2.97</b>	<b>77</b>	51-127	
79-00-5	1,1,2-Trichloroethane	3.92	<b>2.89</b>	<b>74</b>	49-137	
127-18-4	Tetrachloroethene	3.68	<b>2.91</b>	<b>79</b>	58-134	
100-41-4	Ethylbenzene	3.96	<b>3.18</b>	<b>80</b>	56-126	
79-34-5	1,1,2,2-Tetrachloroethane	3.84	<b>3.04</b>	<b>79</b>	53-148	
91-20-3	Naphthalene	3.44	<b>3.08</b>	<b>90</b>	25-149	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
CAS Sample ID: P120410-LCS

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/10/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
75-01-4	Vinyl Chloride	3.88	<b>2.91</b>	<b>75</b>	56-127	
107-06-2	1,2-Dichloroethane	4.00	<b>2.90</b>	<b>73</b>	51-140	
71-43-2	Benzene	3.96	<b>3.49</b>	<b>88</b>	56-125	
79-01-6	Trichloroethene	3.88	<b>2.84</b>	<b>73</b>	51-127	
79-00-5	1,1,2-Trichloroethane	3.92	<b>2.77</b>	<b>71</b>	49-137	
127-18-4	Tetrachloroethene	3.68	<b>2.82</b>	<b>77</b>	58-134	
100-41-4	Ethylbenzene	3.96	<b>3.25</b>	<b>82</b>	56-126	
79-34-5	1,1,2,2-Tetrachloroethane	3.84	<b>2.92</b>	<b>76</b>	53-148	
91-20-3	Naphthalene	3.44	<b>3.12</b>	<b>91</b>	25-149	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY DUPLICATE SUMMARY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 44-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201319  
 CAS Sample ID: P1201319-012DUP

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01183

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -3.92      Final Pressure (psig): 3.76

Canister Dilution Factor: 1.71

CAS #	Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
		µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
75-01-4	Vinyl Chloride	ND	ND	ND	ND	-	-	25	
107-06-2	1,2-Dichloroethane	0.496	0.123	0.491	0.121	0.4935	<b>1</b>	25	
71-43-2	Benzene	0.827	0.259	0.827	0.259	0.827	<b>0</b>	25	
79-01-6	Trichloroethene	ND	ND	ND	ND	-	-	25	
79-00-5	1,1,2-Trichloroethane	ND	ND	ND	ND	-	-	25	
127-18-4	Tetrachloroethene	0.0642	0.00947	0.0641	0.00946	0.06415	<b>0.2</b>	25	
100-41-4	Ethylbenzene	0.790	0.182	0.774	0.178	0.782	<b>2</b>	25	
79-34-5	1,1,2,2-Tetrachloroethane	ND	ND	ND	ND	-	-	25	
91-20-3	Naphthalene	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

Massachusetts APH  
Hydrocarbon Ranges

ICAL Method: M16022712.M

ICAL Date: 2/27/12

Instrument ID: MS16

	areas				masses							
	0.5	1	5	25	50	100	0.5	1	5	25	50	100
<b>C5-C8 Aliphatics</b>												
Isopentane	73866	131100	566770	2255808	4522556	8936172	0.520	1.04	5.20	26.0	52.0	104
n-Hexane	78752	129416	532939	2431156	4850068	9520785	0.490	0.98	4.90	24.5	49.0	98
Cyclohexane	71093	131903	610831	2881536	5778021	11355310	0.500	1.00	5.00	25.0	50.0	100
2,3-Dimethylpentane	71330	134802	622005	2909216	5793262	11304627	0.520	1.04	5.20	26.0	52.0	104
n-Heptane	66345	125793	590534	2785727	5552279	10734061	0.490	0.98	4.90	24.5	49.0	98
n-Octane	77181	142834	669465	3160716	6264037	11839954	0.505	1.01	5.05	25.3	50.5	101
<b>area sum:</b>	<b>438567</b>	<b>795848</b>	<b>3592544</b>	<b>16424159</b>	<b>32760223</b>	<b>63690909</b>	<b>3.025</b>	<b>6.050</b>	<b>30.25</b>	<b>151.30</b>	<b>302.50</b>	<b>605.00</b>

<b>C9-C12 Aliphatics</b>												
2,3-Dimethylheptane	83205	154255	731522	3440887	6755851	12700073	0.485	0.97	4.85	24.3	48.5	97
n-Nonane	77550	145499	691987	3274543	6427207	12148314	0.485	0.97	4.85	24.3	48.5	97
n-Decane	78814	152614	738750	3503011	6899987	12851504	0.490	0.98	4.90	24.5	49.0	98
1,3-Butylcyclohexane	97891	184987	896859	4276591	8376082	15228652	0.495	0.99	4.95	24.8	49.5	99
n-Undecane	79443	152957	755651	3602026	7106272	13091473	0.480	0.96	4.80	24.0	48.0	96
n-Dodecane	75428	146260	753969	3513032	7018288	13181075	0.500	1.00	5.00	25.0	50.0	100
<b>area sum:</b>	<b>492331</b>	<b>936572</b>	<b>4568738</b>	<b>21610090</b>	<b>42583687</b>	<b>79201091</b>	<b>2.935</b>	<b>5.870</b>	<b>29.35</b>	<b>146.90</b>	<b>293.50</b>	<b>587.00</b>

<b>C9-C10 Aromatics</b>												
Isopropylbenzene	12820	24226	114789	564731	1111501	1927344	0.485	0.97	4.85	24.3	48.5	97
3-Ethyltoluene	14043	26099	127717	633290	1241349	2166999	0.470	0.94	4.70	23.5	47.0	94
1,3,5-Trimethylbenzene	19126	35909	175219	859825	1663097	2896573	0.490	0.98	4.90	24.5	49.0	98
1,2,3-Trimethylbenzene	20434	38242	191458	916640	1653887	2575128	0.495	0.99	4.95	24.8	49.5	99
p-Isopropyltoluene	11392	21601	106407	521026	932930	1434539	0.460	0.92	4.60	23.0	46.0	92
<b>area sum:</b>	<b>77815</b>	<b>146077</b>	<b>715590</b>	<b>3495512</b>	<b>6602764</b>	<b>11000583</b>	<b>2.400</b>	<b>4.800</b>	<b>24.00</b>	<b>120.10</b>	<b>240.00</b>	<b>480.00</b>

Massachusetts APH  
Hydrocarbon Ranges

ICAL: M1602712M

ICAL Date: 2/27/12

Instrument ID: MS16

areas

<u>Internal Standards (TIC)</u>	0.5	1	5	25	50	100
Bromochloromethane (IS1)	1821684	1812431	2017469	1784633	1989673	2035509
1,4-Difluorobenzene (IS2)	3489421	3485174	3835800	3398785	3833709	3906245
Chlorobenzene-d5 (IS3)	4171549	4155528	4598058	4065257	4603508	4715823

Internal Standards (EIC)

Bromochloromethane (IS1)	364034	360985	402561	357490	400808	408393
1,4-Difluorobenzene (IS2)	1609720	1606523	1773852	1570016	1771993	1817421
Chlorobenzene-d5 (IS3)	1436269	1437459	1586205	1400955	1581182	1617631

Surrogates (TIC)

1,2-Dichloroethane-d4	1609254	1603178	1763511	1557660	1747893	1788085
o-Toluene-d8	4583566	4557574	5043245	4455345	5014483	5128344
p-Bromofluorobenzene	3615087	3609366	4056000	3584207	4058148	4175322

C5-C8 Aliphatics

	<u>RRFs</u>			<u>RRF<sub>avg</sub></u>		<u>%RSD</u>	
0.5	1	5	25	50	100	1.780	17.44
2.2516	2.0470	1.6738	1.7285	1.5279	1.4481		

C9-C12 Aliphatics

	<u>RRFs</u>			<u>RRF<sub>avg</sub></u>		<u>%RSD</u>	
0.5	1	5	25	50	100	2.525	12.27
2.9198	2.7749	2.4534	2.6251	2.2940	2.0852		

C9-C10 Aromatics

	<u>RRFs</u>			<u>RRF<sub>avg</sub></u>		<u>%RSD</u>	
0.5	1	5	25	50	100	0.479	15.92
0.5644	0.5293	0.4699	0.5194	0.4350	0.3542		

**Massachusetts APH**  
**Continuing Calibration Verification Check Sheet**

Data File Name: 04071203.D  
 Data File Path: J:\MS16\DATA\2012\_04\07\  
 Operator: LH  
 Date Acquired: 4/7/12 1:51  
 Acq. Method File: TO15.M  
 Sample Name: 25ng MAPH CCV STD  
 Misc Info: S25-03291201/S25-03161201  
 Instrument Name: GCMS-16

Enter RRFs from current ICAL!

Internal Standards	RT	Area
7) 1,4-Difluorobenzene (IS2)	13.51	1477594
16) Chlorobenzene-d5 (IS3)	17.46	1283723

C5-C8 Aliphatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
3) Isopentane	7.06	1775206	1.463	124.3	-17.82	-30	30	Pass
4) n-Hexane	11.41	1948862						
9) Cyclohexane	13.44	2333306						
10) 2,3-Dimethylpentane	13.71	2322044	Spike	ICAL				
11) n-Heptane	14.58	2209814	Amt (ng)	RRF				
14) n-Octane	16.81	2491474	151.30	1.78				
		<b>13080706</b>						

C9-C12 Aliphatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
18) 2,3-Dimethylheptane	17.83	2690970	2.293	133.4	-9.18	-30	30	Pass
19) n-Nonane	18.42	2602638						
25) n-Decane	19.66	2736378						
28) Butylcyclohexane	20.16	3274636	Spike	ICAL				
29) n-Undecane	20.68	2881643	Amt (ng)	RRF				
30) n-Dodecane	21.58	3111762	146.90	2.525				
		<b>17298027</b>						

C9-C10 Aromatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
22) Isopropylbenzene	18.73	448849	0.439	110.2	-8.28	-30	30	Pass
23) 1-Methyl-3-ethylbenzene	19.19	504134						
24) 1,3,5-Trimethylbenzene	19.28	676024						
26) p-Isopropyltoluene	19.95	395875	Spike	ICAL				
27) 1,2,3-Trimethylbenzene	19.96	684651	Amt (ng)	RRF				
		<b>2709533</b>	120.1	0.479				

in 4/7/12

**Massachusetts APH  
Continuing Calibration Verification Check Sheet**

Data File Name: 04091203.D  
 Data File Path: J:\MS16\DATA\2012\_04\09\  
 Operator: LH  
 Date Acquired: 4/9/12 10:01  
 Acq. Method File: TO15.M  
 Sample Name: 25ng MAPH CCV STD  
 Misc Info: S25-03291201/S25-03161201  
 Instrument Name: GCMS-16

Enter RRFs from current ICAL!

Internal Standards	RT	Area
7) 1,4-Difluorobenzene (IS2)	13.51	1292971
16) Chlorobenzene-d5 (IS3)	17.46	1099977

C5-C8 Aliphatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
3) Isopentane	7.06	1834143	1.684	143.2	-5.38	-30	30	Pass
4) n-Hexane	11.41	1968429						
9) Cyclohexane	13.44	2344177						
10) 2,3-Dimethylpentane	13.71	2319073	Spike	ICAL				
11) n-Heptane	14.58	2219413	Amt (ng)	RRF				
14) n-Octane	16.81	2494694	151.30	1.78				
		<b>13179929</b>						

C9-C12 Aliphatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
18) 2,3-Dimethylheptane	17.83	2684531	2.630	153.0	4.18	-30	30	Pass
19) n-Nonane	18.42	2597249						
25) n-Decane	19.66	2704647						
28) Butylcyclohexane	20.16	3198559	Spike	ICAL				
29) n-Undecane	20.68	2827213	Amt (ng)	RRF				
30) n-Dodecane	21.58	2989675	146.90	2.525				
		<b>17001874</b>						

C9-C10 Aromatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
22) Isopropylbenzene	18.72	436944	0.500	125.4	4.37	-30	30	Pass
23) 1-Methyl-3-ethylbenzene	19.19	489321						
24) 1,3,5-Trimethylbenzene	19.28	657834						
26) p-Isopropyltoluene	19.95	387057	Spike	ICAL				
27) 1,2,3-Trimethylbenzene	19.96	670727	Amt (ng)	RRF				
		<b>2641883</b>	120.1	0.479				

*u 4/9/12*



Response Factor Report GCMS-16

Method Path : J:\MS16\METHODS\  
 Method File : R16022812.M  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Wed Feb 29 09:45:36 2012  
 Response Via : Initial Calibration

Calibration Files

0.1 =02281209.D 0.2 =02281210.D 0.5 =02281211.D 1.0 =02281212.D 5.0 =02281213.D 25 =02281214.D  
 50 =02281215.D 100 =02281216.D

Compound	0.1	0.2	0.5	1.0	5.0	25	50	100	Avg	%RSD
1) IR Bromochloromethane...				ISTD						
2) T Propene	1.266	1.044	0.941	0.879	1.119	1.182	1.117	0.902	1.056	13.21
3) T Dichlorodifluo...	3.487	3.057	2.439	2.686	2.602	2.568	2.404	1.980	2.653	17.05
4) T Chloromethane	2.142	1.754	1.395	1.560	1.422	1.451	1.319	1.025	1.509	21.81
5) T 1,2-Dichloro-1...	1.855	1.616	1.317	1.450	1.448	1.403	1.296	1.040	1.428	16.72
6) T Vinyl Chloride	1.876	1.726	1.422	1.632	1.615	1.581	1.483	1.238	1.572	12.35
7) T 1,3-Butadiene	1.206	1.049	0.873	0.989	1.035	1.064	1.043	0.883	1.018	10.46
8) T Bromomethane	1.637	1.323	1.022	1.148	1.114	1.146	1.072	0.874	1.167	19.56
9) T Chloroethane	0.941	0.886	0.717	0.799	0.790	0.774	0.724	0.612	0.780	13.07
10) T Ethanol			0.670	0.746	0.735	0.713	0.684	0.589	0.690	8.31
11) T Acetonitrile	2.188	1.812	1.379	1.549	1.515	1.466	1.406	1.238	1.569	19.11
12) T Acrolein	0.816	0.658	0.515	0.568	0.544	0.537	0.522	0.458	0.577	19.38
13) T Acetone			0.698	0.754	0.700	0.674	0.636	0.523	0.664	11.91
14) T Trichlorofluor...	3.048	2.709	2.175	2.469	2.437	2.462	2.331	1.951	2.448	13.50
15) T 2-Propanol (Is...			2.216	2.820	2.081	1.852	1.780	1.573	2.054	21.34
16) T Acrylonitrile	1.076	1.051	0.966	1.110	1.171	1.184	1.148	0.999	1.088	7.35
17) T 1,1-Dichloroet...	1.332	1.160	0.976	1.094	1.109	1.121	1.063	0.906	1.095	11.55
18) T 2-Methyl-2-Pro...		3.293	2.593	2.892	2.865	2.031			2.735	17.04
19) T Methylene Chlo...			1.380	1.317	1.160	1.135	1.071	0.909	1.162	14.64
20) T 3-Chloro-1-pro...	1.417	1.185	1.036	1.175	1.253	1.294	1.263	1.100	1.215	9.77
21) T Trichlorotrifl...	1.439	1.322	1.027	1.156	1.165	1.178	1.112	0.924	1.166	13.76
22) T Carbon Disulfide	5.124	4.418	3.529	3.908	3.928	3.956	3.708	3.097	3.958	15.28
23) T trans-1,2-Dich...	1.818	1.612	1.347	1.547	1.597	1.611	1.524	1.278	1.542	10.88
24) T 1,1-Dichloroet...	2.378	2.110	1.742	1.924	1.931	1.927	1.813	1.526	1.919	13.13
25) T Methyl tert-Bu...	4.447	3.935	3.104	3.445	3.441	3.473	3.268	2.702	3.477	15.13
26) T Vinyl Acetate	0.246	0.233	0.205	0.241	0.261	0.265	0.242	0.186	0.235	11.45
27) T 2-Butanone (MEK)	0.685	0.671	0.601	0.699	0.712	0.720	0.636	0.405	0.641	16.09
28) T cis-1,2-Dichlo...	1.820	1.609	1.313	1.493	1.519	1.526	1.434	1.202	1.489	12.50
29) T Diisopropyl Ether			0.842	0.922	0.913	0.897	0.788	0.579	0.823	15.77
30) T Ethyl Acetate	0.365	0.367	0.331	0.379	0.392	0.377	0.344	0.258	0.352	12.15
31) T n-Hexane	2.337	1.964	1.576	1.715	1.708	1.647	1.492	1.103	1.693	21.11

Method Path : J:\MS16\METHODS\  
 Method File : R16022812.M

Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

Title	2.564	2.289	1.789	2.029	2.039	2.036	1.927	1.626	2.037	14.18
32) T Chloroform	2.564	2.289	1.789	2.029	2.039	2.036	1.927	1.626	2.037	14.18
33) S 1,2-Dichloroet...	1.559	1.561	1.555	1.556	1.534	1.497	1.489	1.489	1.530	2.15
34) T Tetrahydrofura...	0.802	0.750	0.715	0.655	0.625	0.532	0.680	0.680	0.680	14.20
35) T Ethyl tert-But...	1.672	1.566	1.239	1.401	1.406	1.438	1.357	1.125	1.401	12.26
36) T 1,2-Dichloroet...	2.074	1.794	1.470	1.635	1.625	1.623	1.528	1.272	1.628	14.49
37) IR 1,4-Difluorobenzen...	-----ISTD-----									
38) T 1,1,1-Trichlor...	0.581	0.492	0.403	0.447	0.449	0.471	0.449	0.370	0.458	13.70
39) T Isopropyl Acetate	0.183	0.154	0.129	0.150	0.152	0.151	0.141	0.115	0.147	13.39
40) T 1-Butanol	0.299	0.258	0.202	0.232	0.245	0.247	0.231	0.192	0.238	13.98
41) T Benzene	1.522	1.231	0.947	1.003	0.976	0.988	0.923	0.745	1.042	22.55
42) T Carbon Tetrach...	0.398	0.357	0.293	0.346	0.371	0.406	0.388	0.320	0.360	10.87
43) T Cyclohexane	0.515	0.449	0.358	0.398	0.398	0.399	0.360	0.275	0.394	17.79
44) T tert-Amyl Meth...	0.915	0.808	0.654	0.741	0.751	0.769	0.726	0.595	0.745	12.89
45) T 1,2-Dichloropr...	0.311	0.273	0.213	0.249	0.245	0.248	0.233	0.194	0.246	14.49
46) T Bromodichlorom...	0.439	0.377	0.317	0.363	0.375	0.390	0.370	0.301	0.367	11.68
47) T Trichloroethene	0.418	0.370	0.287	0.327	0.332	0.345	0.324	0.260	0.333	14.55
48) T 1,4-Dioxane	0.271	0.230	0.189	0.217	0.218	0.222	0.208	0.165	0.215	14.27
49) T 2,2,4-Trimethy...	1.281	1.127	0.902	0.998	0.991	0.994	0.922	0.755	0.996	15.69
50) T Methyl Methacr...	0.118	0.109	0.093	0.109	0.114	0.118	0.109	0.085	0.107	11.07
51) T n-Heptane	0.318	0.261	0.222	0.247	0.248	0.252	0.237	0.194	0.247	14.31
52) T cis-1,3-Dichlo...	0.438	0.403	0.342	0.399	0.426	0.448	0.421	0.344	0.403	10.01
53) T 4-Methyl-2-pen...	0.219	0.212	0.178	0.211	0.217	0.224	0.212	0.177	0.206	8.91
54) T trans-1,3-Dich...	0.335	0.348	0.292	0.358	0.392	0.424	0.402	0.331	0.360	12.05
55) T 1,1,2-Trichlor...	0.317	0.291	0.238	0.266	0.271	0.282	0.265	0.214	0.268	11.76
56) IR Chlorobenzene-d5 (...)	-----ISTD-----									
57) S Toluene-d8 (SS2)	2.333	2.340	2.320	2.313	2.322	2.300	2.301	2.268	2.312	0.98
58) T Toluene	3.569	3.002	2.316	2.553	2.544	2.520	2.294	1.715	2.564	21.15
59) T 2-Hexanone	1.402	1.216	0.979	1.077	1.077	1.095	1.034	0.845	1.091	15.06
60) T Dibromochlorom...	0.834	0.767	0.621	0.711	0.777	0.829	0.788	0.617	0.743	11.54
61) T 1,2-Dibromoethane	0.826	0.751	0.623	0.717	0.754	0.779	0.734	0.586	0.721	11.04
62) T n-Butyl Acetate	1.568	1.351	1.073	1.255	1.292	1.317	1.279	1.074	1.276	12.40
63) T n-Octane	0.652	0.537	0.418	0.465	0.468	0.473	0.444	0.357	0.477	18.31
64) T Tetrachloroethene	1.126	0.969	0.792	0.889	0.896	0.927	0.860	0.643	0.888	15.64
65) T Chlorobenzene	2.298	2.009	1.568	1.778	1.800	1.798	1.644	1.213	1.764	17.98
66) T Ethylbenzene	3.777	3.256	2.554	2.857	2.923	2.925	2.657	1.940	2.861	18.62
67) T m- & p-Xylenes	2.941	2.593	2.011	2.269	2.337	2.335	2.078	1.453	2.252	19.37
68) T Bromoform	0.747	0.659	0.566	0.659	0.754	0.824	0.778	0.607	0.699	12.84
69) T Styrene	2.113	1.852	1.506	1.723	1.841	1.896	1.727	1.280	1.742	14.60
70) T o-Xylene	3.164	2.681	2.123	2.394	2.454	2.456	2.202	1.559	2.379	19.38

Method Path	J:\MS16\METHODS\	Method File	R16022812.M	Title	EPA TO-15 per SOP	VOA-TO15	(CASS TO-15/GC-MS)	1.425	1.188	0.937	1.041	1.046	1.048	0.976	0.776	1.055	18.04
71) T	n-Nonane	1.385	1.241	0.984	1.136	1.165	1.163	1.056	0.770	1.113	16.41						
72) T	1,1,2,2-Tetrac...	0.935	0.940	0.961	0.955	0.950	0.961	0.957	0.968	0.953	1.18						
73) S	Bromofluoroben...	4.019	3.797	2.823	3.129	3.225	3.210	2.836	1.993	3.129	19.94						
74) T	Cumene	1.863	1.686	1.306	1.478	1.555	1.591	1.452	1.072	1.500	15.96						
75) T	alpha-Pinene	4.604	4.025	3.229	3.671	3.814	3.776	3.339	2.351	3.601	18.29						
76) T	n-Propylbenzene	3.610	3.089	2.518	2.900	3.088	3.218	2.774	2.033	2.904	16.41						
77) T	3-Ethyltoluene	3.148	3.062	2.470	2.820	2.972	2.786	2.522	1.670	2.681	17.67						
78) T	4-Ethyltoluene	3.111	2.802	2.201	2.446	2.549	2.548	2.267	1.615	2.442	18.14						
79) T	1,3,5-Trimethy...	1.550	1.405	1.154	1.348	1.491	1.524	1.364	0.979	1.352	14.53						
80) T	alpha-Methylst...	4.097	3.450	2.686	3.061	3.171	3.135	2.778	1.959	3.042	20.32						
81) T	2-Ethyltoluene	3.214	2.749	2.134	2.476	2.623	2.486	2.032	1.241	2.369	24.66						
82) T	1,2,4-Trimethy...	1.406	1.278	1.042	1.188	1.224	1.254	1.159	0.888	1.180	13.32						
83) T	n-Decane	2.172	1.840	1.536	1.878	2.188	2.375	2.197	1.600	1.973	15.49						
84) T	Benzyl Chloride	1.923	1.688	1.347	1.533	1.630	1.658	1.498	1.047	1.540	16.86						
85) T	1,3-Dichlorobe...	2.134	1.757	1.415	1.594	1.669	1.700	1.514	1.035	1.602	19.53						
86) T	1,4-Dichlorobe...	4.029	3.567	2.848	3.283	3.442	3.412	2.992	2.016	3.198	18.69						
87) T	sec-Butylbenzene	3.919	3.433	2.765	3.224	3.443	3.295	2.711	1.683	3.059	22.14						
88) T	4-Isopropyltol...	3.053	2.720	2.180	2.524	2.699	2.640	2.243	1.459	2.440	19.82						
89) T	1,2,3-Trimethy...	1.856	1.657	1.322	1.504	1.607	1.623	1.439	1.004	1.501	17.06						
90) T	1,2-Dichlorobe...	0.925	0.877	0.732	0.847	0.906	0.925	0.846	0.619	0.835	12.88						
91) T	d-Limonene	0.601	0.532	0.442	0.530	0.612	0.651	0.615	0.473	0.557	13.38						
92) T	1,2-Dibromo-3-...	1.397	1.229	0.991	1.163	1.220	1.246	1.151	0.874	1.159	13.94						
93) T	n-Undecane	1.556	1.297	1.021	1.160	1.263	1.337	1.235	0.893	1.220	16.52						
94) T	1,2,4-Trichlor...	6.159	4.462	3.447	3.848	4.216	4.297	3.885	2.708	4.128	24.04						
95) T	Naphthalene	1.391	1.260	1.019	1.160	1.283	1.280	1.205	0.936	1.192	12.55						
96) T	n-Dodecane	1.077	0.859	0.668	0.727	0.782	0.823	0.776	0.563	0.784	19.20						
97) T	Hexachlorobuta...	1.137	0.944	0.713	0.826	0.829	0.837	0.796	0.654	0.842	17.52						
98) T	Cyclohexanone	3.729	3.267	2.630	3.067	3.276	3.135	2.580	1.602	2.911	22.14						
99) T	tert-Butylbenzene	2.975	2.669	2.172	2.522	2.655	2.653	2.339	1.639	2.453	16.60						
100) T	n-Butylbenzene																

(#) = Out of Range

Evaluate Continuing Calibration Report

Data Path : J:\MS16\DATA\2012\_04\07\  
 Data File : 04071202.D  
 Acq On : 7 Apr 2012 1:17  
 Operator : LH  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03291201/S25-03191203  
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Apr 07 07:29:46 2012  
 Quant Method : J:\MS16\METHODS\R16022812.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Wed Feb 29 09:45:36 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
1	IR Bromochloromethane (IS1)	1.000	1.000	0.0	80	-0.02
2	T Propene	1.056	1.099	-4.1	75	-0.01
3	T Dichlorodifluoromethane (CF	2.653	2.078	21.7	65	0.00
4	T Chloromethane	1.509	1.495	0.9	83	-0.02
5	T 1,2-Dichloro-1,1,2,2-tetra	1.428	1.186	16.9	68	-0.02
6	T Vinyl Chloride	1.572	1.487	5.4	76	-0.02
7	T 1,3-Butadiene	1.018	0.994	2.4	75	-0.02
8	T Bromomethane	1.167	1.021	12.5	72	-0.03
9	T Chloroethane	0.780	0.755	3.2	78	-0.02
10	T Ethanol	0.690	0.702	-1.7	79	0.00
11	T Acetonitrile	1.569	1.445	7.9	79	-0.06
12	T Acrolein	0.577	0.533	7.6	80	-0.03
13	T Acetone	0.664	0.606	8.7	72	-0.06
14	T Trichlorofluoromethane	2.448	2.006	18.1	65	-0.02
15	T 2-Propanol (Isopropanol)	2.054	1.727	15.9	75	-0.06
16	T Acrylonitrile	1.088	1.160	-6.6	79	-0.04
17	T 1,1-Dichloroethene	1.095	1.024	6.5	73	-0.02
18	T 2-Methyl-2-Propanol (tert-B	2.735	1.397	48.9#	55	-0.04
19	T Methylene Chloride	1.162	1.045	10.1	74	-0.02
20	T 3-Chloro-1-propene (Allyl C	1.215	1.238	-1.9	77	-0.02
21	T Trichlorotrifluoroethane	1.166	1.061	9.0	72	-0.02
22	T Carbon Disulfide	3.958	3.758	5.1	76	-0.02
23	T trans-1,2-Dichloroethene	1.542	1.497	2.9	75	-0.02
24	T 1,1-Dichloroethane	1.919	1.813	5.5	76	-0.02
25	T Methyl tert-Butyl Ether	3.477	2.841	18.3	66	-0.02
26	T Vinyl Acetate	0.235	0.217	7.7	66	-0.04
27	T 2-Butanone (MEK)	0.641	0.640	0.2	71	-0.03
28	T cis-1,2-Dichloroethene	1.489	1.377	7.5	72	-0.02
29	T Diisopropyl Ether	0.823	0.747	9.2	67	-0.01
30	T Ethyl Acetate	0.352	0.327	7.1	70	-0.03
31	T n-Hexane	1.693	1.394	17.7	68	0.00
32	T Chloroform	2.037	1.776	12.8	70	-0.03
33	S 1,2-Dichloroethane-d4 (SS1)	1.530	1.360	11.1	73	-0.02
34	T Tetrahydrofuran (THF)	0.680	0.639	6.0	78	-0.02
35	T Ethyl tert-Butyl Ether	1.401	1.195	14.7	67	-0.02
36	T 1,2-Dichloroethane	1.628	1.310	19.5	65	-0.02
37	IR 1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	81	-0.01
38	T 1,1,1-Trichloroethane	0.458	0.394	14.0	68	-0.01

Evaluate Continuing Calibration Report

Data Path : J:\MS16\DATA\2012\_04\07\  
 Data File : 04071202.D  
 Acq On : 7 Apr 2012 1:17  
 Operator : LH  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03291201/S25-03191203  
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Apr 07 07:29:46 2012  
 Quant Method : J:\MS16\METHODS\R16022812.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Wed Feb 29 09:45:36 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
39 T	Isopropyl Acetate	0.147	0.134	8.8	72	-0.02
40 T	1-Butanol	0.238	0.218	8.4	71	-0.04
41 T	Benzene	1.042	0.898	13.8	74	-0.01
42 T	Carbon Tetrachloride	0.360	0.335	6.9	67	-0.01
43 T	Cyclohexane	0.394	0.348	11.7	71	-0.02
44 T	tert-Amyl Methyl Ether	0.745	0.649	12.9	68	-0.01
45 T	1,2-Dichloropropane	0.246	0.235	4.5	77	-0.01
46 T	Bromodichloromethane	0.367	0.327	10.9	68	-0.01
47 T	Trichloroethene	0.333	0.306	8.1	72	-0.01
48 T	1,4-Dioxane	0.215	0.204	5.1	74	-0.02
49 T	2,2,4-Trimethylpentane (Iso	0.996	0.957	3.9	78	-0.01
50 T	Methyl Methacrylate	0.107	0.105	1.9	72	-0.02
51 T	n-Heptane	0.247	0.230	6.9	74	-0.01
52 T	cis-1,3-Dichloropropene	0.403	0.399	1.0	72	-0.01
53 T	4-Methyl-2-pentanone	0.206	0.208	-1.0	75	-0.02
54 T	trans-1,3-Dichloropropene	0.360	0.368	-2.2	70	0.00
55 T	1,1,2-Trichloroethane	0.268	0.255	4.9	73	0.00
56 IR	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	83	0.00
57 S	Toluene-d8 (SS2)	2.312	2.259	2.3	82	0.00
58 T	Toluene	2.564	2.191	14.5	72	-0.01
59 T	2-Hexanone	1.091	0.900	17.5	68	-0.01
60 T	Dibromochloromethane	0.743	0.678	8.7	68	0.00
61 T	1,2-Dibromoethane	0.721	0.660	8.5	70	0.00
62 T	n-Butyl Acetate	1.276	1.134	11.1	72	-0.01
63 T	n-Octane	0.477	0.423	11.3	74	-0.01
64 T	Tetrachloroethene	0.888	0.808	9.0	72	0.00
65 T	Chlorobenzene	1.764	1.550	12.1	72	0.00
66 T	Ethylbenzene	2.861	2.493	12.9	71	0.00
67 T	m- & p-Xylenes	2.252	1.921	14.7	68	-0.01
68 T	Bromoform	0.699	0.675	3.4	68	0.00
69 T	Styrene	1.742	1.579	9.4	69	0.00
70 T	o-Xylene	2.379	2.042	14.2	69	-0.01
71 T	n-Nonane	1.055	0.887	15.9	70	0.00
72 T	1,1,2,2-Tetrachloroethane	1.113	0.964	13.4	69	-0.01
73 S	Bromofluorobenzene (SS3)	0.953	0.991	-4.0	86	0.00
74 T	Cumene	3.129	2.704	13.6	70	-0.01
75 T	alpha-Pinene	1.500	1.316	12.3	69	0.00
76 T	n-Propylbenzene	3.601	3.111	13.6	68	0.00

Evaluate Continuing Calibration Report

Data Path : J:\MS16\DATA\2012\_04\07\  
 Data File : 04071202.D  
 Acq On : 7 Apr 2012 1:17  
 Operator : LH  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03291201/S25-03191203  
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Apr 07 07:29:46 2012  
 Quant Method : J:\MS16\METHODS\R16022812.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Wed Feb 29 09:45:36 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
77 T	3-Ethyltoluene	2.904	2.489	14.3	64	-0.01
78 T	4-Ethyltoluene	2.681	2.428	9.4	72	0.00
79 T	1,3,5-Trimethylbenzene	2.442	2.098	14.1	68	0.00
80 T	alpha-Methylstyrene	1.352	1.189	12.1	65	0.00
81 T	2-Ethyltoluene	3.042	2.529	16.9	67	-0.01
82 T	1,2,4-Trimethylbenzene	2.369	1.962	17.2	66	-0.01
83 T	n-Decane	1.180	1.038	12.0	69	0.00
84 T	Benzyl Chloride	1.973	1.839	6.8	64	-0.01
85 T	1,3-Dichlorobenzene	1.540	1.361	11.6	68	-0.01
86 T	1,4-Dichlorobenzene	1.602	1.399	12.7	68	0.00
87 T	sec-Butylbenzene	3.198	2.783	13.0	68	-0.01
88 T	4-Isopropyltoluene (p-Cymen	3.059	2.675	12.6	67	0.00
89 T	1,2,3-Trimethylbenzene	2.440	2.093	14.2	66	-0.02
90 T	1,2-Dichlorobenzene	1.501	1.315	12.4	67	0.00
91 T	d-Limonene	0.835	0.715	14.4	64	0.00
92 T	1,2-Dibromo-3-Chloropropane	0.557	0.546	2.0	70	0.00
93 T	n-Undecane	1.159	1.108	4.4	74	0.00
94 T	1,2,4-Trichlorobenzene	1.220	1.211	0.7	75	0.00
95 T	Naphthalene	4.128	3.797	8.0	73	0.00
96 T	n-Dodecane	1.192	1.249	-4.8	81	0.00
97 T	Hexachlorobutadiene	0.784	0.768	2.0	77	0.00
98 T	Cyclohexanone	0.842	0.742	11.9	74	-0.01
99 T	tert-Butylbenzene	2.911	2.545	12.6	67	0.00
100 T	n-Butylbenzene	2.453	2.158	12.0	68	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*UH 4/7/12*

Evaluate Continuing Calibration Report

Data Path : J:\MS16\DATA\2012\_04\09\  
 Data File : 04091202.D  
 Acq On : 9 Apr 2012 9:28  
 Operator : LH  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03291201/S25-03191203  
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Apr 09 11:30:10 2012  
 Quant Method : J:\MS16\METHODS\R16022812.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Wed Feb 29 09:45:36 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	IR Bromochloromethane (IS1)	1.000	1.000	0.0	75	-0.01
2	T Propene	1.056	1.192	-12.9	76	0.00
3	T Dichlorodifluoromethane (CF	2.653	2.243	15.5	66	0.00
4	T Chloromethane	1.509	1.657	-9.8	86	-0.02
5	T 1,2-Dichloro-1,1,2,2-tetra	1.428	1.262	11.6	67	-0.02
6	T Vinyl Chloride	1.572	1.636	-4.1	78	-0.02
7	T 1,3-Butadiene	1.018	1.124	-10.4	79	-0.02
8	T Bromomethane	1.167	1.085	7.0	71	-0.02
9	T Chloroethane	0.780	0.847	-8.6	82	-0.02
10	T Ethanol	0.690	0.808	-17.1	85	0.00
11	T Acetonitrile	1.569	1.685	-7.4	86	-0.06
12	T Acrolein	0.577	0.604	-4.7	84	-0.03
13	T Acetone	0.664	0.676	-1.8	75	-0.05
14	T Trichlorofluoromethane	2.448	2.159	11.8	66	-0.01
15	T 2-Propanol (Isopropanol)	2.054	2.187	-6.5	89	-0.06
16	T Acrylonitrile	1.088	1.266	-16.4	80	-0.04
17	T 1,1-Dichloroethene	1.095	1.093	0.2	73	-0.02
18	T 2-Methyl-2-Propanol (tert-B	2.735	1.473	46.1#	54	-0.03
19	T Methylene Chloride	1.162	1.116	4.0	74	-0.02
20	T 3-Chloro-1-propene (Allyl C	1.215	1.424	-17.2	83	-0.02
21	T Trichlorotrifluoroethane	1.166	1.129	3.2	72	-0.02
22	T Carbon Disulfide	3.958	4.050	-2.3	77	-0.02
23	T trans-1,2-Dichloroethene	1.542	1.641	-6.4	76	-0.02
24	T 1,1-Dichloroethane	1.919	1.969	-2.6	77	-0.02
25	T Methyl tert-Butyl Ether	3.477	3.100	10.8	67	-0.01
26	T Vinyl Acetate	0.235	0.234	0.4	66	-0.04
27	T 2-Butanone (MEK)	0.641	0.691	-7.8	72	-0.03
28	T cis-1,2-Dichloroethene	1.489	1.501	-0.8	74	-0.02
29	T Diisopropyl Ether	0.823	0.809	1.7	68	0.00
30	T Ethyl Acetate	0.352	0.358	-1.7	71	-0.03
31	T n-Hexane	1.693	1.556	8.1	71	0.00
32	T Chloroform	2.037	1.892	7.1	70	-0.02
33	S 1,2-Dichloroethane-d4 (SS1)	1.530	1.413	7.6	71	-0.01
34	T Tetrahydrofuran (THF)	0.680	0.694	-2.1	80	-0.01
35	T Ethyl tert-Butyl Ether	1.401	1.282	8.5	67	-0.01
36	T 1,2-Dichloroethane	1.628	1.429	12.2	66	-0.02
37	IR 1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	78	-0.01
38	T 1,1,1-Trichloroethane	0.458	0.408	10.9	67	-0.01

Evaluate Continuing Calibration Report

Data Path : J:\MS16\DATA\2012\_04\09\  
 Data File : 04091202.D  
 Acq On : 9 Apr 2012 9:28  
 Operator : LH  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03291201/S25-03191203  
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Apr 09 11:30:10 2012  
 Quant Method : J:\MS16\METHODS\R16022812.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Wed Feb 29 09:45:36 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
39 T	Isopropyl Acetate	0.147	0.143	2.7	73	-0.02
40 T	1-Butanol	0.238	0.234	1.7	73	-0.04
41 T	Benzene	1.042	0.934	10.4	73	-0.01
42 T	Carbon Tetrachloride	0.360	0.345	4.2	66	-0.01
43 T	Cyclohexane	0.394	0.366	7.1	71	-0.02
44 T	tert-Amyl Methyl Ether	0.745	0.685	8.1	69	-0.01
45 T	1,2-Dichloropropane	0.246	0.250	-1.6	78	-0.01
46 T	Bromodichloromethane	0.367	0.340	7.4	67	-0.01
47 T	Trichloroethene	0.333	0.312	6.3	70	-0.01
48 T	1,4-Dioxane	0.215	0.211	1.9	73	-0.02
49 T	2,2,4-Trimethylpentane (Iso	0.996	1.026	-3.0	80	-0.01
50 T	Methyl Methacrylate	0.107	0.109	-1.9	72	-0.02
51 T	n-Heptane	0.247	0.241	2.4	74	-0.01
52 T	cis-1,3-Dichloropropene	0.403	0.419	-4.0	72	0.00
53 T	4-Methyl-2-pentanone	0.206	0.221	-7.3	76	-0.01
54 T	trans-1,3-Dichloropropene	0.360	0.389	-8.1	71	0.00
55 T	1,1,2-Trichloroethane	0.268	0.260	3.0	71	0.00
56 IR	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	77	0.00
57 S	Toluene-d8 (SS2)	2.312	2.308	0.2	78	0.00
58 T	Toluene	2.564	2.344	8.6	72	-0.01
59 T	2-Hexanone	1.091	0.995	8.8	70	-0.01
60 T	Dibromochloromethane	0.743	0.714	3.9	67	0.00
61 T	1,2-Dibromoethane	0.721	0.689	4.4	69	0.00
62 T	n-Butyl Acetate	1.276	1.264	0.9	74	-0.01
63 T	n-Octane	0.477	0.461	3.4	76	0.00
64 T	Tetrachloroethene	0.888	0.850	4.3	71	0.00
65 T	Chlorobenzene	1.764	1.630	7.6	70	0.00
66 T	Ethylbenzene	2.861	2.650	7.4	70	0.00
67 T	m- & p-Xylenes	2.252	2.058	8.6	68	-0.01
68 T	Bromoform	0.699	0.715	-2.3	67	0.00
69 T	Styrene	1.742	1.684	3.3	69	0.00
70 T	o-Xylene	2.379	2.191	7.9	69	-0.01
71 T	n-Nonane	1.055	0.986	6.5	73	0.00
72 T	1,1,2,2-Tetrachloroethane	1.113	1.025	7.9	68	-0.01
73 S	Bromofluorobenzene (SS3)	0.953	0.959	-0.6	77	0.00
74 T	Cumene	3.129	2.877	8.1	69	0.00
75 T	alpha-Pinene	1.500	1.410	6.0	69	0.00
76 T	n-Propylbenzene	3.601	3.311	8.1	68	0.00



Evaluate Continuing Calibration Report

Data Path : J:\MS16\DATA\2012\_04\09\  
 Data File : 04091202.D  
 Acq On : 9 Apr 2012 9:28  
 Operator : LH  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-03291201/S25-03191203  
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Apr 09 11:30:10 2012  
 Quant Method : J:\MS16\METHODS\R16022812.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Wed Feb 29 09:45:36 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
77 T	3-Ethyltoluene	2.904	2.828	2.6	68	-0.01
78 T	4-Ethyltoluene	2.681	2.428	9.4	68	0.00
79 T	1,3,5-Trimethylbenzene	2.442	2.244	8.1	68	0.00
80 T	alpha-Methylstyrene	1.352	1.263	6.6	64	0.00
81 T	2-Ethyltoluene	3.042	2.698	11.3	67	-0.01
82 T	1,2,4-Trimethylbenzene	2.369	2.102	11.3	65	-0.01
83 T	n-Decane	1.180	1.129	4.3	70	0.00
84 T	Benzyl Chloride	1.973	1.998	-1.3	65	-0.01
85 T	1,3-Dichlorobenzene	1.540	1.445	6.2	68	-0.01
86 T	1,4-Dichlorobenzene	1.602	1.490	7.0	68	0.00
87 T	sec-Butylbenzene	3.198	2.970	7.1	67	-0.01
88 T	4-Isopropyltoluene (p-Cymen	3.059	2.875	6.0	68	0.00
89 T	1,2,3-Trimethylbenzene	2.440	2.243	8.1	66	-0.01
90 T	1,2-Dichlorobenzene	1.501	1.399	6.8	67	0.00
91 T	d-Limonene	0.835	0.778	6.8	65	0.00
92 T	1,2-Dibromo-3-Chloropropane	0.557	0.582	-4.5	69	0.00
93 T	n-Undecane	1.159	1.204	-3.9	75	0.00
94 T	1,2,4-Trichlorobenzene	1.220	1.274	-4.4	74	0.00
95 T	Naphthalene	4.128	4.004	3.0	72	0.00
96 T	n-Dodecane	1.192	1.339	-12.3	81	0.00
97 T	Hexachlorobutadiene	0.784	0.807	-2.9	76	0.00
98 T	Cyclohexanone	0.842	0.813	3.4	75	-0.01
99 T	tert-Butylbenzene	2.911	2.735	6.0	68	0.00
100 T	n-Butylbenzene	2.453	2.299	6.3	67	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

Response Factor Report MS07

Method : J:\MS07\METHODS\X7021712B.M (RTE Integrator)  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
 1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	Avg	%RSD
1) I Bromochloromethan					ISTD						
2) T Dichlorodifluorom	3.784	3.527	3.467	3.822	3.279	2.986	2.875	3.155	2.920	3.313	10.83
3) T Chloromethane		1.470	0.956	1.190	0.904	0.778	0.569	0.897	0.827	0.949	28.81
4) T Vinyl Chloride	2.802	2.749	2.588	3.044	2.597	2.406	2.314	2.588	2.446	2.615	8.57
5) T Bromomethane	1.894	1.496	1.211	1.613	1.335	1.208	1.105	1.433	1.367	1.407	17.11
6) T Chloroethane	1.511	1.416	1.357	1.517	1.295	1.237	1.195	1.312	1.224	1.340	8.92
7) T Acetone				1.548	1.214	1.122	1.080	1.193	1.090	1.208	14.52
8) T Trichlorofluorome	3.051	2.887	2.831	3.145	2.692	2.565	2.447	2.691	2.536	2.761	8.59
9) T 1,1-Dichloroethen	1.591	1.440	1.366	1.547	1.322	1.229	1.214	1.385	1.295	1.376	9.52
10) T Methylene Chlorid		2.134	1.795	1.866	1.536	1.426	1.404	1.494	1.411	1.633	16.43
11) T Trichlorotrifluor	1.408	1.325	1.295	1.376	1.169	1.081	1.062	1.170	1.097	1.220	10.87
12) T trans-1,2-Dichlor	1.972	1.710	1.589	1.762	1.505	1.406	1.393	1.512	1.439	1.588	12.18
13) T 1,1-Dichloroethan	3.265	3.020	3.016	3.383	2.925	2.756	2.681	2.937	2.841	2.980	7.60
14) T Methyl tert-Butyl	5.168	4.740	4.762	5.045	4.335	4.137	4.166	4.690	4.471	4.613	7.90
15) T cis-1,2-Dichloroe	1.885	1.626	1.630	1.776	1.524	1.434	1.388	1.575	1.482	1.591	10.09
16) T Chloroform			3.041	3.257	2.586	2.390	2.280	2.524	2.374	2.636	14.03
17) S 1,2-Dichloroethan	2.000	2.002	2.032	2.018	1.989	1.973	1.960	1.938	1.932	1.983	1.75
18) T 1,2-Dichloroethan	2.582	2.460	2.375	2.604	2.222	2.090	1.988	2.222	2.094	2.293	9.73
19) T 1,1,1-Trichloroet	2.587	2.431	2.340	2.573	2.191	2.059	2.007	2.278	2.150	2.291	9.18
20) T Benzene			7.536	7.534	6.055	5.665	5.395	5.903	5.468	6.222	14.88
21) T Carbon Tetrachlor	2.280	1.842	1.769	1.950	1.676	1.579	1.549	1.771	1.692	1.790	12.41
22) I 1,4-Difluorobenze					ISTD						
23) T 1,2-Dichloropropa	0.417	0.397	0.373	0.409	0.354	0.331	0.319	0.358	0.336	0.366	9.68
24) T Bromodichlorometh	0.491	0.464	0.436	0.489	0.420	0.398	0.394	0.447	0.416	0.439	8.26
25) T Trichloroethene	0.418	0.379	0.349	0.390	0.324	0.305	0.287	0.326	0.307	0.343	12.96
26) T 1,4-Dioxane	0.324	0.290	0.270	0.278	0.242	0.231	0.227	0.258	0.243	0.263	11.99
27) T cis-1,3-Dichlorop	0.567	0.523	0.508	0.540	0.499	0.484	0.488	0.559	0.522	0.521	5.67
28) T trans-1,3-Dichlor	0.457	0.434	0.431	0.444	0.419	0.409	0.417	0.505	0.482	0.444	7.22

(#) = Out of Range ### Number of calibration levels exceeded format ###  
 X7021712B.M Mon Apr 09 08:16:11 2012

Response Factor Report MS07

Method : J:\Ms07\METHODS\X7021712B.M (RTE Integrator)  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
 1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	Avg	%RSD
29) T 1,1,2-Trichloroet	0.315	0.296	0.286	0.305	0.263	0.248	0.242	0.274	0.257	0.276	9.32
30) S Toluene-d8 (SS2)	1.053	1.071	1.054	1.047	1.051	1.052	1.071	1.060	1.049	1.056	0.84
31) T Toluene	1.782	1.541	1.420	1.485	1.269	1.198	1.177	1.307	1.186	1.374	14.76
32) T 1,2-Dibromoethane	0.379	0.362	0.345	0.362	0.323	0.307	0.304	0.351	0.331	0.340	7.65
33) T Tetrachloroethene	0.423	0.394	0.371	0.400	0.341	0.320	0.311	0.347	0.326	0.359	11.03
-----ISTD-----											
34) I Chlorobenzene-d5											
35) T Chlorobenzene	4.581	4.172	4.017	4.308	3.688	3.469	3.311	3.705	3.345	3.844	11.68
36) T Ethylbenzene	7.604	7.082	6.707	7.235	6.297	6.070	5.915	6.682	5.878	6.608	9.30
37) T m,p-Xylene	5.861	5.347	5.132	5.536	4.930	4.777	4.719	5.268	4.558	5.125	8.24
38) T o-Xylene	6.505	5.826	5.526	5.889	5.188	4.995	4.950	5.594	4.932	5.490	9.67
39) T 1,1,2,2-Tetrachlo	2.655	2.674	2.710	2.646	2.437	2.371	2.298	2.761	2.498	2.561	6.40
40) S Bromofluorobenzene	1.712	1.701	1.695	1.697	1.683	1.680	1.670	1.642	1.593	1.675	2.20
41) T 1,3-Dichlorobenzene	3.304	3.261	3.222	3.204	2.835	2.719	2.608	3.138	2.846	3.015	8.71
42) T 1,4-Dichlorobenzene	3.470	3.303	3.208	3.208	2.803	2.671	2.579	3.174	2.891	3.034	10.13
43) T 1,2-Dichlorobenzene	3.271	3.123	3.112	3.026	2.662	2.536	2.418	3.053	2.772	2.886	10.31
44) T 1,2,4-Trichlorobene	2.184	2.118	2.169	1.753	1.641	1.608	1.575	2.139	1.978	1.907	13.66
45) T Naphthalene	6.857	6.355	6.704	5.143	5.107	5.172	5.297	8.126	7.328	6.232	17.82
46) T Hexachlorobutadie	1.486	1.423	1.403	1.217	1.055	1.013	0.975	1.227	1.145	1.216	15.40

(#) = Out of Range ### Number of calibration levels exceeded format ###  
 X7021712B.M Mon Apr 09 08:16:11 2012

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_04\09\  
 Data File : 04091203.D  
 Acq On : 9 Apr 2012 7:01 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00163  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Apr 09 08:15:00 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane (IS1)	1.000	1.000	0.0	82	0.00
2 T	Dichlorodifluoromethane (CF)	3.313	2.539	23.4	63	0.00
3 T	Chloromethane	0.949	0.738	22.2	67	0.00
4 T	Vinyl Chloride	2.615	2.065	21.0	65	0.00
5 T	Bromomethane	1.407	1.190	15.4	73	0.00
6 T	Chloroethane	1.340	1.065	20.5	67	0.00
7 T	Acetone	1.208	1.043	13.7	70	0.00
8 T	Trichlorofluoromethane	2.761	2.169	21.4	66	0.00
9 T	1,1-Dichloroethene	1.376	1.147	16.6	71	0.00
10 T	Methylene Chloride	1.633	1.261	22.8	67	0.00
11 T	Trichlorotrifluoroethane	1.220	1.046	14.3	73	0.00
12 T	trans-1,2-Dichloroethene	1.588	1.249	21.3	68	0.00
13 T	1,1-Dichloroethane	2.980	2.362	20.7	66	0.00
14 T	Methyl tert-Butyl Ether	4.613	3.609	21.8	68	0.00
15 T	cis-1,2-Dichloroethene	1.591	1.330	16.4	71	0.00
16 T	Chloroform	2.636	2.237	15.1	71	0.00
17 S	1,2-Dichloroethane-d4 (SS1)	1.983	1.917	3.3	79	0.00
18 T	1,2-Dichloroethane	2.293	1.706	25.6	63	0.00
19 T	1,1,1-Trichloroethane	2.291	1.879	18.0	70	0.00
20 T	Benzene	6.222	5.361	13.8	72	0.00
21 T	Carbon Tetrachloride	1.790	1.425	20.4	69	0.00
22 I	1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	88	0.00
23 T	1,2-Dichloropropane	0.366	0.268	26.8	67	0.00
24 T	Bromodichloromethane	0.439	0.322	26.7	68	0.00
25 T	Trichloroethene	0.343	0.255	25.7	70	0.00
26 T	1,4-Dioxane	0.263	0.198	24.7	73	0.00
27 T	cis-1,3-Dichloropropene	0.521	0.396	24.0	70	0.00
28 T	trans-1,3-Dichloropropene	0.444	0.343	22.7	72	0.00
29 T	1,1,2-Trichloroethane	0.276	0.207	25.0	70	0.00
30 S	Toluene-d8 (SS2)	1.056	1.179	-11.6	99	0.00
31 T	Toluene	1.374	1.051	23.5	73	0.00
32 T	1,2-Dibromoethane	0.340	0.264	22.4	72	0.00
33 T	Tetrachloroethene	0.359	0.287	20.1	75	0.00
34 I	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	85	0.00
35 T	Chlorobenzene	3.844	3.219	16.3	75	0.00
36 T	Ethylbenzene	6.608	5.416	18.0	74	0.00
37 T	m,p-Xylene	5.125	4.308	15.9	75	0.00

*Handwritten signature/initials*

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_04\09\  
 Data File : 04091203.D  
 Acq On : 9 Apr 2012 7:01 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00163  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Apr 09 08:15:00 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
38 T	o-Xylene	5.490	4.561	16.9	75	0.00
39 T	1,1,2,2-Tetrachloroethane	2.561	2.098	18.1	74	0.00
40 S	Bromofluorobenzene (SS3)	1.675	1.861	-11.1	95	0.00
41 T	1,3-Dichlorobenzene	3.015	2.704	10.3	82	0.00
42 T	1,4-Dichlorobenzene	3.034	2.684	11.5	82	0.00
43 T	1,2-Dichlorobenzene	2.886	2.603	9.8	84	0.00
44 T	1,2,4-Trichlorobenzene	1.907	1.808	5.2	94	0.00
45 T	Naphthalene	6.232	6.283	-0.8	105	0.00
46 T	Hexachlorobutadiene	1.216	1.083	10.9	88	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*KRA/9/12*

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_04\10\  
 Data File : 04101203.D  
 Acq On : 10 Apr 2012 9:21 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00163  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Apr 10 10:05:18 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane (IS1)	1.000	1.000	0.0	78	0.00
2 T	Dichlorodifluoromethane (CF)	3.313	2.565	22.6	61	0.00
3 T	Chloromethane	0.949	0.735	22.6	63	0.00
4 T	Vinyl Chloride	2.615	2.063	21.1	62	0.00
5 T	Bromomethane	1.407	1.209	14.1	71	0.00
6 T	Chloroethane	1.340	1.091	18.6	66	0.00
7 T	Acetone	1.208	1.070	11.4	69	0.00
8 T	Trichlorofluoromethane	2.761	2.217	19.7	64	0.00
9 T	1,1-Dichloroethene	1.376	1.163	15.5	69	0.00
10 T	Methylene Chloride	1.633	1.296	20.6	66	0.00
11 T	Trichlorotrifluoroethane	1.220	1.078	11.6	72	0.00
12 T	trans-1,2-Dichloroethene	1.588	1.296	18.4	67	0.00
13 T	1,1-Dichloroethane	2.980	2.472	17.0	66	0.00
14 T	Methyl tert-Butyl Ether	4.613	3.623	21.5	65	0.00
15 T	cis-1,2-Dichloroethene	1.591	1.314	17.4	67	0.00
16 T	Chloroform	2.636	2.183	17.2	66	0.00
17 S	1,2-Dichloroethane-d4 (SS1)	1.983	1.970	0.7	77	0.00
18 T	1,2-Dichloroethane	2.293	1.741	24.1	61	0.00
19 T	1,1,1-Trichloroethane	2.291	1.853	19.1	66	0.00
20 T	Benzene	6.222	5.270	15.3	68	0.00
21 T	Carbon Tetrachloride	1.790	1.442	19.4	67	0.00
22 I	1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	88	0.00
23 T	1,2-Dichloropropane	0.366	0.260	29.0	65	0.00
24 T	Bromodichloromethane	0.439	0.306	30.3#	64	0.00
25 T	Trichloroethene	0.343	0.253	26.2	69	0.00
26 T	1,4-Dioxane	0.263	0.199	24.3	73	0.00
27 T	cis-1,3-Dichloropropene	0.521	0.396	24.0	70	0.00
28 T	trans-1,3-Dichloropropene	0.444	0.339	23.6	71	0.00
29 T	1,1,2-Trichloroethane	0.276	0.200	27.5	67	0.00
30 S	Toluene-d8 (SS2)	1.056	1.154	-9.3	97	0.00
31 T	Toluene	1.374	1.027	25.3	72	0.00
32 T	1,2-Dibromoethane	0.340	0.257	24.4	70	0.00
33 T	Tetrachloroethene	0.359	0.284	20.9	74	0.00
34 I	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	84	0.00
35 T	Chlorobenzene	3.844	3.248	15.5	74	0.00
36 T	Ethylbenzene	6.608	5.549	16.0	74	0.00
37 T	m,p-Xylene	5.125	4.350	15.1	74	0.00

*KR 4/10/12*

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_04\10\  
 Data File : 04101203.D  
 Acq On : 10 Apr 2012 9:21 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00163  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Apr 10 10:05:18 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
38 T	o-Xylene	5.490	4.592	16.4	74	0.00
39 T	1,1,2,2-Tetrachloroethane	2.561	2.087	18.5	72	0.00
40 S	Bromofluorobenzene (SS3)	1.675	1.858	-10.9	93	0.00
41 T	1,3-Dichlorobenzene	3.015	2.719	9.8	80	0.00
42 T	1,4-Dichlorobenzene	3.034	2.710	10.7	81	0.00
43 T	1,2-Dichlorobenzene	2.886	2.620	9.2	83	0.00
44 T	1,2,4-Trichlorobenzene	1.907	1.807	5.2	92	0.00
45 T	Naphthalene	6.232	6.463	-3.7	106	0.00
46 T	Hexachlorobutadiene	1.216	1.083	10.9	86	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

## LABORATORY REPORT

April 18, 2012

Christopher Cote  
Montana DEQ  
1100 N. Last Chance Gulch  
Helena, MT 59601

### **RE: Background Indoor Air Study**

Dear Christopher:

Enclosed are the results of the samples submitted to our laboratory on April 4, 2012. For your reference, these analyses have been assigned our service request number P1201320.

All analyses were performed according to our laboratory's NELAP and DoD-ELAP-approved quality assurance program. The test results meet requirements of the current NELAP and DoD-ELAP standards, where applicable, and except as noted in the laboratory case narrative provided. For a specific list of NELAP and DoD-ELAP-accredited analytes, refer to the certifications section at [www.caslab.com](http://www.caslab.com). Results are intended to be considered in their entirety and apply only to the samples analyzed and reported herein.

Columbia Analytical Services, Inc. is certified by the California Department of Health Services, NELAP Laboratory Certificate No. 02115CA; Arizona Department of Health Services, Certificate No. AZ0694; Florida Department of Health, NELAP Certification E871020; New Jersey Department of Environmental Protection, NELAP Laboratory Certification ID #CA009; New York State Department of Health, NELAP NY Lab ID No: 11221; Oregon Environmental Laboratory Accreditation Program, NELAP ID: CA20007; The American Industrial Hygiene Association, Laboratory #101661; United States Department of Defense Environmental Laboratory Accreditation Program (DoD-ELAP), Certificate No. L11-203; Pennsylvania Registration No. 68-03307; TX Commission of Environmental Quality, NELAP ID T104704413-11-2; Minnesota Department of Health, NELAP Certificate No. 362188; Washington State Department of Ecology, ELAP Lab ID: C946, State of Utah Department of Health, NELAP Certificate No. CA015272011-1; Los Angeles Department of Building and Safety, Approval No: TA00001. Each of the certifications listed above have an explicit Scope of Accreditation that applies to specific matrices/methods/analytes; therefore, please contact me for information corresponding to a particular certification.

If you have any questions, please call me at (805) 526-7161.

Respectfully submitted,

**ALS | Environmental**

Kate Aguilera  
Project Manager



Client: Montana DEQ  
Project: Background Indoor Air Study

Service Request No: P1201320

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## CASE NARRATIVE

The samples were received intact under chain of custody on April 4, 2012 and were stored in accordance with the analytical method requirements. Please refer to the sample acceptance check form for additional information. The results reported herein are applicable only to the condition of the samples at the time of sample receipt.

### Air-Phase Petroleum Hydrocarbons (APH) Analysis

The samples were analyzed for total aliphatic and aromatic gasoline range hydrocarbons by gas chromatography/mass spectrometry according to the Method for the Determination of Air-Phase Petroleum Hydrocarbons (APH), Massachusetts Department of Environmental Protection, Revision 1, December, 2009.

### Volatile Organic Compound Analysis

The samples were also analyzed in scan and SIM mode for selected volatile organic compounds in accordance with EPA Method TO-15 from the Compendium of Methods for the Determination of Toxic Organic Compounds in Ambient Air, Second Edition (EPA/625/R-96/010b), January, 1999. The analytical system was comprised of a gas chromatograph / mass spectrometer (GC/MS) interfaced to a whole-air preconcentrator.

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*The results of analyses are given in the attached laboratory report. All results are intended to be considered in their entirety, and Columbia Analytical Services, Inc. (CAS) is not responsible for utilization of less than the complete report.*

*Use of Columbia Analytical Services, Inc. (CAS) Name. Client shall not use CAS's name or trademark in any marketing or reporting materials, press releases or in any other manner ("Materials") whatsoever and shall not attribute to CAS any test result, tolerance or specification derived from CAS's data ("Attribution") without CAS's prior written consent, which may be withheld by CAS for any reason in its sole discretion. To request CAS's consent, Client shall provide copies of the proposed Materials or Attribution and describe in writing Client's proposed use of such Materials or Attribution. If CAS has not provided written approval of the Materials or Attribution within ten (10) days of receipt from Client, Client's request to use CAS's name or trademark in any Materials or Attribution shall be deemed denied. CAS may, in its discretion, reasonably charge Client for its time in reviewing Materials or Attribution requests. Client acknowledges and agrees that the unauthorized use of CAS's name or trademark may cause CAS to incur irreparable harm for which the recovery of money damages will be inadequate. Accordingly, Client acknowledges and agrees that a violation shall justify preliminary injunctive relief. For questions contact the laboratory.*

DETAIL SUMMARY REPORT

Client: Montana DEQ  
 Project ID: Background Indoor Air Study

Service Request: P1201320

Date Received: 4/4/2012  
 Time Received: 09:50

Client Sample ID	Lab Code	Matrix	Date Collected	Time Collected	Container ID	Pi1 (psig)	Pf1 (psig)	MA APH 1.0 - MA VOC PH Can	TO-15 - VOC Cans	TO-15 - VOC SIM
31-1F-032412	P1201320-001	Air	3/25/2012	15:35	AC01571	-1.72	3.69	X	X	X
37-1F-032612	P1201320-002	Air	3/27/2012	08:12	AC01118	-4.87	3.59	X	X	X
29-1F-032612	P1201320-003	Air	3/27/2012	14:45	AC01127	-2.68	3.65	X	X	X
DUP-54	P1201320-004	Air	3/27/2012	14:45	AC01749	-5.09	3.72	X	X	X
30-1F-032612	P1201320-005	Air	3/27/2012	16:19	AS00147	-5.30	3.68	X	X	X
33-1B-032712	P1201320-006	Air	3/28/2012	17:52	AC00781	-2.17	3.68	X	X	X
32-1F-032812	P1201320-007	Air	3/29/2012	08:15	AC01142	-2.83	3.72	X	X	X
36-1F-032812	P1201320-008	Air	3/29/2012	11:20	AC01413	-4.51	3.73	X	X	X
34-1F-032912	P1201320-009	Air	3/30/2012	13:36	AC01533	-1.39	3.57	X	X	X
35-1F-032912	P1201320-010	Air	3/30/2012	13:03	AC00943	-4.85	3.66	X	X	X

**Air - Chain of Custody Record & Analytical Service Request**

Requested Turnaround Time in Business Days (Surcharges) please circle  
 1 Day (100%) 2 Day (75%) 3 Day (50%) 4 Day (35%) 5 Day (25%) 10 Day-Standard

CAS Project No. **PR013720**

Company Name & Address (Reporting Information)		Project Name		P.O. # / Billing Information		Analysis Method		Comments e.g. Actual Preservative or specific instructions		
MONTANA DEPARTMENT OF ENVIRONMENTAL QUALITY PO Box 200901 1100 N LAST CHANCE GULCH HELENA MT 59620-0901		BACKGROUND INDOOR AIR STUDY		REFERENCE AS: BACKGROUND INDOOR AIR STUDY BUILT TO CHRIS COTE OF MDEQ		TO-15 TO-15 SIM MA-APH				
Project Manager CHRIS COTE		Project Number		Sampler (Print & Sign) SHANNON GALE & JESSICA GUTTING		CAS Contact:				
Phone (406) 841-5078		Fax		Canister ID (Bar code # - AC, SC, etc.)		Flow Controller ID (Bar code # - FC #)		Sample Volume		
Email Address for Result Reporting ccote2@mt.gov; MARTICHBJ@edm.com		Laboratory ID Number		Date Collected		Time Collected		Canister Start Pressure "Hg	Canister End Pressure "Hg/psig	
31-1F-032412	01-176	03/25/12	1535	AC01571	FCAD00136	-28.5	-9.0	19.5	X	ANALYZE
37-1F-032012	02-484	3/27/12	0812	AC01118	FCAD00526	-28.0	-9.0	19.0	X	PROJECT SPECIFIC
29-1F-032612	02-265	3/27/12	1445	AC01127	FCAD00473	-27.0	-3.5	23.5	X	SPECIFIC
DUP-54	04-514	3/27/12	1445	AC01749	FCAD00483	-29.25	-10.0	19.25	X	LIST OF COMPOUNDS
30-1F-032612	05-531	3/27/12	1019	AC00147	FCAD00490	-28.0	-9.5	18.5	X	TO PROJECT
33-1B-032712	06-216	3/28/12	1752	AC00781	FCAD00050	-28.0	-3.5	24.5	X	SPECIFIC REPORTING LIMITS
32-1F-032812	07-282	3/29/12	0815	AC01142	FCAD00291	-28.0	-5.0	23.0	X	
36-1F-032812	08-452	3/29/12	1120	AC01413	FCAD00133	-29.0	-9.5	19.5	X	
34-1F-032812	09-135	3/30/12	1330	AC01533	FCAD00381	-29.5	-3.5	26.0	X	
35-1F-032912	10-484	3/30/12	1303	AC00943	FCAD00500	-25.5	-10.5	19.0	X	

**Report Tier Levels - please select**  
 Tier I - Results (Default if not specified) \_\_\_\_\_  
 Tier II (Results + QC Summaries) \_\_\_\_\_  
 Tier III (Results + QC & Calibration Summaries) **X**  
 Tier IV (Data Validation Package) 10% Surcharge \_\_\_\_\_

Relinquished by: (Signature) *Shannon Gale* Date: **04/03/12** Time: **0945**  
 Received by: (Signature) *[Signature]* Date: **4/16/12** Time: **0830**  
 Project Requirements (MRLS, GAPP)  
 Cooler / Blank Temperature \_\_\_\_\_ °C

**Sample Acceptance Check Form**

Client: Montana DEQ Work order: P1201320  
 Project: Background Indoor Air Study  
 Sample(s) received on: 4/4/12 Date opened: 4/4/12 by: MZAMORA

**Note:** This form is used for all samples received by CAS. The use of this form for custody seals is strictly meant to indicate presence/absence and not as an indication of compliance or nonconformity. Thermal preservation and pH will only be evaluated either at the request of the client and/or as required by the method/SOP.

- |  | <u>Yes</u>                          | <u>No</u>                           | <u>N/A</u>                          |
|--|-------------------------------------|-------------------------------------|-------------------------------------|
| 1  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 2  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 3  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 4  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 5  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 6  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 7  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 8  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 9  | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| 10   | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| Location of seal(s)? _____ Sealing Lid?            | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were signature and date included?                  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were seals intact?                                 | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were custody seals on outside of sample container? | <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| Location of seal(s)? _____ Sealing Lid?            | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were signature and date included?                  | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| Were seals intact?                                 | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 11   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 12   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 13   | <input type="checkbox"/>            | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |

Lab Sample ID	Container Description	Required pH *	Received pH	Adjusted pH	VOA Headspace (Presence/Absence)	Receipt / Preservation Comments
P1201320-001.01	6.0 L Ambient Can					
P1201320-002.01	6.0 L Ambient Can					
P1201320-003.01	6.0 L Ambient Can					
P1201320-004.01	6.0 L Ambient Can					
P1201320-005.01	6.0 L Silonite Can					
P1201320-006.01	6.0 L Ambient Can					
P1201320-007.01	6.0 L Ambient Can					
P1201320-008.01	6.0 L Ambient Can					

Explain any discrepancies: (include lab sample ID numbers): \_\_\_\_\_



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 31-1F-032412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-001

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01571

Date Collected: 3/25/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.72      Final Pressure (psig): 3.69

Canister Dilution Factor: 1.42

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	94	28	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	44	14	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	7.1	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 37-1F-032612  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-002

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01118

Date Collected: 3/27/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.87      Final Pressure (psig): 3.59

Canister Dilution Factor: 1.86

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	71	37	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	62	19	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	9.3	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 29-1F-032612  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-003

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01127

Date Collected: 3/27/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.68      Final Pressure (psig): 3.65

Canister Dilution Factor: 1.53

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	370	31	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	30	15	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	34	7.7	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** DUP-54  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-004

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01749

Date Collected: 3/27/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.09      Final Pressure (psig): 3.72

Canister Dilution Factor: 1.92

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	280	38	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	24	19	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	28	9.6	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 30-1F-032612  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-005

Test Code: Massachusetts APH, Revision 1, December 2009  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00147

Date Collected: 3/27/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.30 Final Pressure (psig): 3.68

Canister Dilution Factor: 1.96

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	47	39	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	23	20	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	9.8	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 33-1B-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-006

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00781

Date Collected: 3/28/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 0.50 Liter(s)

Initial Pressure (psig): -2.17      Final Pressure (psig): 3.68

Canister Dilution Factor: 1.47

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	79	59	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	29	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	15	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 32-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-007

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01142

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.83      Final Pressure (psig): 3.72

Canister Dilution Factor: 1.55

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	110	31	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	46	16	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	8.6	7.8	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 36-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-008

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01413

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.51      Final Pressure (psig): 3.73

Canister Dilution Factor: 1.81

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	60	36	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	18	18	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	9.1	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 34-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-009

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01533

Date Collected: 3/30/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.39      Final Pressure (psig): 3.57

Canister Dilution Factor: 1.37

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	46	27	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	14	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	6.9	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 35-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-010

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00943

Date Collected: 3/30/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.85      Final Pressure (psig): 3.66

Canister Dilution Factor: 1.86

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	81	37	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	29	19	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	12	9.3	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P120407-MB

Test Code: Massachusetts APH, Revision 1, December 2009  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/07/12  
Volume(s) Analyzed: 1.00 Liter(s)

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	20	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	10	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	5.0	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P120409-MB

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/09/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Data Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	ND	20	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	ND	10	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	ND	5.0	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P120407-LCS

Test Code: Massachusetts APH, Revision 1, December 2009  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/07/12  
 Volume(s) Analyzed: NA Liter(s)

Compound	Spike Amount ng	Result ng	% Recovery	CAS	Data Qualifier
				Acceptance Limits	
C5 - C8 Aliphatic Hydrocarbons	24.5	25.1	102	70-130	
C9 - C12 Aliphatic Hydrocarbons	25.0	25.2	101	70-130	
C9 - C10 Aromatic Hydrocarbons	48.5	46.3	95	70-130	

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P120409-LCS

Test Code: Massachusetts APH, Revision 1, December 2009  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/09/12  
Volume(s) Analyzed: NA Liter(s)

Compound	Spike Amount ng	Result ng	% Recovery	CAS	Data Qualifier
				Acceptance Limits	
C5 - C8 Aliphatic Hydrocarbons	24.5	24.6	100	70-130	
C9 - C12 Aliphatic Hydrocarbons	25.0	24.8	99	70-130	
C9 - C10 Aromatic Hydrocarbons	48.5	45.7	94	70-130	

LABORATORY DUPLICATE SUMMARY RESULTS

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**Client:** Montana DEQ

**Client Sample ID:** DUP-54

**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320

CAS Sample ID: P1201320-004DUP

Test Code: Massachusetts APH, Revision 1, December 2009

Date Collected: 3/27/12

Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8

Date Received: 4/4/12

Analyst: Elsa Moctezuma

Date Analyzed: 4/9/12

Sampling Media: 6.0 L Summa Canister

Volume(s) Analyzed: 1.00 Liter(s)

Test Notes:

Container ID: AC01749

Initial Pressure (psig): -5.09 Final Pressure (psig): 3.72

Canister Dilution Factor: 1.92

Compound	Sample Result	Duplicate Sample Result	Average	% RPD	RPD	Data
	$\mu\text{g}/\text{m}^3$	$\mu\text{g}/\text{m}^3$	$\mu\text{g}/\text{m}^3$		Limit	Qualifier
C <sub>5</sub> - C <sub>8</sub> Aliphatic Hydrocarbons <sup>1,2</sup>	282	280	281	<b>0.7</b>	30	
C <sub>9</sub> - C <sub>12</sub> Aliphatic Hydrocarbons <sup>1,3</sup>	23.7	21.4	22.55	<b>10</b>	30	
C <sub>9</sub> - C <sub>10</sub> Aromatic Hydrocarbons	28.3	28.5	28.4	<b>0.7</b>	30	

Significant non-petroleum related peaks (i.e. halogenated, oxygenated, terpenes, etc.) are subtracted from the hydrocarbon range areas when present.

<sup>1</sup>Hydrocarbon Range data from total ion chromatogram excluding any internal/tuning standards eluting in that range.

<sup>2</sup>C<sub>5</sub>-C<sub>8</sub> Aliphatic Hydrocarbons exclude the concentration of Target APH analytes eluting in that range.

<sup>3</sup>C<sub>9</sub>-C<sub>12</sub> Aliphatic Hydrocarbons exclude concentration of Target APH Analytes eluting in that range and concentration of C<sub>9</sub>-C<sub>10</sub> Aromatic Hydrocarbons.

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 31-1F-032412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-001

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01571

Date Collected: 3/25/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12 & 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)  
0.10 Liter(s)

Initial Pressure (psig): -1.72 Final Pressure (psig): 3.69

Canister Dilution Factor: 1.42

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	2.4	0.71	1.4	0.41	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.1	0.71	0.43	0.14	
74-87-3	Chloromethane	ND	0.71	ND	0.34	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.71	ND	0.10	
106-99-0	1,3-Butadiene	ND	0.71	ND	0.32	
74-83-9	Bromomethane	ND	0.71	ND	0.18	
75-00-3	Chloroethane	ND	0.71	ND	0.27	
64-17-5	Ethanol	1,300	71	700	38	D
75-05-8	Acetonitrile	0.79	0.71	0.47	0.42	
107-02-8	Acrolein	4.3	2.8	1.9	1.2	
67-64-1	Acetone	83	7.1	35	3.0	
75-69-4	Trichlorofluoromethane	1.1	0.71	0.20	0.13	
67-63-0	2-Propanol (Isopropyl Alcohol)	7.3	1.4	3.0	0.58	
107-13-1	Acrylonitrile	ND	0.71	ND	0.33	
75-35-4	1,1-Dichloroethene	ND	0.71	ND	0.18	
75-09-2	Methylene Chloride	ND	0.71	ND	0.20	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.71	ND	0.23	
76-13-1	Trichlorotrifluoroethane	ND	0.71	ND	0.093	
75-15-0	Carbon Disulfide	ND	7.1	ND	2.3	
156-60-5	trans-1,2-Dichloroethene	ND	0.71	ND	0.18	
75-34-3	1,1-Dichloroethane	ND	0.71	ND	0.18	
1634-04-4	Methyl tert-Butyl Ether	ND	0.71	ND	0.20	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 31-1F-032412  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201320  
**CAS Sample ID:** P1201320-001

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01571

**Date Collected:** 3/25/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/7/12 & 4/9/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -1.72      Final Pressure (psig): 3.69

Canister Dilution Factor: 1.42

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	7.1	ND	2.0	
156-59-2	cis-1,2-Dichloroethene	ND	0.71	ND	0.18	
78-93-3	2-Butanone (MEK)	ND	7.1	ND	2.4	
141-78-6	Ethyl Acetate	<b>28</b>	1.4	<b>7.9</b>	0.39	
110-54-3	n-Hexane	<b>2.2</b>	0.71	<b>0.63</b>	0.20	
67-66-3	Chloroform	<b>1.2</b>	0.71	<b>0.25</b>	0.15	
109-99-9	Tetrahydrofuran (THF)	<b>0.80</b>	0.71	<b>0.27</b>	0.24	
71-55-6	1,1,1-Trichloroethane	ND	0.71	ND	0.13	
56-23-5	Carbon Tetrachloride	<b>0.88</b>	0.71	<b>0.14</b>	0.11	
110-82-7	Cyclohexane	ND	1.4	ND	0.41	
78-87-5	1,2-Dichloropropane	ND	0.71	ND	0.15	
75-27-4	Bromodichloromethane	ND	0.71	ND	0.11	
123-91-1	1,4-Dioxane	ND	0.71	ND	0.20	
80-62-6	Methyl Methacrylate	ND	1.4	ND	0.35	
142-82-5	n-Heptane	<b>1.3</b>	0.71	<b>0.33</b>	0.17	
10061-01-5	cis-1,3-Dichloropropene	ND	0.71	ND	0.16	
108-10-1	4-Methyl-2-pentanone	<b>0.87</b>	0.71	<b>0.21</b>	0.17	
10061-02-6	trans-1,3-Dichloropropene	ND	0.71	ND	0.16	
108-88-3	Toluene	<b>6.1</b>	0.71	<b>1.6</b>	0.19	
591-78-6	2-Hexanone	<b>0.96</b>	0.71	<b>0.24</b>	0.17	
124-48-1	Dibromochloromethane	ND	0.71	ND	0.083	
106-93-4	1,2-Dibromoethane	ND	0.71	ND	0.092	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 31-1F-032412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-001

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01571

Date Collected: 3/25/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/7/12 & 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -1.72 Final Pressure (psig): 3.69

Canister Dilution Factor: 1.42

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	5.9	0.71	1.2	0.15	
111-65-9	n-Octane	ND	0.71	ND	0.15	
108-90-7	Chlorobenzene	ND	0.71	ND	0.15	
179601-23-1	m,p-Xylenes	2.3	1.4	0.52	0.33	
75-25-2	Bromoform	ND	0.71	ND	0.069	
100-42-5	Styrene	1.6	0.71	0.38	0.17	
95-47-6	o-Xylene	ND	0.71	ND	0.16	
111-84-2	n-Nonane	ND	0.71	ND	0.14	
98-82-8	Cumene	ND	0.71	ND	0.14	
80-56-8	alpha-Pinene	3.3	0.71	0.60	0.13	
103-65-1	n-Propylbenzene	ND	0.71	ND	0.14	
622-96-8	4-Ethyltoluene	ND	0.71	ND	0.14	
108-67-8	1,3,5-Trimethylbenzene	ND	0.71	ND	0.14	
95-63-6	1,2,4-Trimethylbenzene	ND	0.71	ND	0.14	
100-44-7	Benzyl Chloride	ND	0.71	ND	0.14	
541-73-1	1,3-Dichlorobenzene	ND	0.71	ND	0.12	
106-46-7	1,4-Dichlorobenzene	ND	0.71	ND	0.12	
95-50-1	1,2-Dichlorobenzene	ND	0.71	ND	0.12	
5989-27-5	d-Limonene	38	0.71	6.7	0.13	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.71	ND	0.073	
120-82-1	1,2,4-Trichlorobenzene	ND	0.71	ND	0.096	
87-68-3	Hexachlorobutadiene	ND	0.71	ND	0.067	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 37-1F-032612  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-002

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01118

Date Collected: 3/27/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.87 Final Pressure (psig): 3.59

Canister Dilution Factor: 1.86

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	1.3	0.93	0.77	0.54	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.1	0.93	0.42	0.19	
74-87-3	Chloromethane	ND	0.93	ND	0.45	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.93	ND	0.13	
106-99-0	1,3-Butadiene	ND	0.93	ND	0.42	
74-83-9	Bromomethane	ND	0.93	ND	0.24	
75-00-3	Chloroethane	ND	0.93	ND	0.35	
64-17-5	Ethanol	320	9.3	170	4.9	
75-05-8	Acetonitrile	ND	0.93	ND	0.55	
107-02-8	Acrolein	ND	3.7	ND	1.6	
67-64-1	Acetone	37	9.3	16	3.9	
75-69-4	Trichlorofluoromethane	1.1	0.93	0.20	0.17	
67-63-0	2-Propanol (Isopropyl Alcohol)	3.3	1.9	1.4	0.76	
107-13-1	Acrylonitrile	ND	0.93	ND	0.43	
75-35-4	1,1-Dichloroethene	ND	0.93	ND	0.23	
75-09-2	Methylene Chloride	ND	0.93	ND	0.27	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.93	ND	0.30	
76-13-1	Trichlorotrifluoroethane	ND	0.93	ND	0.12	
75-15-0	Carbon Disulfide	ND	9.3	ND	3.0	
156-60-5	trans-1,2-Dichloroethene	ND	0.93	ND	0.23	
75-34-3	1,1-Dichloroethane	ND	0.93	ND	0.23	
1634-04-4	Methyl tert-Butyl Ether	ND	0.93	ND	0.26	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 37-1F-032612  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-002

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01118

Date Collected: 3/27/12  
Date Received: 4/4/12  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.87 Final Pressure (psig): 3.59

Canister Dilution Factor: 1.86

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	9.3	ND	2.6	
156-59-2	cis-1,2-Dichloroethene	ND	0.93	ND	0.23	
78-93-3	2-Butanone (MEK)	ND	9.3	ND	3.2	
141-78-6	Ethyl Acetate	<b>3.9</b>	1.9	<b>1.1</b>	0.52	
110-54-3	n-Hexane	<b>1.3</b>	0.93	<b>0.36</b>	0.26	
67-66-3	Chloroform	ND	0.93	ND	0.19	
109-99-9	Tetrahydrofuran (THF)	ND	0.93	ND	0.32	
71-55-6	1,1,1-Trichloroethane	ND	0.93	ND	0.17	
56-23-5	Carbon Tetrachloride	ND	0.93	ND	0.15	
110-82-7	Cyclohexane	ND	1.9	ND	0.54	
78-87-5	1,2-Dichloropropane	ND	0.93	ND	0.20	
75-27-4	Bromodichloromethane	ND	0.93	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.93	ND	0.26	
80-62-6	Methyl Methacrylate	ND	1.9	ND	0.45	
142-82-5	n-Heptane	<b>2.6</b>	0.93	<b>0.63</b>	0.23	
10061-01-5	cis-1,3-Dichloropropene	ND	0.93	ND	0.20	
108-10-1	4-Methyl-2-pentanone	<b>1.7</b>	0.93	<b>0.42</b>	0.23	
10061-02-6	trans-1,3-Dichloropropene	ND	0.93	ND	0.20	
108-88-3	Toluene	<b>26</b>	0.93	<b>7.0</b>	0.25	
591-78-6	2-Hexanone	ND	0.93	ND	0.23	
124-48-1	Dibromochloromethane	ND	0.93	ND	0.11	
106-93-4	1,2-Dibromoethane	ND	0.93	ND	0.12	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 37-1F-032612  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201320  
**CAS Sample ID:** P1201320-002

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01118

**Date Collected:** 3/27/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/7/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -4.87      Final Pressure (psig): 3.59

Canister Dilution Factor: 1.86

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.93	ND	0.20	
111-65-9	n-Octane	<b>1.0</b>	0.93	<b>0.22</b>	0.20	
108-90-7	Chlorobenzene	ND	0.93	ND	0.20	
179601-23-1	m,p-Xylenes	<b>2.8</b>	1.9	<b>0.64</b>	0.43	
75-25-2	Bromoform	ND	0.93	ND	0.090	
100-42-5	Styrene	ND	0.93	ND	0.22	
95-47-6	o-Xylene	ND	0.93	ND	0.21	
111-84-2	n-Nonane	<b>4.3</b>	0.93	<b>0.81</b>	0.18	
98-82-8	Cumene	ND	0.93	ND	0.19	
80-56-8	alpha-Pinene	<b>4.9</b>	0.93	<b>0.88</b>	0.17	
103-65-1	n-Propylbenzene	ND	0.93	ND	0.19	
622-96-8	4-Ethyltoluene	ND	0.93	ND	0.19	
108-67-8	1,3,5-Trimethylbenzene	ND	0.93	ND	0.19	
95-63-6	1,2,4-Trimethylbenzene	ND	0.93	ND	0.19	
100-44-7	Benzyl Chloride	ND	0.93	ND	0.18	
541-73-1	1,3-Dichlorobenzene	ND	0.93	ND	0.15	
106-46-7	1,4-Dichlorobenzene	ND	0.93	ND	0.15	
95-50-1	1,2-Dichlorobenzene	ND	0.93	ND	0.15	
5989-27-5	d-Limonene	<b>33</b>	0.93	<b>6.0</b>	0.17	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.93	ND	0.096	
120-82-1	1,2,4-Trichlorobenzene	ND	0.93	ND	0.13	
87-68-3	Hexachlorobutadiene	ND	0.93	ND	0.087	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 29-1F-032612  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-003

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01127

Date Collected: 3/27/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.68 Final Pressure (psig): 3.65

Canister Dilution Factor: 1.53

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	5.0	0.77	2.9	0.44	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.1	0.77	0.42	0.15	
74-87-3	Chloromethane	0.77	0.77	0.37	0.37	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.77	ND	0.11	
106-99-0	1,3-Butadiene	ND	0.77	ND	0.35	
74-83-9	Bromomethane	ND	0.77	ND	0.20	
75-00-3	Chloroethane	ND	0.77	ND	0.29	
64-17-5	Ethanol	69	7.7	37	4.1	
75-05-8	Acetonitrile	ND	0.77	ND	0.46	
107-02-8	Acrolein	ND	3.1	ND	1.3	
67-64-1	Acetone	54	7.7	23	3.2	
75-69-4	Trichlorofluoromethane	1.1	0.77	0.20	0.14	
67-63-0	2-Propanol (Isopropyl Alcohol)	21	1.5	8.6	0.62	
107-13-1	Acrylonitrile	ND	0.77	ND	0.35	
75-35-4	1,1-Dichloroethene	ND	0.77	ND	0.19	
75-09-2	Methylene Chloride	ND	0.77	ND	0.22	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.77	ND	0.24	
76-13-1	Trichlorotrifluoroethane	ND	0.77	ND	0.10	
75-15-0	Carbon Disulfide	ND	7.7	ND	2.5	
156-60-5	trans-1,2-Dichloroethene	ND	0.77	ND	0.19	
75-34-3	1,1-Dichloroethane	ND	0.77	ND	0.19	
1634-04-4	Methyl tert-Butyl Ether	ND	0.77	ND	0.21	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 29-1F-032612  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-003

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01127

Date Collected: 3/27/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.68 Final Pressure (psig): 3.65

Canister Dilution Factor: 1.53

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	7.7	ND	2.2	
156-59-2	cis-1,2-Dichloroethene	ND	0.77	ND	0.19	
78-93-3	2-Butanone (MEK)	ND	7.7	ND	2.6	
141-78-6	Ethyl Acetate	1.6	1.5	0.45	0.42	
110-54-3	n-Hexane	16	0.77	4.5	0.22	
67-66-3	Chloroform	ND	0.77	ND	0.16	
109-99-9	Tetrahydrofuran (THF)	2.9	0.77	1.0	0.26	
71-55-6	1,1,1-Trichloroethane	ND	0.77	ND	0.14	
56-23-5	Carbon Tetrachloride	ND	0.77	ND	0.12	
110-82-7	Cyclohexane	6.0	1.5	1.7	0.44	
78-87-5	1,2-Dichloropropane	ND	0.77	ND	0.17	
75-27-4	Bromodichloromethane	ND	0.77	ND	0.11	
123-91-1	1,4-Dioxane	ND	0.77	ND	0.21	
80-62-6	Methyl Methacrylate	ND	1.5	ND	0.37	
142-82-5	n-Heptane	4.2	0.77	1.0	0.19	
10061-01-5	cis-1,3-Dichloropropene	ND	0.77	ND	0.17	
108-10-1	4-Methyl-2-pentanone	ND	0.77	ND	0.19	
10061-02-6	trans-1,3-Dichloropropene	ND	0.77	ND	0.17	
108-88-3	Toluene	56	0.77	15	0.20	
591-78-6	2-Hexanone	ND	0.77	ND	0.19	
124-48-1	Dibromochloromethane	ND	0.77	ND	0.090	
106-93-4	1,2-Dibromoethane	ND	0.77	ND	0.10	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 29-1F-032612  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-003

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01127

Date Collected: 3/27/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.68      Final Pressure (psig): 3.65

Canister Dilution Factor: 1.53

CAS #	Compound	Result μg/m <sup>3</sup>	MRL μg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.77	ND	0.16	
111-65-9	n-Octane	<b>1.5</b>	0.77	<b>0.31</b>	0.16	
108-90-7	Chlorobenzene	ND	0.77	ND	0.17	
179601-23-1	m,p-Xylenes	<b>26</b>	1.5	<b>6.0</b>	0.35	
75-25-2	Bromoform	ND	0.77	ND	0.074	
100-42-5	Styrene	<b>2.0</b>	0.77	<b>0.47</b>	0.18	
95-47-6	o-Xylene	<b>7.5</b>	0.77	<b>1.7</b>	0.18	
111-84-2	n-Nonane	<b>0.88</b>	0.77	<b>0.17</b>	0.15	
98-82-8	Cumene	ND	0.77	ND	0.16	
80-56-8	alpha-Pinene	<b>7.6</b>	0.77	<b>1.4</b>	0.14	
103-65-1	n-Propylbenzene	<b>1.6</b>	0.77	<b>0.32</b>	0.16	
622-96-8	4-Ethyltoluene	<b>2.8</b>	0.77	<b>0.57</b>	0.16	
108-67-8	1,3,5-Trimethylbenzene	<b>2.7</b>	0.77	<b>0.54</b>	0.16	
95-63-6	1,2,4-Trimethylbenzene	<b>8.5</b>	0.77	<b>1.7</b>	0.16	
100-44-7	Benzyl Chloride	ND	0.77	ND	0.15	
541-73-1	1,3-Dichlorobenzene	ND	0.77	ND	0.13	
106-46-7	1,4-Dichlorobenzene	ND	0.77	ND	0.13	
95-50-1	1,2-Dichlorobenzene	ND	0.77	ND	0.13	
5989-27-5	d-Limonene	<b>11</b>	0.77	<b>2.0</b>	0.14	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.77	ND	0.079	
120-82-1	1,2,4-Trichlorobenzene	ND	0.77	ND	0.10	
87-68-3	Hexachlorobutadiene	ND	0.77	ND	0.072	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** DUP-54  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-004

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01749

**Date Collected:** 3/27/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/9/12  
**Volume(s) Analyzed:** 1.00 Liter(s)

Initial Pressure (psig): -5.09      Final Pressure (psig): 3.72

Canister Dilution Factor: 1.92

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	12	0.96	6.8	0.56	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.0	0.96	0.41	0.19	
74-87-3	Chloromethane	ND	0.96	ND	0.47	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.96	ND	0.14	
106-99-0	1,3-Butadiene	ND	0.96	ND	0.43	
74-83-9	Bromomethane	ND	0.96	ND	0.25	
75-00-3	Chloroethane	ND	0.96	ND	0.36	
64-17-5	Ethanol	65	9.6	34	5.1	
75-05-8	Acetonitrile	ND	0.96	ND	0.57	
107-02-8	Acrolein	ND	3.8	ND	1.7	
67-64-1	Acetone	52	9.6	22	4.0	
75-69-4	Trichlorofluoromethane	1.1	0.96	0.20	0.17	
67-63-0	2-Propanol (Isopropyl Alcohol)	67	1.9	27	0.78	
107-13-1	Acrylonitrile	ND	0.96	ND	0.44	
75-35-4	1,1-Dichloroethene	ND	0.96	ND	0.24	
75-09-2	Methylene Chloride	ND	0.96	ND	0.28	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.96	ND	0.31	
76-13-1	Trichlorotrifluoroethane	ND	0.96	ND	0.13	
75-15-0	Carbon Disulfide	ND	9.6	ND	3.1	
156-60-5	trans-1,2-Dichloroethene	ND	0.96	ND	0.24	
75-34-3	1,1-Dichloroethane	ND	0.96	ND	0.24	
1634-04-4	Methyl tert-Butyl Ether	ND	0.96	ND	0.27	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** DUP-54  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-004

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01749

Date Collected: 3/27/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.09 Final Pressure (psig): 3.72

Canister Dilution Factor: 1.92

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	9.6	ND	2.7	
156-59-2	cis-1,2-Dichloroethene	ND	0.96	ND	0.24	
78-93-3	2-Butanone (MEK)	ND	9.6	ND	3.3	
141-78-6	Ethyl Acetate	ND	1.9	ND	0.53	
110-54-3	n-Hexane	<b>12</b>	0.96	<b>3.4</b>	0.27	
67-66-3	Chloroform	ND	0.96	ND	0.20	
109-99-9	Tetrahydrofuran (THF)	<b>2.2</b>	0.96	<b>0.74</b>	0.33	
71-55-6	1,1,1-Trichloroethane	ND	0.96	ND	0.18	
56-23-5	Carbon Tetrachloride	ND	0.96	ND	0.15	
110-82-7	Cyclohexane	<b>4.5</b>	1.9	<b>1.3</b>	0.56	
78-87-5	1,2-Dichloropropane	ND	0.96	ND	0.21	
75-27-4	Bromodichloromethane	ND	0.96	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.96	ND	0.27	
80-62-6	Methyl Methacrylate	ND	1.9	ND	0.47	
142-82-5	n-Heptane	<b>3.2</b>	0.96	<b>0.79</b>	0.23	
10061-01-5	cis-1,3-Dichloropropene	ND	0.96	ND	0.21	
108-10-1	4-Methyl-2-pentanone	ND	0.96	ND	0.23	
10061-02-6	trans-1,3-Dichloropropene	ND	0.96	ND	0.21	
108-88-3	Toluene	<b>45</b>	0.96	<b>12</b>	0.25	
591-78-6	2-Hexanone	ND	0.96	ND	0.23	
124-48-1	Dibromochloromethane	ND	0.96	ND	0.11	
106-93-4	1,2-Dibromoethane	ND	0.96	ND	0.12	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** DUP-54  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-004

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01749

Date Collected: 3/27/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.09 Final Pressure (psig): 3.72

Canister Dilution Factor: 1.92

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.96	ND	0.20	
111-65-9	n-Octane	1.2	0.96	0.26	0.21	
108-90-7	Chlorobenzene	ND	0.96	ND	0.21	
179601-23-1	m,p-Xylenes	21	1.9	4.9	0.44	
75-25-2	Bromoform	ND	0.96	ND	0.093	
100-42-5	Styrene	1.8	0.96	0.42	0.23	
95-47-6	o-Xylene	6.2	0.96	1.4	0.22	
111-84-2	n-Nonane	ND	0.96	ND	0.18	
98-82-8	Cumene	ND	0.96	ND	0.20	
80-56-8	alpha-Pinene	6.0	0.96	1.1	0.17	
103-65-1	n-Propylbenzene	1.4	0.96	0.28	0.20	
622-96-8	4-Ethyltoluene	2.4	0.96	0.48	0.20	
108-67-8	1,3,5-Trimethylbenzene	2.3	0.96	0.46	0.20	
95-63-6	1,2,4-Trimethylbenzene	7.2	0.96	1.5	0.20	
100-44-7	Benzyl Chloride	ND	0.96	ND	0.19	
541-73-1	1,3-Dichlorobenzene	ND	0.96	ND	0.16	
106-46-7	1,4-Dichlorobenzene	ND	0.96	ND	0.16	
95-50-1	1,2-Dichlorobenzene	ND	0.96	ND	0.16	
5989-27-5	d-Limonene	7.0	0.96	1.3	0.17	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.96	ND	0.099	
120-82-1	1,2,4-Trichlorobenzene	ND	0.96	ND	0.13	
87-68-3	Hexachlorobutadiene	ND	0.96	ND	0.090	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 30-1F-032612  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-005

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00147

Date Collected: 3/27/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.30 Final Pressure (psig): 3.68

Canister Dilution Factor: 1.96

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	1.4	0.98	0.82	0.57	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.5	0.98	0.50	0.20	
74-87-3	Chloromethane	ND	0.98	ND	0.47	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.98	ND	0.14	
106-99-0	1,3-Butadiene	ND	0.98	ND	0.44	
74-83-9	Bromomethane	ND	0.98	ND	0.25	
75-00-3	Chloroethane	ND	0.98	ND	0.37	
64-17-5	Ethanol	130	9.8	70	5.2	
75-05-8	Acetonitrile	ND	0.98	ND	0.58	
107-02-8	Acrolein	ND	3.9	ND	1.7	
67-64-1	Acetone	20	9.8	8.5	4.1	
75-69-4	Trichlorofluoromethane	1.1	0.98	0.20	0.17	
67-63-0	2-Propanol (Isopropyl Alcohol)	6.0	2.0	2.4	0.80	
107-13-1	Acrylonitrile	ND	0.98	ND	0.45	
75-35-4	1,1-Dichloroethene	ND	0.98	ND	0.25	
75-09-2	Methylene Chloride	ND	0.98	ND	0.28	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.98	ND	0.31	
76-13-1	Trichlorotrifluoroethane	ND	0.98	ND	0.13	
75-15-0	Carbon Disulfide	ND	9.8	ND	3.1	
156-60-5	trans-1,2-Dichloroethene	ND	0.98	ND	0.25	
75-34-3	1,1-Dichloroethane	ND	0.98	ND	0.24	
1634-04-4	Methyl tert-Butyl Ether	ND	0.98	ND	0.27	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 30-1F-032612  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-005

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AS00147

Date Collected: 3/27/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.30 Final Pressure (psig): 3.68

Canister Dilution Factor: 1.96

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	9.8	ND	2.8	
156-59-2	cis-1,2-Dichloroethene	ND	0.98	ND	0.25	
78-93-3	2-Butanone (MEK)	ND	9.8	ND	3.3	
141-78-6	Ethyl Acetate	ND	2.0	ND	0.54	
110-54-3	n-Hexane	<b>2.0</b>	0.98	<b>0.57</b>	0.28	
67-66-3	Chloroform	ND	0.98	ND	0.20	
109-99-9	Tetrahydrofuran (THF)	ND	0.98	ND	0.33	
71-55-6	1,1,1-Trichloroethane	ND	0.98	ND	0.18	
56-23-5	Carbon Tetrachloride	ND	0.98	ND	0.16	
110-82-7	Cyclohexane	ND	2.0	ND	0.57	
78-87-5	1,2-Dichloropropane	ND	0.98	ND	0.21	
75-27-4	Bromodichloromethane	ND	0.98	ND	0.15	
123-91-1	1,4-Dioxane	ND	0.98	ND	0.27	
80-62-6	Methyl Methacrylate	ND	2.0	ND	0.48	
142-82-5	n-Heptane	ND	0.98	ND	0.24	
10061-01-5	cis-1,3-Dichloropropene	ND	0.98	ND	0.22	
108-10-1	4-Methyl-2-pentanone	ND	0.98	ND	0.24	
10061-02-6	trans-1,3-Dichloropropene	ND	0.98	ND	0.22	
108-88-3	Toluene	<b>7.7</b>	0.98	<b>2.0</b>	0.26	
591-78-6	2-Hexanone	ND	0.98	ND	0.24	
124-48-1	Dibromochloromethane	ND	0.98	ND	0.12	
106-93-4	1,2-Dibromoethane	ND	0.98	ND	0.13	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 30-1F-032612  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-005

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00147

Date Collected: 3/27/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.30 Final Pressure (psig): 3.68

Canister Dilution Factor: 1.96

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.98	ND	0.21	
111-65-9	n-Octane	ND	0.98	ND	0.21	
108-90-7	Chlorobenzene	ND	0.98	ND	0.21	
179601-23-1	m,p-Xylenes	<b>3.9</b>	2.0	<b>0.90</b>	0.45	
75-25-2	Bromoform	ND	0.98	ND	0.095	
100-42-5	Styrene	ND	0.98	ND	0.23	
95-47-6	o-Xylene	<b>1.2</b>	0.98	<b>0.28</b>	0.23	
111-84-2	n-Nonane	ND	0.98	ND	0.19	
98-82-8	Cumene	ND	0.98	ND	0.20	
80-56-8	alpha-Pinene	<b>2.0</b>	0.98	<b>0.36</b>	0.18	
103-65-1	n-Propylbenzene	ND	0.98	ND	0.20	
622-96-8	4-Ethyltoluene	ND	0.98	ND	0.20	
108-67-8	1,3,5-Trimethylbenzene	ND	0.98	ND	0.20	
95-63-6	1,2,4-Trimethylbenzene	<b>1.2</b>	0.98	<b>0.23</b>	0.20	
100-44-7	Benzyl Chloride	ND	0.98	ND	0.19	
541-73-1	1,3-Dichlorobenzene	ND	0.98	ND	0.16	
106-46-7	1,4-Dichlorobenzene	ND	0.98	ND	0.16	
95-50-1	1,2-Dichlorobenzene	ND	0.98	ND	0.16	
5989-27-5	d-Limonene	<b>22</b>	0.98	<b>4.0</b>	0.18	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.98	ND	0.10	
120-82-1	1,2,4-Trichlorobenzene	ND	0.98	ND	0.13	
87-68-3	Hexachlorobutadiene	ND	0.98	ND	0.092	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 33-1B-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-006

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC00781

**Date Collected:** 3/28/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/9/12  
**Volume(s) Analyzed:** 0.50 Liter(s)  
 0.050 Liter(s)

Initial Pressure (psig): -2.17 Final Pressure (psig): 3.68

Canister Dilution Factor: 1.47

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
115-07-1	Propene	850	15	500	8.5	D
75-71-8	Dichlorodifluoromethane (CFC 12)	2.2	1.5	0.45	0.30	
74-87-3	Chloromethane	ND	1.5	ND	0.71	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	1.5	ND	0.21	
106-99-0	1,3-Butadiene	ND	1.5	ND	0.66	
74-83-9	Bromomethane	ND	1.5	ND	0.38	
75-00-3	Chloroethane	ND	1.5	ND	0.56	
64-17-5	Ethanol	2,000	150	1,000	78	D
75-05-8	Acetonitrile	1.7	1.5	1.0	0.88	
107-02-8	Acrolein	ND	5.9	ND	2.6	
67-64-1	Acetone	130	15	53	6.2	
75-69-4	Trichlorofluoromethane	ND	1.5	ND	0.26	
67-63-0	2-Propanol (Isopropyl Alcohol)	12	2.9	4.9	1.2	
107-13-1	Acrylonitrile	ND	1.5	ND	0.68	
75-35-4	1,1-Dichloroethene	ND	1.5	ND	0.37	
75-09-2	Methylene Chloride	2.2	1.5	0.63	0.42	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	1.5	ND	0.47	
76-13-1	Trichlorotrifluoroethane	ND	1.5	ND	0.19	
75-15-0	Carbon Disulfide	ND	15	ND	4.7	
156-60-5	trans-1,2-Dichloroethene	ND	1.5	ND	0.37	
75-34-3	1,1-Dichloroethane	ND	1.5	ND	0.36	
1634-04-4	Methyl tert-Butyl Ether	ND	1.5	ND	0.41	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 33-1B-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-006

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00781

Date Collected: 3/28/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 0.50 Liter(s)  
0.050 Liter(s)

Initial Pressure (psig): -2.17 Final Pressure (psig): 3.68

Canister Dilution Factor: 1.47

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	15	ND	4.2	
156-59-2	cis-1,2-Dichloroethene	ND	1.5	ND	0.37	
78-93-3	2-Butanone (MEK)	ND	15	ND	5.0	
141-78-6	Ethyl Acetate	4.0	2.9	1.1	0.82	
110-54-3	n-Hexane	ND	1.5	ND	0.42	
67-66-3	Chloroform	2.1	1.5	0.43	0.30	
109-99-9	Tetrahydrofuran (THF)	2.5	1.5	0.84	0.50	
71-55-6	1,1,1-Trichloroethane	ND	1.5	ND	0.27	
56-23-5	Carbon Tetrachloride	ND	1.5	ND	0.23	
110-82-7	Cyclohexane	3.7	2.9	1.1	0.85	
78-87-5	1,2-Dichloropropane	ND	1.5	ND	0.32	
75-27-4	Bromodichloromethane	ND	1.5	ND	0.22	
123-91-1	1,4-Dioxane	ND	1.5	ND	0.41	
80-62-6	Methyl Methacrylate	ND	2.9	ND	0.72	
142-82-5	n-Heptane	1.7	1.5	0.42	0.36	
10061-01-5	cis-1,3-Dichloropropene	ND	1.5	ND	0.32	
108-10-1	4-Methyl-2-pentanone	ND	1.5	ND	0.36	
10061-02-6	trans-1,3-Dichloropropene	ND	1.5	ND	0.32	
108-88-3	Toluene	8.7	1.5	2.3	0.39	
591-78-6	2-Hexanone	ND	1.5	ND	0.36	
124-48-1	Dibromochloromethane	ND	1.5	ND	0.17	
106-93-4	1,2-Dibromoethane	ND	1.5	ND	0.19	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 33-1B-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-006

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC00781

Date Collected: 3/28/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 0.50 Liter(s)  
0.050 Liter(s)

Initial Pressure (psig): -2.17 Final Pressure (psig): 3.68

Canister Dilution Factor: 1.47

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	4.2	1.5	0.88	0.31	
111-65-9	n-Octane	ND	1.5	ND	0.31	
108-90-7	Chlorobenzene	ND	1.5	ND	0.32	
179601-23-1	m,p-Xylenes	ND	2.9	ND	0.68	
75-25-2	Bromoform	ND	1.5	ND	0.14	
100-42-5	Styrene	ND	1.5	ND	0.35	
95-47-6	o-Xylene	ND	1.5	ND	0.34	
111-84-2	n-Nonane	ND	1.5	ND	0.28	
98-82-8	Cumene	ND	1.5	ND	0.30	
80-56-8	alpha-Pinene	52	1.5	9.3	0.26	
103-65-1	n-Propylbenzene	ND	1.5	ND	0.30	
622-96-8	4-Ethyltoluene	ND	1.5	ND	0.30	
108-67-8	1,3,5-Trimethylbenzene	ND	1.5	ND	0.30	
95-63-6	1,2,4-Trimethylbenzene	ND	1.5	ND	0.30	
100-44-7	Benzyl Chloride	ND	1.5	ND	0.28	
541-73-1	1,3-Dichlorobenzene	ND	1.5	ND	0.24	
106-46-7	1,4-Dichlorobenzene	ND	1.5	ND	0.24	
95-50-1	1,2-Dichlorobenzene	ND	1.5	ND	0.24	
5989-27-5	d-Limonene	46	1.5	8.3	0.26	
96-12-8	1,2-Dibromo-3-chloropropane	ND	1.5	ND	0.15	
120-82-1	1,2,4-Trichlorobenzene	ND	1.5	ND	0.20	
87-68-3	Hexachlorobutadiene	ND	1.5	ND	0.14	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 32-1F-032812  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201320  
**CAS Sample ID:** P1201320-007

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01142

**Date Collected:** 3/29/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/9/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -2.83      Final Pressure (psig): 3.72

Canister Dilution Factor: 1.55

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	14	0.78	8.4	0.45	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.1	0.78	0.42	0.16	
74-87-3	Chloromethane	1.0	0.78	0.51	0.38	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.78	ND	0.11	
106-99-0	1,3-Butadiene	ND	0.78	ND	0.35	
74-83-9	Bromomethane	ND	0.78	ND	0.20	
75-00-3	Chloroethane	ND	0.78	ND	0.29	
64-17-5	Ethanol	1,300	78	670	41	D
75-05-8	Acetonitrile	ND	0.78	ND	0.46	
107-02-8	Acrolein	ND	3.1	ND	1.4	
67-64-1	Acetone	130	7.8	56	3.3	
75-69-4	Trichlorofluoromethane	1.2	0.78	0.22	0.14	
67-63-0	2-Propanol (Isopropyl Alcohol)	57	1.6	23	0.63	
107-13-1	Acrylonitrile	ND	0.78	ND	0.36	
75-35-4	1,1-Dichloroethene	ND	0.78	ND	0.20	
75-09-2	Methylene Chloride	ND	0.78	ND	0.22	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.78	ND	0.25	
76-13-1	Trichlorotrifluoroethane	ND	0.78	ND	0.10	
75-15-0	Carbon Disulfide	ND	7.8	ND	2.5	
156-60-5	trans-1,2-Dichloroethene	ND	0.78	ND	0.20	
75-34-3	1,1-Dichloroethane	ND	0.78	ND	0.19	
1634-04-4	Methyl tert-Butyl Ether	ND	0.78	ND	0.22	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.

## RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 32-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-007

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC01142

**Date Collected:** 3/29/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/9/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -2.83      Final Pressure (psig): 3.72

Canister Dilution Factor: 1.55

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	7.8	ND	2.2	
156-59-2	cis-1,2-Dichloroethene	ND	0.78	ND	0.20	
78-93-3	2-Butanone (MEK)	<b>26</b>	7.8	<b>8.8</b>	2.6	
141-78-6	Ethyl Acetate	<b>11</b>	1.6	<b>3.0</b>	0.43	
110-54-3	n-Hexane	<b>11</b>	0.78	<b>3.2</b>	0.22	
67-66-3	Chloroform	<b>5.1</b>	0.78	<b>1.1</b>	0.16	
109-99-9	Tetrahydrofuran (THF)	ND	0.78	ND	0.26	
71-55-6	1,1,1-Trichloroethane	ND	0.78	ND	0.14	
56-23-5	Carbon Tetrachloride	<b>0.82</b>	0.78	<b>0.13</b>	0.12	
110-82-7	Cyclohexane	<b>2.4</b>	1.6	<b>0.71</b>	0.45	
78-87-5	1,2-Dichloropropane	ND	0.78	ND	0.17	
75-27-4	Bromodichloromethane	ND	0.78	ND	0.12	
123-91-1	1,4-Dioxane	ND	0.78	ND	0.22	
80-62-6	Methyl Methacrylate	ND	1.6	ND	0.38	
142-82-5	n-Heptane	<b>4.9</b>	0.78	<b>1.2</b>	0.19	
10061-01-5	cis-1,3-Dichloropropene	ND	0.78	ND	0.17	
108-10-1	4-Methyl-2-pentanone	<b>0.80</b>	0.78	<b>0.20</b>	0.19	
10061-02-6	trans-1,3-Dichloropropene	ND	0.78	ND	0.17	
108-88-3	Toluene	<b>49</b>	0.78	<b>13</b>	0.21	
591-78-6	2-Hexanone	ND	0.78	ND	0.19	
124-48-1	Dibromochloromethane	ND	0.78	ND	0.091	
106-93-4	1,2-Dibromoethane	ND	0.78	ND	0.10	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 32-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-007

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01142

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -2.83 Final Pressure (psig): 3.72

Canister Dilution Factor: 1.55

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	1.1	0.78	0.22	0.16	
111-65-9	n-Octane	1.1	0.78	0.24	0.17	
108-90-7	Chlorobenzene	ND	0.78	ND	0.17	
179601-23-1	m,p-Xylenes	4.4	1.6	1.0	0.36	
75-25-2	Bromoform	ND	0.78	ND	0.075	
100-42-5	Styrene	2.3	0.78	0.55	0.18	
95-47-6	o-Xylene	1.4	0.78	0.32	0.18	
111-84-2	n-Nonane	1.0	0.78	0.19	0.15	
98-82-8	Cumene	ND	0.78	ND	0.16	
80-56-8	alpha-Pinene	9.1	0.78	1.6	0.14	
103-65-1	n-Propylbenzene	ND	0.78	ND	0.16	
622-96-8	4-Ethyltoluene	ND	0.78	ND	0.16	
108-67-8	1,3,5-Trimethylbenzene	ND	0.78	ND	0.16	
95-63-6	1,2,4-Trimethylbenzene	1.4	0.78	0.29	0.16	
100-44-7	Benzyl Chloride	ND	0.78	ND	0.15	
541-73-1	1,3-Dichlorobenzene	ND	0.78	ND	0.13	
106-46-7	1,4-Dichlorobenzene	ND	0.78	ND	0.13	
95-50-1	1,2-Dichlorobenzene	ND	0.78	ND	0.13	
5989-27-5	d-Limonene	15	0.78	2.7	0.14	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.78	ND	0.080	
120-82-1	1,2,4-Trichlorobenzene	ND	0.78	ND	0.10	
87-68-3	Hexachlorobutadiene	ND	0.78	ND	0.073	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 36-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-008

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01413

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.51 Final Pressure (psig): 3.73

Canister Dilution Factor: 1.81

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	1.9	0.91	1.1	0.53	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.2	0.91	0.44	0.18	
74-87-3	Chloromethane	ND	0.91	ND	0.44	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.91	ND	0.13	
106-99-0	1,3-Butadiene	ND	0.91	ND	0.41	
74-83-9	Bromomethane	ND	0.91	ND	0.23	
75-00-3	Chloroethane	ND	0.91	ND	0.34	
64-17-5	Ethanol	450	9.1	240	4.8	
75-05-8	Acetonitrile	ND	0.91	ND	0.54	
107-02-8	Acrolein	8.5	3.6	3.7	1.6	
67-64-1	Acetone	44	9.1	18	3.8	
75-69-4	Trichlorofluoromethane	1.3	0.91	0.23	0.16	
67-63-0	2-Propanol (Isopropyl Alcohol)	6.4	1.8	2.6	0.74	
107-13-1	Acrylonitrile	ND	0.91	ND	0.42	
75-35-4	1,1-Dichloroethene	ND	0.91	ND	0.23	
75-09-2	Methylene Chloride	ND	0.91	ND	0.26	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.91	ND	0.29	
76-13-1	Trichlorotrifluoroethane	ND	0.91	ND	0.12	
75-15-0	Carbon Disulfide	ND	9.1	ND	2.9	
156-60-5	trans-1,2-Dichloroethene	ND	0.91	ND	0.23	
75-34-3	1,1-Dichloroethane	ND	0.91	ND	0.22	
1634-04-4	Methyl tert-Butyl Ether	ND	0.91	ND	0.25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 36-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-008

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01413

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.51 Final Pressure (psig): 3.73

Canister Dilution Factor: 1.81

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	9.1	ND	2.6	
156-59-2	cis-1,2-Dichloroethene	ND	0.91	ND	0.23	
78-93-3	2-Butanone (MEK)	<b>12</b>	9.1	<b>4.1</b>	3.1	
141-78-6	Ethyl Acetate	<b>2.7</b>	1.8	<b>0.75</b>	0.50	
110-54-3	n-Hexane	ND	0.91	ND	0.26	
67-66-3	Chloroform	ND	0.91	ND	0.19	
109-99-9	Tetrahydrofuran (THF)	<b>2.7</b>	0.91	<b>0.92</b>	0.31	
71-55-6	1,1,1-Trichloroethane	ND	0.91	ND	0.17	
56-23-5	Carbon Tetrachloride	ND	0.91	ND	0.14	
110-82-7	Cyclohexane	ND	1.8	ND	0.53	
78-87-5	1,2-Dichloropropane	ND	0.91	ND	0.20	
75-27-4	Bromodichloromethane	ND	0.91	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.91	ND	0.25	
80-62-6	Methyl Methacrylate	ND	1.8	ND	0.44	
142-82-5	n-Heptane	<b>1.3</b>	0.91	<b>0.31</b>	0.22	
10061-01-5	cis-1,3-Dichloropropene	ND	0.91	ND	0.20	
108-10-1	4-Methyl-2-pentanone	ND	0.91	ND	0.22	
10061-02-6	trans-1,3-Dichloropropene	ND	0.91	ND	0.20	
108-88-3	Toluene	<b>7.9</b>	0.91	<b>2.1</b>	0.24	
591-78-6	2-Hexanone	ND	0.91	ND	0.22	
124-48-1	Dibromochloromethane	ND	0.91	ND	0.11	
106-93-4	1,2-Dibromoethane	ND	0.91	ND	0.12	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 36-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-008

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01413

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.51 Final Pressure (psig): 3.73

Canister Dilution Factor: 1.81

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.91	ND	0.19	
111-65-9	n-Octane	<b>1.1</b>	0.91	<b>0.23</b>	0.19	
108-90-7	Chlorobenzene	ND	0.91	ND	0.20	
179601-23-1	m,p-Xylenes	ND	1.8	ND	0.42	
75-25-2	Bromoform	ND	0.91	ND	0.088	
100-42-5	Styrene	ND	0.91	ND	0.21	
95-47-6	o-Xylene	ND	0.91	ND	0.21	
111-84-2	n-Nonane	ND	0.91	ND	0.17	
98-82-8	Cumene	ND	0.91	ND	0.18	
80-56-8	alpha-Pinene	<b>3.5</b>	0.91	<b>0.63</b>	0.16	
103-65-1	n-Propylbenzene	ND	0.91	ND	0.18	
622-96-8	4-Ethyltoluene	ND	0.91	ND	0.18	
108-67-8	1,3,5-Trimethylbenzene	ND	0.91	ND	0.18	
95-63-6	1,2,4-Trimethylbenzene	ND	0.91	ND	0.18	
100-44-7	Benzyl Chloride	ND	0.91	ND	0.17	
541-73-1	1,3-Dichlorobenzene	ND	0.91	ND	0.15	
106-46-7	1,4-Dichlorobenzene	ND	0.91	ND	0.15	
95-50-1	1,2-Dichlorobenzene	ND	0.91	ND	0.15	
5989-27-5	d-Limonene	<b>24</b>	0.91	<b>4.3</b>	0.16	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.91	ND	0.094	
120-82-1	1,2,4-Trichlorobenzene	ND	0.91	ND	0.12	
87-68-3	Hexachlorobutadiene	ND	0.91	ND	0.085	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 34-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-009

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01533

Date Collected: 3/30/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.39 Final Pressure (psig): 3.57

Canister Dilution Factor: 1.37

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	2.1	0.69	1.2	0.40	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.0	0.69	0.41	0.14	
74-87-3	Chloromethane	ND	0.69	ND	0.33	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.69	ND	0.098	
106-99-0	1,3-Butadiene	ND	0.69	ND	0.31	
74-83-9	Bromomethane	ND	0.69	ND	0.18	
75-00-3	Chloroethane	ND	0.69	ND	0.26	
64-17-5	Ethanol	120	6.9	65	3.6	
75-05-8	Acetonitrile	ND	0.69	ND	0.41	
107-02-8	Acrolein	ND	2.7	ND	1.2	
67-64-1	Acetone	19	6.9	7.9	2.9	
75-69-4	Trichlorofluoromethane	1.1	0.69	0.19	0.12	
67-63-0	2-Propanol (Isopropyl Alcohol)	8.0	1.4	3.2	0.56	
107-13-1	Acrylonitrile	ND	0.69	ND	0.32	
75-35-4	1,1-Dichloroethene	ND	0.69	ND	0.17	
75-09-2	Methylene Chloride	ND	0.69	ND	0.20	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.69	ND	0.22	
76-13-1	Trichlorotrifluoroethane	ND	0.69	ND	0.089	
75-15-0	Carbon Disulfide	ND	6.9	ND	2.2	
156-60-5	trans-1,2-Dichloroethene	ND	0.69	ND	0.17	
75-34-3	1,1-Dichloroethane	ND	0.69	ND	0.17	
1634-04-4	Methyl tert-Butyl Ether	ND	0.69	ND	0.19	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 34-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-009

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01533

Date Collected: 3/30/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.39 Final Pressure (psig): 3.57

Canister Dilution Factor: 1.37

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	6.9	ND	1.9	
156-59-2	cis-1,2-Dichloroethene	ND	0.69	ND	0.17	
78-93-3	2-Butanone (MEK)	ND	6.9	ND	2.3	
141-78-6	Ethyl Acetate	3.5	1.4	0.98	0.38	
110-54-3	n-Hexane	1.5	0.69	0.43	0.19	
67-66-3	Chloroform	ND	0.69	ND	0.14	
109-99-9	Tetrahydrofuran (THF)	3.8	0.69	1.3	0.23	
71-55-6	1,1,1-Trichloroethane	ND	0.69	ND	0.13	
56-23-5	Carbon Tetrachloride	ND	0.69	ND	0.11	
110-82-7	Cyclohexane	ND	1.4	ND	0.40	
78-87-5	1,2-Dichloropropane	ND	0.69	ND	0.15	
75-27-4	Bromodichloromethane	ND	0.69	ND	0.10	
123-91-1	1,4-Dioxane	ND	0.69	ND	0.19	
80-62-6	Methyl Methacrylate	ND	1.4	ND	0.33	
142-82-5	n-Heptane	ND	0.69	ND	0.17	
10061-01-5	cis-1,3-Dichloropropene	ND	0.69	ND	0.15	
108-10-1	4-Methyl-2-pentanone	ND	0.69	ND	0.17	
10061-02-6	trans-1,3-Dichloropropene	ND	0.69	ND	0.15	
108-88-3	Toluene	4.5	0.69	1.2	0.18	
591-78-6	2-Hexanone	ND	0.69	ND	0.17	
124-48-1	Dibromochloromethane	ND	0.69	ND	0.080	
106-93-4	1,2-Dibromoethane	ND	0.69	ND	0.089	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 34-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-009

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01533

Date Collected: 3/30/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.39 Final Pressure (psig): 3.57

Canister Dilution Factor: 1.37

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.69	ND	0.14	
111-65-9	n-Octane	ND	0.69	ND	0.15	
108-90-7	Chlorobenzene	ND	0.69	ND	0.15	
179601-23-1	m,p-Xylenes	1.5	1.4	0.35	0.32	
75-25-2	Bromoform	ND	0.69	ND	0.066	
100-42-5	Styrene	0.85	0.69	0.20	0.16	
95-47-6	o-Xylene	ND	0.69	ND	0.16	
111-84-2	n-Nonane	ND	0.69	ND	0.13	
98-82-8	Cumene	ND	0.69	ND	0.14	
80-56-8	alpha-Pinene	1.3	0.69	0.24	0.12	
103-65-1	n-Propylbenzene	ND	0.69	ND	0.14	
622-96-8	4-Ethyltoluene	ND	0.69	ND	0.14	
108-67-8	1,3,5-Trimethylbenzene	ND	0.69	ND	0.14	
95-63-6	1,2,4-Trimethylbenzene	ND	0.69	ND	0.14	
100-44-7	Benzyl Chloride	ND	0.69	ND	0.13	
541-73-1	1,3-Dichlorobenzene	ND	0.69	ND	0.11	
106-46-7	1,4-Dichlorobenzene	ND	0.69	ND	0.11	
95-50-1	1,2-Dichlorobenzene	ND	0.69	ND	0.11	
5989-27-5	d-Limonene	9.6	0.69	1.7	0.12	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.69	ND	0.071	
120-82-1	1,2,4-Trichlorobenzene	ND	0.69	ND	0.092	
87-68-3	Hexachlorobutadiene	ND	0.69	ND	0.064	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** 35-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-010

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00943

Date Collected: 3/30/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9 - 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -4.85 Final Pressure (psig): 3.66

Canister Dilution Factor: 1.86

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	6.5	0.93	3.8	0.54	
75-71-8	Dichlorodifluoromethane (CFC 12)	2.2	0.93	0.45	0.19	
74-87-3	Chloromethane	ND	0.93	ND	0.45	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.93	ND	0.13	
106-99-0	1,3-Butadiene	ND	0.93	ND	0.42	
74-83-9	Bromomethane	ND	0.93	ND	0.24	
75-00-3	Chloroethane	ND	0.93	ND	0.35	
64-17-5	Ethanol	1,500	93	790	49	D
75-05-8	Acetonitrile	ND	0.93	ND	0.55	
107-02-8	Acrolein	ND	3.7	ND	1.6	
67-64-1	Acetone	49	9.3	21	3.9	
75-69-4	Trichlorofluoromethane	1.5	0.93	0.27	0.17	
67-63-0	2-Propanol (Isopropyl Alcohol)	46	1.9	19	0.76	
107-13-1	Acrylonitrile	ND	0.93	ND	0.43	
75-35-4	1,1-Dichloroethene	ND	0.93	ND	0.23	
75-09-2	Methylene Chloride	ND	0.93	ND	0.27	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.93	ND	0.30	
76-13-1	Trichlorotrifluoroethane	ND	0.93	ND	0.12	
75-15-0	Carbon Disulfide	ND	9.3	ND	3.0	
156-60-5	trans-1,2-Dichloroethene	ND	0.93	ND	0.23	
75-34-3	1,1-Dichloroethane	ND	0.93	ND	0.23	
1634-04-4	Methyl tert-Butyl Ether	ND	0.93	ND	0.26	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

D = The reported result is from a dilution.



## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 35-1F-032912  
**Client Project ID:** Background Indoor Air Study

**CAS Project ID:** P1201320  
**CAS Sample ID:** P1201320-010

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC00943

**Date Collected:** 3/30/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/9 - 4/10/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -4.85 Final Pressure (psig): 3.66

Canister Dilution Factor: 1.86

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	9.3	ND	2.6	
156-59-2	cis-1,2-Dichloroethene	ND	0.93	ND	0.23	
78-93-3	2-Butanone (MEK)	ND	9.3	ND	3.2	
141-78-6	Ethyl Acetate	16	1.9	4.5	0.52	
110-54-3	n-Hexane	3.2	0.93	0.90	0.26	
67-66-3	Chloroform	1.3	0.93	0.28	0.19	
109-99-9	Tetrahydrofuran (THF)	ND	0.93	ND	0.32	
71-55-6	1,1,1-Trichloroethane	ND	0.93	ND	0.17	
56-23-5	Carbon Tetrachloride	ND	0.93	ND	0.15	
110-82-7	Cyclohexane	ND	1.9	ND	0.54	
78-87-5	1,2-Dichloropropane	ND	0.93	ND	0.20	
75-27-4	Bromodichloromethane	ND	0.93	ND	0.14	
123-91-1	1,4-Dioxane	ND	0.93	ND	0.26	
80-62-6	Methyl Methacrylate	ND	1.9	ND	0.45	
142-82-5	n-Heptane	1.5	0.93	0.37	0.23	
10061-01-5	cis-1,3-Dichloropropene	ND	0.93	ND	0.20	
108-10-1	4-Methyl-2-pentanone	1.4	0.93	0.34	0.23	
10061-02-6	trans-1,3-Dichloropropene	ND	0.93	ND	0.20	
108-88-3	Toluene	16	0.93	4.2	0.25	
591-78-6	2-Hexanone	ND	0.93	ND	0.23	
124-48-1	Dibromochloromethane	ND	0.93	ND	0.11	
106-93-4	1,2-Dibromoethane	ND	0.93	ND	0.12	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 35-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-010

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**  
**Container ID:** AC00943

**Date Collected:** 3/30/12  
**Date Received:** 4/4/12  
**Date Analyzed:** 4/9 - 4/10/12  
**Volume(s) Analyzed:** 1.00 Liter(s)  
 0.10 Liter(s)

Initial Pressure (psig): -4.85      Final Pressure (psig): 3.66

Canister Dilution Factor: 1.86

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	1.3	0.93	0.27	0.20	
111-65-9	n-Octane	ND	0.93	ND	0.20	
108-90-7	Chlorobenzene	ND	0.93	ND	0.20	
179601-23-1	m,p-Xylenes	5.4	1.9	1.3	0.43	
75-25-2	Bromoform	ND	0.93	ND	0.090	
100-42-5	Styrene	ND	0.93	ND	0.22	
95-47-6	o-Xylene	1.9	0.93	0.43	0.21	
111-84-2	n-Nonane	ND	0.93	ND	0.18	
98-82-8	Cumene	ND	0.93	ND	0.19	
80-56-8	alpha-Pinene	5.1	0.93	0.92	0.17	
103-65-1	n-Propylbenzene	ND	0.93	ND	0.19	
622-96-8	4-Ethyltoluene	ND	0.93	ND	0.19	
108-67-8	1,3,5-Trimethylbenzene	ND	0.93	ND	0.19	
95-63-6	1,2,4-Trimethylbenzene	2.3	0.93	0.47	0.19	
100-44-7	Benzyl Chloride	ND	0.93	ND	0.18	
541-73-1	1,3-Dichlorobenzene	ND	0.93	ND	0.15	
106-46-7	1,4-Dichlorobenzene	ND	0.93	ND	0.15	
95-50-1	1,2-Dichlorobenzene	ND	0.93	ND	0.15	
5989-27-5	d-Limonene	28	0.93	5.0	0.17	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.93	ND	0.096	
120-82-1	1,2,4-Trichlorobenzene	ND	0.93	ND	0.13	
87-68-3	Hexachlorobutadiene	ND	0.93	ND	0.087	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P120407-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	ND	0.50	ND	0.29	
75-71-8	Dichlorodifluoromethane (CFC 12)	ND	0.50	ND	0.10	
74-87-3	Chloromethane	ND	0.50	ND	0.24	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.50	ND	0.072	
106-99-0	1,3-Butadiene	ND	0.50	ND	0.23	
74-83-9	Bromomethane	ND	0.50	ND	0.13	
75-00-3	Chloroethane	ND	0.50	ND	0.19	
64-17-5	Ethanol	ND	5.0	ND	2.7	
75-05-8	Acetonitrile	ND	0.50	ND	0.30	
107-02-8	Acrolein	ND	2.0	ND	0.87	
67-64-1	Acetone	ND	5.0	ND	2.1	
75-69-4	Trichlorofluoromethane	ND	0.50	ND	0.089	
67-63-0	2-Propanol (Isopropyl Alcohol)	ND	1.0	ND	0.41	
107-13-1	Acrylonitrile	ND	0.50	ND	0.23	
75-35-4	1,1-Dichloroethene	ND	0.50	ND	0.13	
75-09-2	Methylene Chloride	ND	0.50	ND	0.14	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.50	ND	0.16	
76-13-1	Trichlorotrifluoroethane	ND	0.50	ND	0.065	
75-15-0	Carbon Disulfide	ND	5.0	ND	1.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.50	ND	0.13	
75-34-3	1,1-Dichloroethane	ND	0.50	ND	0.12	
1634-04-4	Methyl tert-Butyl Ether	ND	0.50	ND	0.14	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P120407-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/7/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	5.0	ND	1.4	
156-59-2	cis-1,2-Dichloroethene	ND	0.50	ND	0.13	
78-93-3	2-Butanone (MEK)	ND	5.0	ND	1.7	
141-78-6	Ethyl Acetate	ND	1.0	ND	0.28	
110-54-3	n-Hexane	ND	0.50	ND	0.14	
67-66-3	Chloroform	ND	0.50	ND	0.10	
109-99-9	Tetrahydrofuran (THF)	ND	0.50	ND	0.17	
71-55-6	1,1,1-Trichloroethane	ND	0.50	ND	0.092	
56-23-5	Carbon Tetrachloride	ND	0.50	ND	0.080	
110-82-7	Cyclohexane	ND	1.0	ND	0.29	
78-87-5	1,2-Dichloropropane	ND	0.50	ND	0.11	
75-27-4	Bromodichloromethane	ND	0.50	ND	0.075	
123-91-1	1,4-Dioxane	ND	0.50	ND	0.14	
80-62-6	Methyl Methacrylate	ND	1.0	ND	0.24	
142-82-5	n-Heptane	ND	0.50	ND	0.12	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-10-1	4-Methyl-2-pentanone	ND	0.50	ND	0.12	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-88-3	Toluene	ND	0.50	ND	0.13	
591-78-6	2-Hexanone	ND	0.50	ND	0.12	
124-48-1	Dibromochloromethane	ND	0.50	ND	0.059	
106-93-4	1,2-Dibromoethane	ND	0.50	ND	0.065	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P120407-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/7/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.50	ND	0.11	
111-65-9	n-Octane	ND	0.50	ND	0.11	
108-90-7	Chlorobenzene	ND	0.50	ND	0.11	
179601-23-1	m,p-Xylenes	ND	1.0	ND	0.23	
75-25-2	Bromoform	ND	0.50	ND	0.048	
100-42-5	Styrene	ND	0.50	ND	0.12	
95-47-6	o-Xylene	ND	0.50	ND	0.12	
111-84-2	n-Nonane	ND	0.50	ND	0.095	
98-82-8	Cumene	ND	0.50	ND	0.10	
80-56-8	alpha-Pinene	ND	0.50	ND	0.090	
103-65-1	n-Propylbenzene	ND	0.50	ND	0.10	
622-96-8	4-Ethyltoluene	ND	0.50	ND	0.10	
108-67-8	1,3,5-Trimethylbenzene	ND	0.50	ND	0.10	
95-63-6	1,2,4-Trimethylbenzene	ND	0.50	ND	0.10	
100-44-7	Benzyl Chloride	ND	0.50	ND	0.097	
541-73-1	1,3-Dichlorobenzene	ND	0.50	ND	0.083	
106-46-7	1,4-Dichlorobenzene	ND	0.50	ND	0.083	
95-50-1	1,2-Dichlorobenzene	ND	0.50	ND	0.083	
5989-27-5	d-Limonene	ND	0.50	ND	0.090	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.50	ND	0.052	
120-82-1	1,2,4-Trichlorobenzene	ND	0.50	ND	0.067	
87-68-3	Hexachlorobutadiene	ND	0.50	ND	0.047	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P120409-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
115-07-1	Propene	ND	0.50	ND	0.29	
75-71-8	Dichlorodifluoromethane (CFC 12)	ND	0.50	ND	0.10	
74-87-3	Chloromethane	ND	0.50	ND	0.24	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.50	ND	0.072	
106-99-0	1,3-Butadiene	ND	0.50	ND	0.23	
74-83-9	Bromomethane	ND	0.50	ND	0.13	
75-00-3	Chloroethane	ND	0.50	ND	0.19	
64-17-5	Ethanol	ND	5.0	ND	2.7	
75-05-8	Acetonitrile	ND	0.50	ND	0.30	
107-02-8	Acrolein	ND	2.0	ND	0.87	
67-64-1	Acetone	ND	5.0	ND	2.1	
75-69-4	Trichlorofluoromethane	ND	0.50	ND	0.089	
67-63-0	2-Propanol (Isopropyl Alcohol)	ND	1.0	ND	0.41	
107-13-1	Acrylonitrile	ND	0.50	ND	0.23	
75-35-4	1,1-Dichloroethene	ND	0.50	ND	0.13	
75-09-2	Methylene Chloride	ND	0.50	ND	0.14	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.50	ND	0.16	
76-13-1	Trichlorotrifluoroethane	ND	0.50	ND	0.065	
75-15-0	Carbon Disulfide	ND	5.0	ND	1.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.50	ND	0.13	
75-34-3	1,1-Dichloroethane	ND	0.50	ND	0.12	
1634-04-4	Methyl tert-Butyl Ether	ND	0.50	ND	0.14	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P120409-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	5.0	ND	1.4	
156-59-2	cis-1,2-Dichloroethene	ND	0.50	ND	0.13	
78-93-3	2-Butanone (MEK)	ND	5.0	ND	1.7	
141-78-6	Ethyl Acetate	ND	1.0	ND	0.28	
110-54-3	n-Hexane	ND	0.50	ND	0.14	
67-66-3	Chloroform	ND	0.50	ND	0.10	
109-99-9	Tetrahydrofuran (THF)	ND	0.50	ND	0.17	
71-55-6	1,1,1-Trichloroethane	ND	0.50	ND	0.092	
56-23-5	Carbon Tetrachloride	ND	0.50	ND	0.080	
110-82-7	Cyclohexane	ND	1.0	ND	0.29	
78-87-5	1,2-Dichloropropane	ND	0.50	ND	0.11	
75-27-4	Bromodichloromethane	ND	0.50	ND	0.075	
123-91-1	1,4-Dioxane	ND	0.50	ND	0.14	
80-62-6	Methyl Methacrylate	ND	1.0	ND	0.24	
142-82-5	n-Heptane	ND	0.50	ND	0.12	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-10-1	4-Methyl-2-pentanone	ND	0.50	ND	0.12	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-88-3	Toluene	ND	0.50	ND	0.13	
591-78-6	2-Hexanone	ND	0.50	ND	0.12	
124-48-1	Dibromochloromethane	ND	0.50	ND	0.059	
106-93-4	1,2-Dibromoethane	ND	0.50	ND	0.065	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P120409-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.50	ND	0.11	
111-65-9	n-Octane	ND	0.50	ND	0.11	
108-90-7	Chlorobenzene	ND	0.50	ND	0.11	
179601-23-1	m,p-Xylenes	ND	1.0	ND	0.23	
75-25-2	Bromoform	ND	0.50	ND	0.048	
100-42-5	Styrene	ND	0.50	ND	0.12	
95-47-6	o-Xylene	ND	0.50	ND	0.12	
111-84-2	n-Nonane	ND	0.50	ND	0.095	
98-82-8	Cumene	ND	0.50	ND	0.10	
80-56-8	alpha-Pinene	ND	0.50	ND	0.090	
103-65-1	n-Propylbenzene	ND	0.50	ND	0.10	
622-96-8	4-Ethyltoluene	ND	0.50	ND	0.10	
108-67-8	1,3,5-Trimethylbenzene	ND	0.50	ND	0.10	
95-63-6	1,2,4-Trimethylbenzene	ND	0.50	ND	0.10	
100-44-7	Benzyl Chloride	ND	0.50	ND	0.097	
541-73-1	1,3-Dichlorobenzene	ND	0.50	ND	0.083	
106-46-7	1,4-Dichlorobenzene	ND	0.50	ND	0.083	
95-50-1	1,2-Dichlorobenzene	ND	0.50	ND	0.083	
5989-27-5	d-Limonene	ND	0.50	ND	0.090	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.50	ND	0.052	
120-82-1	1,2,4-Trichlorobenzene	ND	0.50	ND	0.067	
87-68-3	Hexachlorobutadiene	ND	0.50	ND	0.047	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P120410-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result		MRL		Data Qualifier
		$\mu\text{g}/\text{m}^3$	$\mu\text{g}/\text{m}^3$	ppbV	ppbV	
115-07-1	Propene	ND	0.50	ND	0.29	
75-71-8	Dichlorodifluoromethane (CFC 12)	ND	0.50	ND	0.10	
74-87-3	Chloromethane	ND	0.50	ND	0.24	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	0.50	ND	0.072	
106-99-0	1,3-Butadiene	ND	0.50	ND	0.23	
74-83-9	Bromomethane	ND	0.50	ND	0.13	
75-00-3	Chloroethane	ND	0.50	ND	0.19	
64-17-5	Ethanol	ND	5.0	ND	2.7	
75-05-8	Acetonitrile	ND	0.50	ND	0.30	
107-02-8	Acrolein	ND	2.0	ND	0.87	
67-64-1	Acetone	ND	5.0	ND	2.1	
75-69-4	Trichlorofluoromethane	ND	0.50	ND	0.089	
67-63-0	2-Propanol (Isopropyl Alcohol)	ND	1.0	ND	0.41	
107-13-1	Acrylonitrile	ND	0.50	ND	0.23	
75-35-4	1,1-Dichloroethene	ND	0.50	ND	0.13	
75-09-2	Methylene Chloride	ND	0.50	ND	0.14	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	ND	0.50	ND	0.16	
76-13-1	Trichlorotrifluoroethane	ND	0.50	ND	0.065	
75-15-0	Carbon Disulfide	ND	5.0	ND	1.6	
156-60-5	trans-1,2-Dichloroethene	ND	0.50	ND	0.13	
75-34-3	1,1-Dichloroethane	ND	0.50	ND	0.12	
1634-04-4	Methyl tert-Butyl Ether	ND	0.50	ND	0.14	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P120410-MB

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/10/12  
Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
108-05-4	Vinyl Acetate	ND	5.0	ND	1.4	
156-59-2	cis-1,2-Dichloroethene	ND	0.50	ND	0.13	
78-93-3	2-Butanone (MEK)	ND	5.0	ND	1.7	
141-78-6	Ethyl Acetate	ND	1.0	ND	0.28	
110-54-3	n-Hexane	ND	0.50	ND	0.14	
67-66-3	Chloroform	ND	0.50	ND	0.10	
109-99-9	Tetrahydrofuran (THF)	ND	0.50	ND	0.17	
71-55-6	1,1,1-Trichloroethane	ND	0.50	ND	0.092	
56-23-5	Carbon Tetrachloride	ND	0.50	ND	0.080	
110-82-7	Cyclohexane	ND	1.0	ND	0.29	
78-87-5	1,2-Dichloropropane	ND	0.50	ND	0.11	
75-27-4	Bromodichloromethane	ND	0.50	ND	0.075	
123-91-1	1,4-Dioxane	ND	0.50	ND	0.14	
80-62-6	Methyl Methacrylate	ND	1.0	ND	0.24	
142-82-5	n-Heptane	ND	0.50	ND	0.12	
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-10-1	4-Methyl-2-pentanone	ND	0.50	ND	0.12	
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	ND	0.11	
108-88-3	Toluene	ND	0.50	ND	0.13	
591-78-6	2-Hexanone	ND	0.50	ND	0.12	
124-48-1	Dibromochloromethane	ND	0.50	ND	0.059	
106-93-4	1,2-Dibromoethane	ND	0.50	ND	0.065	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P120410-MB

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
123-86-4	n-Butyl Acetate	ND	0.50	ND	0.11	
111-65-9	n-Octane	ND	0.50	ND	0.11	
108-90-7	Chlorobenzene	ND	0.50	ND	0.11	
179601-23-1	m,p-Xylenes	ND	1.0	ND	0.23	
75-25-2	Bromoform	ND	0.50	ND	0.048	
100-42-5	Styrene	ND	0.50	ND	0.12	
95-47-6	o-Xylene	ND	0.50	ND	0.12	
111-84-2	n-Nonane	ND	0.50	ND	0.095	
98-82-8	Cumene	ND	0.50	ND	0.10	
80-56-8	alpha-Pinene	ND	0.50	ND	0.090	
103-65-1	n-Propylbenzene	ND	0.50	ND	0.10	
622-96-8	4-Ethyltoluene	ND	0.50	ND	0.10	
108-67-8	1,3,5-Trimethylbenzene	ND	0.50	ND	0.10	
95-63-6	1,2,4-Trimethylbenzene	ND	0.50	ND	0.10	
100-44-7	Benzyl Chloride	ND	0.50	ND	0.097	
541-73-1	1,3-Dichlorobenzene	ND	0.50	ND	0.083	
106-46-7	1,4-Dichlorobenzene	ND	0.50	ND	0.083	
95-50-1	1,2-Dichlorobenzene	ND	0.50	ND	0.083	
5989-27-5	d-Limonene	ND	0.50	ND	0.090	
96-12-8	1,2-Dibromo-3-chloropropane	ND	0.50	ND	0.052	
120-82-1	1,2,4-Trichlorobenzene	ND	0.50	ND	0.067	
87-68-3	Hexachlorobutadiene	ND	0.50	ND	0.047	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister(s)  
 Test Notes:

Date(s) Collected: 3/25 - 3/30/12  
 Date(s) Received: 4/4/12  
 Date(s) Analyzed: 4/7 - 4/10/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		Percent Recovered	Percent Recovered	Percent Recovered		
Method Blank	P120407-MB	106	99	97	70-130	
Method Blank	P120409-MB	105	99	96	70-130	
Method Blank	P120410-MB	104	100	95	70-130	
Lab Control Sample	P120407-LCS	103	99	100	70-130	
Lab Control Sample	P120409-LCS	104	99	99	70-130	
Lab Control Sample	P120410-LCS	112	98	96	70-130	
31-1F-032412	P1201320-001	110	99	94	70-130	
37-1F-032612	P1201320-002	105	99	96	70-130	
29-1F-032612	P1201320-003	106	98	95	70-130	
DUP-54	P1201320-004	106	98	97	70-130	
DUP-54	P1201320-004DUP	106	98	97	70-130	
30-1F-032612	P1201320-005	106	100	92	70-130	
33-1B-032712	P1201320-006	107	101	92	70-130	
32-1F-032812	P1201320-007	104	99	95	70-130	
36-1F-032812	P1201320-008	110	99	90	70-130	
34-1F-032912	P1201320-009	108	97	91	70-130	
35-1F-032912	P1201320-010	110	97	94	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.

Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P120407-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/07/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
115-07-1	Propene	198	215	109	59-137	
75-71-8	Dichlorodifluoromethane (CFC 12)	196	170	87	63-115	
74-87-3	Chloromethane	190	173	91	59-124	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	200	170	85	65-113	
106-99-0	1,3-Butadiene	204	205	100	60-138	
74-83-9	Bromomethane	194	174	90	69-129	
75-00-3	Chloroethane	196	174	89	60-120	
64-17-5	Ethanol	928	861	93	58-121	
75-05-8	Acetonitrile	194	194	100	64-129	
107-02-8	Acrolein	198	183	92	54-127	
67-64-1	Acetone	1,010	922	91	59-114	
75-69-4	Trichlorofluoromethane	202	177	88	66-108	
67-63-0	2-Propanol (Isopropyl Alcohol)	382	299	78	50-113	
107-13-1	Acrylonitrile	198	199	101	72-135	
75-35-4	1,1-Dichloroethene	212	198	93	70-117	
75-09-2	Methylene Chloride	206	183	89	61-108	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	208	217	104	70-131	
76-13-1	Trichlorotrifluoroethane	206	191	93	70-113	
75-15-0	Carbon Disulfide	208	186	89	65-112	
156-60-5	trans-1,2-Dichloroethene	196	188	96	71-119	
75-34-3	1,1-Dichloroethane	200	192	96	71-116	
1634-04-4	Methyl tert-Butyl Ether	198	193	97	67-116	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result. Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P120407-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 4/07/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
108-05-4	Vinyl Acetate	952	992	104	59-142	
156-59-2	cis-1,2-Dichloroethene	206	195	95	69-119	
78-93-3	2-Butanone (MEK)	206	205	100	68-125	
141-78-6	Ethyl Acetate	398	391	98	63-130	
110-54-3	n-Hexane	198	180	91	57-120	
67-66-3	Chloroform	214	192	90	69-111	
109-99-9	Tetrahydrofuran (THF)	202	191	95	57-123	
71-55-6	1,1,1-Trichloroethane	198	186	94	73-119	
56-23-5	Carbon Tetrachloride	202	194	96	74-129	
110-82-7	Cyclohexane	390	355	91	70-113	
78-87-5	1,2-Dichloropropane	198	183	92	69-118	
75-27-4	Bromodichloromethane	198	188	95	75-124	
123-91-1	1,4-Dioxane	200	189	95	71-123	
80-62-6	Methyl Methacrylate	400	390	98	72-127	
142-82-5	n-Heptane	196	188	96	68-120	
10061-01-5	cis-1,3-Dichloropropene	188	183	97	71-130	
108-10-1	4-Methyl-2-pentanone	204	204	100	69-130	
10061-02-6	trans-1,3-Dichloropropene	210	212	101	76-133	
108-88-3	Toluene	202	179	89	67-111	
591-78-6	2-Hexanone	222	213	96	70-123	
124-48-1	Dibromochloromethane	206	197	96	75-129	
106-93-4	1,2-Dibromoethane	200	180	90	73-122	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P120407-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/07/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
123-86-4	n-Butyl Acetate	220	211	96	68-132	
111-65-9	n-Octane	200	195	98	68-116	
108-90-7	Chlorobenzene	202	177	88	69-113	
179601-23-1	m,p-Xylenes	392	360	92	70-116	
75-25-2	Bromoform	208	188	90	69-127	
100-42-5	Styrene	200	188	94	71-125	
95-47-6	o-Xylene	194	176	91	70-116	
111-84-2	n-Nonane	196	184	94	68-116	
98-82-8	Cumene	190	174	92	70-116	
80-56-8	alpha-Pinene	186	174	94	71-119	
103-65-1	n-Propylbenzene	192	172	90	71-119	
622-96-8	4-Ethyltoluene	198	174	88	71-119	
108-67-8	1,3,5-Trimethylbenzene	200	181	91	71-121	
95-63-6	1,2,4-Trimethylbenzene	194	180	93	73-127	
100-44-7	Benzyl Chloride	200	206	103	65-137	
541-73-1	1,3-Dichlorobenzene	200	164	82	68-123	
106-46-7	1,4-Dichlorobenzene	206	161	78	65-120	
95-50-1	1,2-Dichlorobenzene	198	163	82	67-121	
5989-27-5	d-Limonene	200	189	95	67-130	
96-12-8	1,2-Dibromo-3-chloropropane	196	181	92	72-133	
120-82-1	1,2,4-Trichlorobenzene	196	163	83	62-133	
87-68-3	Hexachlorobutadiene	202	166	82	60-128	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P120409-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/09/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
115-07-1	Propene	198	207	105	59-137	
75-71-8	Dichlorodifluoromethane (CFC 12)	196	168	86	63-115	
74-87-3	Chloromethane	190	171	90	59-124	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	200	170	85	65-113	
106-99-0	1,3-Butadiene	204	205	100	60-138	
74-83-9	Bromomethane	194	171	88	69-129	
75-00-3	Chloroethane	196	170	87	60-120	
64-17-5	Ethanol	928	895	96	58-121	
75-05-8	Acetonitrile	194	192	99	64-129	
107-02-8	Acrolein	198	180	91	54-127	
67-64-1	Acetone	1,010	909	90	59-114	
75-69-4	Trichlorofluoromethane	202	173	86	66-108	
67-63-0	2-Propanol (Isopropyl Alcohol)	382	313	82	50-113	
107-13-1	Acrylonitrile	198	198	100	72-135	
75-35-4	1,1-Dichloroethene	212	195	92	70-117	
75-09-2	Methylene Chloride	206	181	88	61-108	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	208	215	103	70-131	
76-13-1	Trichlorotrifluoroethane	206	185	90	70-113	
75-15-0	Carbon Disulfide	208	183	88	65-112	
156-60-5	trans-1,2-Dichloroethene	196	185	94	71-119	
75-34-3	1,1-Dichloroethane	200	191	96	71-116	
1634-04-4	Methyl tert-Butyl Ether	198	191	96	67-116	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result. Reported results are shown in concentration units and as a result of the calculation, may vary slightly.



## LABORATORY CONTROL SAMPLE SUMMARY

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P120409-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 4/09/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
108-05-4	Vinyl Acetate	952	981	103	59-142	
156-59-2	cis-1,2-Dichloroethene	206	193	94	69-119	
78-93-3	2-Butanone (MEK)	206	202	98	68-125	
141-78-6	Ethyl Acetate	398	389	98	63-130	
110-54-3	n-Hexane	198	180	91	57-120	
67-66-3	Chloroform	214	188	88	69-111	
109-99-9	Tetrahydrofuran (THF)	202	187	93	57-123	
71-55-6	1,1,1-Trichloroethane	198	181	91	73-119	
56-23-5	Carbon Tetrachloride	202	189	94	74-129	
110-82-7	Cyclohexane	390	349	89	70-113	
78-87-5	1,2-Dichloropropane	198	180	91	69-118	
75-27-4	Bromodichloromethane	198	185	93	75-124	
123-91-1	1,4-Dioxane	200	184	92	71-123	
80-62-6	Methyl Methacrylate	400	385	96	72-127	
142-82-5	n-Heptane	196	183	93	68-120	
10061-01-5	cis-1,3-Dichloropropene	188	179	95	71-130	
108-10-1	4-Methyl-2-pentanone	204	202	99	69-130	
10061-02-6	trans-1,3-Dichloropropene	210	207	99	76-133	
108-88-3	Toluene	202	176	87	67-111	
591-78-6	2-Hexanone	222	210	95	70-123	
124-48-1	Dibromochloromethane	206	195	95	75-129	
106-93-4	1,2-Dibromoethane	200	178	89	73-122	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 3 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P120409-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/09/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
123-86-4	n-Butyl Acetate	220	207	94	68-132	
111-65-9	n-Octane	200	194	97	68-116	
108-90-7	Chlorobenzene	202	173	86	69-113	
179601-23-1	m,p-Xylenes	392	352	90	70-116	
75-25-2	Bromoform	208	184	88	69-127	
100-42-5	Styrene	200	184	92	71-125	
95-47-6	o-Xylene	194	173	89	70-116	
111-84-2	n-Nonane	196	182	93	68-116	
98-82-8	Cumene	190	171	90	70-116	
80-56-8	alpha-Pinene	186	170	91	71-119	
103-65-1	n-Propylbenzene	192	169	88	71-119	
622-96-8	4-Ethyltoluene	198	173	87	71-119	
108-67-8	1,3,5-Trimethylbenzene	200	178	89	71-121	
95-63-6	1,2,4-Trimethylbenzene	194	177	91	73-127	
100-44-7	Benzyl Chloride	200	204	102	65-137	
541-73-1	1,3-Dichlorobenzene	200	161	81	68-123	
106-46-7	1,4-Dichlorobenzene	206	157	76	65-120	
95-50-1	1,2-Dichlorobenzene	198	160	81	67-121	
5989-27-5	d-Limonene	200	184	92	67-130	
96-12-8	1,2-Dibromo-3-chloropropane	196	179	91	72-133	
120-82-1	1,2,4-Trichlorobenzene	196	158	81	62-133	
87-68-3	Hexachlorobutadiene	202	161	80	60-128	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

## LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P120410-LCS

**Test Code:** EPA TO-15  
**Instrument ID:** Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
**Analyst:** Elsa Moctezuma  
**Sampling Media:** 6.0 L Summa Canister  
**Test Notes:**

**Date Collected:** NA  
**Date Received:** NA  
**Date Analyzed:** 4/10/12  
**Volume(s) Analyzed:** 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
115-07-1	Propene	198	234	118	59-137	
75-71-8	Dichlorodifluoromethane (CFC 12)	196	185	94	63-115	
74-87-3	Chloromethane	190	184	97	59-124	
76-14-2	1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	200	170	85	65-113	
106-99-0	1,3-Butadiene	204	214	105	60-138	
74-83-9	Bromomethane	194	174	90	69-129	
75-00-3	Chloroethane	196	179	91	60-120	
64-17-5	Ethanol	928	891	96	58-121	
75-05-8	Acetonitrile	194	206	106	64-129	
107-02-8	Acrolein	198	188	95	54-127	
67-64-1	Acetone	1,010	945	94	59-114	
75-69-4	Trichlorofluoromethane	202	181	90	66-108	
67-63-0	2-Propanol (Isopropyl Alcohol)	382	327	86	50-113	
107-13-1	Acrylonitrile	198	206	104	72-135	
75-35-4	1,1-Dichloroethene	212	202	95	70-117	
75-09-2	Methylene Chloride	206	184	89	61-108	
107-05-1	3-Chloro-1-propene (Allyl Chloride)	208	230	111	70-131	
76-13-1	Trichlorotrifluoroethane	206	186	90	70-113	
75-15-0	Carbon Disulfide	208	192	92	65-112	
156-60-5	trans-1,2-Dichloroethene	196	196	100	71-119	
75-34-3	1,1-Dichloroethane	200	202	101	71-116	
1634-04-4	Methyl tert-Butyl Ether	198	200	101	67-116	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
 Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

Page 2 of 3

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P120410-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/10/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
108-05-4	Vinyl Acetate	952	1010	106	59-142	
156-59-2	cis-1,2-Dichloroethene	206	203	99	69-119	
78-93-3	2-Butanone (MEK)	206	209	101	68-125	
141-78-6	Ethyl Acetate	398	404	102	63-130	
110-54-3	n-Hexane	198	191	96	57-120	
67-66-3	Chloroform	214	197	92	69-111	
109-99-9	Tetrahydrofuran (THF)	202	195	97	57-123	
71-55-6	1,1,1-Trichloroethane	198	186	94	73-119	
56-23-5	Carbon Tetrachloride	202	191	95	74-129	
110-82-7	Cyclohexane	390	350	90	70-113	
78-87-5	1,2-Dichloropropane	198	185	93	69-118	
75-27-4	Bromodichloromethane	198	191	96	75-124	
123-91-1	1,4-Dioxane	200	185	93	71-123	
80-62-6	Methyl Methacrylate	400	376	94	72-127	
142-82-5	n-Heptane	196	187	95	68-120	
10061-01-5	cis-1,3-Dichloropropene	188	185	98	71-130	
108-10-1	4-Methyl-2-pentanone	204	207	101	69-130	
10061-02-6	trans-1,3-Dichloropropene	210	213	101	76-133	
108-88-3	Toluene	202	174	86	67-111	
591-78-6	2-Hexanone	222	218	98	70-123	
124-48-1	Dibromochloromethane	206	192	93	75-129	
106-93-4	1,2-Dibromoethane	200	176	88	73-122	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY CONTROL SAMPLE SUMMARY

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**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P120410-LCS

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/10/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
123-86-4	n-Butyl Acetate	220	215	98	68-132	
111-65-9	n-Octane	200	196	98	68-116	
108-90-7	Chlorobenzene	202	172	85	69-113	
179601-23-1	m,p-Xylenes	392	352	90	70-116	
75-25-2	Bromoform	208	179	86	69-127	
100-42-5	Styrene	200	182	91	71-125	
95-47-6	o-Xylene	194	171	88	70-116	
111-84-2	n-Nonane	196	188	96	68-116	
98-82-8	Cumene	190	169	89	70-116	
80-56-8	alpha-Pinene	186	170	91	71-119	
103-65-1	n-Propylbenzene	192	170	89	71-119	
622-96-8	4-Ethyltoluene	198	171	86	71-119	
108-67-8	1,3,5-Trimethylbenzene	200	176	88	71-121	
95-63-6	1,2,4-Trimethylbenzene	194	177	91	73-127	
100-44-7	Benzyl Chloride	200	203	102	65-137	
541-73-1	1,3-Dichlorobenzene	200	158	79	68-123	
106-46-7	1,4-Dichlorobenzene	206	156	76	65-120	
95-50-1	1,2-Dichlorobenzene	198	159	80	67-121	
5989-27-5	d-Limonene	200	187	94	67-130	
96-12-8	1,2-Dibromo-3-chloropropane	196	177	90	72-133	
120-82-1	1,2,4-Trichlorobenzene	196	157	80	62-133	
87-68-3	Hexachlorobutadiene	202	162	80	60-128	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

LABORATORY DUPLICATE SUMMARY RESULTS

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**Client:** Montana DEQ  
**Client Sample ID:** DUP-54  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-004DUP

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01749

Date Collected: 3/27/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.09

Final Pressure (psig): 3.72

Canister Dilution Factor: 1.92

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
Propene	11.7	6.78	10.5	6.13	11.1	<b>11</b>	25	
Dichlorodifluoromethane (CFC 12)	2.00	0.405	2.02	0.409	2.01	<b>1</b>	25	
Chloromethane	ND	ND	ND	ND	-	-	25	
1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)	ND	ND	ND	ND	-	-	25	
1,3-Butadiene	ND	ND	ND	ND	-	-	25	
Bromomethane	ND	ND	ND	ND	-	-	25	
Chloroethane	ND	ND	ND	ND	-	-	25	
Ethanol	64.6	34.3	63.7	33.8	64.15	<b>1</b>	25	
Acetonitrile	ND	ND	ND	ND	-	-	25	
Acrolein	ND	ND	ND	ND	-	-	25	
Acetone	52.4	22.1	52.3	22.0	52.35	<b>0.2</b>	25	
Trichlorofluoromethane	1.11	0.197	1.08	0.192	1.095	<b>3</b>	25	
2-Propanol (Isopropyl Alcohol)	66.7	27.2	68.4	27.8	67.55	<b>3</b>	25	
Acrylonitrile	ND	ND	ND	ND	-	-	25	
1,1-Dichloroethene	ND	ND	ND	ND	-	-	25	
Methylene Chloride	ND	ND	ND	ND	-	-	25	
3-Chloro-1-propene (Allyl Chloride)	ND	ND	ND	ND	-	-	25	
Trichlorotrifluoroethane	ND	ND	ND	ND	-	-	25	
Carbon Disulfide	ND	ND	ND	ND	-	-	25	
trans-1,2-Dichloroethene	ND	ND	ND	ND	-	-	25	
1,1-Dichloroethane	ND	ND	ND	ND	-	-	25	
Methyl tert-Butyl Ether	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

LABORATORY DUPLICATE SUMMARY RESULTS

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**Client:** Montana DEQ  
**Client Sample ID:** DUP-54  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-004DUP

Test Code: EPA TO-15  
Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
Analyst: Elsa Moctezuma  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01749

Date Collected: 3/27/12  
Date Received: 4/4/12  
Date Analyzed: 4/9/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.09

Final Pressure (psig): 3.72

Canister Dilution Factor: 1.92

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
Vinyl Acetate	ND	ND	ND	ND	-	-	25	
cis-1,2-Dichloroethene	ND	ND	ND	ND	-	-	25	
2-Butanone (MEK)	ND	ND	ND	ND	-	-	25	
Ethyl Acetate	ND	ND	ND	ND	-	-	25	
n-Hexane	12.1	3.43	12.2	3.45	12.15	<b>0.8</b>	25	
Chloroform	ND	ND	ND	ND	-	-	25	
Tetrahydrofuran (THF)	2.19	0.743	2.34	0.795	2.265	<b>7</b>	25	
1,1,1-Trichloroethane	ND	ND	ND	ND	-	-	25	
Carbon Tetrachloride	ND	ND	ND	ND	-	-	25	
Cyclohexane	4.54	1.32	4.60	1.34	4.57	<b>1</b>	25	
1,2-Dichloropropane	ND	ND	ND	ND	-	-	25	
Bromodichloromethane	ND	ND	ND	ND	-	-	25	
1,4-Dioxane	ND	ND	ND	ND	-	-	25	
Methyl Methacrylate	ND	ND	ND	ND	-	-	25	
n-Heptane	3.23	0.789	3.25	0.793	3.24	<b>0.6</b>	25	
cis-1,3-Dichloropropene	ND	ND	ND	ND	-	-	25	
4-Methyl-2-pentanone	ND	ND	ND	ND	-	-	25	
trans-1,3-Dichloropropene	ND	ND	ND	ND	-	-	25	
Toluene	44.9	11.9	45.5	12.1	45.2	<b>1</b>	25	
2-Hexanone	ND	ND	ND	ND	-	-	25	
Dibromochloromethane	ND	ND	ND	ND	-	-	25	
1,2-Dibromoethane	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

LABORATORY DUPLICATE SUMMARY RESULTS

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**Client:** Montana DEQ  
**Client Sample ID:** DUP-54  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-004DUP

Test Code: EPA TO-15  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973inert/6890N/MS8  
 Analyst: Elsa Moctezuma  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01749

Date Collected: 3/27/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/9/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.09

Final Pressure (psig): 3.72

Canister Dilution Factor: 1.92

Compound	Sample Result		Duplicate Sample Result		Average µg/m <sup>3</sup>	% RPD	RPD Limit	Data Qualifier
	µg/m <sup>3</sup>	ppbV	µg/m <sup>3</sup>	ppbV				
n-Butyl Acetate	ND	ND	ND	ND	-	-	25	
n-Octane	1.20	0.257	1.17	0.251	1.185	<b>3</b>	25	
Chlorobenzene	ND	ND	ND	ND	-	-	25	
m,p-Xylenes	21.4	4.93	21.6	4.98	21.5	<b>0.9</b>	25	
Bromoform	ND	ND	ND	ND	-	-	25	
Styrene	1.79	0.420	1.78	0.419	1.785	<b>0.6</b>	25	
o-Xylene	6.18	1.42	6.26	1.44	6.22	<b>1</b>	25	
n-Nonane	ND	ND	ND	ND	-	-	25	
Cumene	ND	ND	ND	ND	-	-	25	
alpha-Pinene	6.03	1.08	6.06	1.09	6.045	<b>0.5</b>	25	
n-Propylbenzene	1.39	0.283	1.41	0.286	1.4	<b>1</b>	25	
4-Ethyltoluene	2.35	0.479	2.26	0.459	2.305	<b>4</b>	25	
1,3,5-Trimethylbenzene	2.26	0.460	2.25	0.459	2.255	<b>0.4</b>	25	
1,2,4-Trimethylbenzene	7.22	1.47	7.25	1.48	7.235	<b>0.4</b>	25	
Benzyl Chloride	ND	ND	ND	ND	-	-	25	
1,3-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
1,4-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
1,2-Dichlorobenzene	ND	ND	ND	ND	-	-	25	
d-Limonene	7.02	1.26	6.83	1.23	6.925	<b>3</b>	25	
1,2-Dibromo-3-chloropropane	ND	ND	ND	ND	-	-	25	
1,2,4-Trichlorobenzene	ND	ND	ND	ND	-	-	25	
Hexachlorobutadiene	ND	ND	ND	ND	-	-	25	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.



RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 31-1F-032412  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-001

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01571

Date Collected: 3/25/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.72      Final Pressure (psig): 3.69

Canister Dilution Factor: 1.42

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.036	ND	0.014	
107-06-2	1,2-Dichloroethane	<b>0.15</b>	0.036	<b>0.037</b>	0.0088	
71-43-2	Benzene	<b>1.8</b>	0.11	<b>0.57</b>	0.033	
79-01-6	Trichloroethene	<b>0.097</b>	0.036	<b>0.018</b>	0.0066	
79-00-5	1,1,2-Trichloroethane	ND	0.14	ND	0.026	
127-18-4	Tetrachloroethene	<b>0.051</b>	0.036	<b>0.0075</b>	0.0052	
100-41-4	Ethylbenzene	<b>0.86</b>	0.14	<b>0.20</b>	0.033	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.036	ND	0.0052	
91-20-3	Naphthalene	<b>0.51</b>	0.14	<b>0.097</b>	0.027	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 37-1F-032612  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-002

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01118

Date Collected: 3/27/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.87      Final Pressure (psig): 3.59

Canister Dilution Factor: 1.86

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.047	ND	0.018	
107-06-2	1,2-Dichloroethane	<b>0.074</b>	0.047	<b>0.018</b>	0.011	
71-43-2	Benzene	<b>0.87</b>	0.14	<b>0.27</b>	0.044	
79-01-6	Trichloroethene	<b>1.3</b>	0.047	<b>0.23</b>	0.0087	
79-00-5	1,1,2-Trichloroethane	ND	0.19	ND	0.034	
127-18-4	Tetrachloroethene	ND	0.047	ND	0.0069	
100-41-4	Ethylbenzene	<b>0.72</b>	0.19	<b>0.16</b>	0.043	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.047	ND	0.0068	
91-20-3	Naphthalene	<b>0.58</b>	0.19	<b>0.11</b>	0.035	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 29-1F-032612  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-003

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01127

Date Collected: 3/27/12  
Date Received: 4/4/12  
Date Analyzed: 4/10/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.68      Final Pressure (psig): 3.65

Canister Dilution Factor: 1.53

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.038	ND	0.015	
107-06-2	1,2-Dichloroethane	<b>0.15</b>	0.038	<b>0.037</b>	0.0095	
71-43-2	Benzene	<b>13</b>	0.11	<b>4.1</b>	0.036	
79-01-6	Trichloroethene	ND	0.038	ND	0.0071	
79-00-5	1,1,2-Trichloroethane	ND	0.15	ND	0.028	
127-18-4	Tetrachloroethene	ND	0.038	ND	0.0056	
100-41-4	Ethylbenzene	<b>6.1</b>	0.15	<b>1.4</b>	0.035	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.038	ND	0.0056	
91-20-3	Naphthalene	<b>0.92</b>	0.15	<b>0.18</b>	0.029	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** DUP-54  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-004

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01749

Date Collected: 3/27/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.09      Final Pressure (psig): 3.72

Canister Dilution Factor: 1.92

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.048	ND	0.019	
107-06-2	1,2-Dichloroethane	<b>0.17</b>	0.048	<b>0.042</b>	0.012	
71-43-2	Benzene	<b>10</b>	0.14	<b>3.1</b>	0.045	
79-01-6	Trichloroethene	ND	0.048	ND	0.0089	
79-00-5	1,1,2-Trichloroethane	ND	0.19	ND	0.035	
127-18-4	Tetrachloroethene	ND	0.048	ND	0.0071	
100-41-4	Ethylbenzene	<b>5.1</b>	0.19	<b>1.2</b>	0.044	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.048	ND	0.0070	
91-20-3	Naphthalene	<b>0.63</b>	0.19	<b>0.12</b>	0.037	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 30-1F-032612  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-005

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AS00147

Date Collected: 3/27/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -5.30      Final Pressure (psig): 3.68

Canister Dilution Factor: 1.96

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.049	ND	0.019	
107-06-2	1,2-Dichloroethane	<b>0.095</b>	0.049	<b>0.024</b>	0.012	
71-43-2	Benzene	<b>1.7</b>	0.15	<b>0.54</b>	0.046	
79-01-6	Trichloroethene	ND	0.049	ND	0.0091	
79-00-5	1,1,2-Trichloroethane	ND	0.20	ND	0.036	
127-18-4	Tetrachloroethene	ND	0.049	ND	0.0072	
100-41-4	Ethylbenzene	<b>0.93</b>	0.20	<b>0.21</b>	0.045	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.049	ND	0.0071	
91-20-3	Naphthalene	<b>0.27</b>	0.20	<b>0.052</b>	0.037	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

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**Client:** Montana DEQ  
**Client Sample ID:** 33-1B-032712  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-006

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00781

Date Collected: 3/28/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 0.50 Liter(s)

Initial Pressure (psig): -2.17      Final Pressure (psig): 3.68

Canister Dilution Factor: 1.47

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.074	ND	0.029	
107-06-2	1,2-Dichloroethane	<b>0.80</b>	0.074	<b>0.20</b>	0.018	
71-43-2	Benzene	<b>0.70</b>	0.22	<b>0.22</b>	0.069	
79-01-6	Trichloroethene	ND	0.074	ND	0.014	
79-00-5	1,1,2-Trichloroethane	ND	0.29	ND	0.054	
127-18-4	Tetrachloroethene	<b>0.31</b>	0.074	<b>0.045</b>	0.011	
100-41-4	Ethylbenzene	<b>1.3</b>	0.29	<b>0.29</b>	0.068	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.074	ND	0.011	
91-20-3	Naphthalene	<b>0.36</b>	0.29	<b>0.068</b>	0.056	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 32-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P1201320-007

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:  
Container ID: AC01142

Date Collected: 3/29/12  
Date Received: 4/4/12  
Date Analyzed: 4/10/12  
Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -2.83      Final Pressure (psig): 3.72

Canister Dilution Factor: 1.55

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.039	ND	0.015	
107-06-2	1,2-Dichloroethane	<b>0.40</b>	0.039	<b>0.10</b>	0.0096	
71-43-2	Benzene	<b>2.7</b>	0.12	<b>0.84</b>	0.036	
79-01-6	Trichloroethene	<b>0.056</b>	0.039	<b>0.010</b>	0.0072	
79-00-5	1,1,2-Trichloroethane	ND	0.16	ND	0.028	
127-18-4	Tetrachloroethene	<b>0.045</b>	0.039	<b>0.0067</b>	0.0057	
100-41-4	Ethylbenzene	<b>1.4</b>	0.16	<b>0.33</b>	0.036	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.039	ND	0.0056	
91-20-3	Naphthalene	<b>0.81</b>	0.16	<b>0.16</b>	0.030	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 36-1F-032812  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-008

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01413

Date Collected: 3/29/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.51      Final Pressure (psig): 3.73

Canister Dilution Factor: 1.81

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.045	ND	0.018	
107-06-2	1,2-Dichloroethane	<b>0.58</b>	0.045	<b>0.14</b>	0.011	
71-43-2	Benzene	<b>0.85</b>	0.14	<b>0.27</b>	0.043	
79-01-6	Trichloroethene	ND	0.045	ND	0.0084	
79-00-5	1,1,2-Trichloroethane	ND	0.18	ND	0.033	
127-18-4	Tetrachloroethene	ND	0.045	ND	0.0067	
100-41-4	Ethylbenzene	<b>0.40</b>	0.18	<b>0.093</b>	0.042	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.045	ND	0.0066	
91-20-3	Naphthalene	<b>0.21</b>	0.18	<b>0.040</b>	0.035	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.



RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 34-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-009

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC01533

Date Collected: 3/30/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -1.39      Final Pressure (psig): 3.57

Canister Dilution Factor: 1.37

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.034	ND	0.013	
107-06-2	1,2-Dichloroethane	<b>0.22</b>	0.034	<b>0.053</b>	0.0085	
71-43-2	Benzene	<b>1.1</b>	0.10	<b>0.35</b>	0.032	
79-01-6	Trichloroethene	ND	0.034	ND	0.0064	
79-00-5	1,1,2-Trichloroethane	ND	0.14	ND	0.025	
127-18-4	Tetrachloroethene	ND	0.034	ND	0.0051	
100-41-4	Ethylbenzene	<b>0.48</b>	0.14	<b>0.11</b>	0.032	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.034	ND	0.0050	
91-20-3	Naphthalene	ND	0.14	ND	0.026	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** 35-1F-032912  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P1201320-010

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:  
 Container ID: AC00943

Date Collected: 3/30/12  
 Date Received: 4/4/12  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Initial Pressure (psig): -4.85      Final Pressure (psig): 3.66

Canister Dilution Factor: 1.86

CAS #	Compound	Result µg/m <sup>3</sup>	MRL µg/m <sup>3</sup>	Result ppbV	MRL ppbV	Data Qualifier
75-01-4	Vinyl Chloride	ND	0.047	ND	0.018	
107-06-2	1,2-Dichloroethane	<b>0.35</b>	0.047	<b>0.087</b>	0.011	
71-43-2	Benzene	<b>2.7</b>	0.14	<b>0.84</b>	0.044	
79-01-6	Trichloroethene	ND	0.047	ND	0.0087	
79-00-5	1,1,2-Trichloroethane	ND	0.19	ND	0.034	
127-18-4	Tetrachloroethene	<b>2.4</b>	0.047	<b>0.36</b>	0.0069	
100-41-4	Ethylbenzene	<b>1.4</b>	0.19	<b>0.33</b>	0.043	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.047	ND	0.0068	
91-20-3	Naphthalene	<b>0.54</b>	0.19	<b>0.10</b>	0.035	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Method Blank  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
 CAS Sample ID: P120410-MB

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister  
 Test Notes:

Date Collected: NA  
 Date Received: NA  
 Date Analyzed: 4/10/12  
 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

CAS #	Compound	Result	MRL	Result	MRL	Data Qualifier
		µg/m <sup>3</sup>	µg/m <sup>3</sup>	ppbV	ppbV	
75-01-4	Vinyl Chloride	ND	0.025	ND	0.0098	
107-06-2	1,2-Dichloroethane	ND	0.025	ND	0.0062	
71-43-2	Benzene	ND	0.075	ND	0.023	
79-01-6	Trichloroethene	ND	0.025	ND	0.0047	
79-00-5	1,1,2-Trichloroethane	ND	0.10	ND	0.018	
127-18-4	Tetrachloroethene	ND	0.025	ND	0.0037	
100-41-4	Ethylbenzene	ND	0.10	ND	0.023	
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.025	ND	0.0036	
91-20-3	Naphthalene	ND	0.10	ND	0.019	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** Montana DEQ  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320

Test Code: EPA TO-15 SIM  
 Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
 Analyst: Karen Ryan  
 Sampling Media: 6.0 L Summa Canister(s)  
 Test Notes:

Date(s) Collected: 3/25 - 3/30/12  
 Date(s) Received: 4/4/12  
 Date(s) Analyzed: 4/10/12

Client Sample ID	CAS Sample ID	1,2-Dichloroethane-d4	Toluene-d8	Bromofluorobenzene	Acceptance Limits	Data Qualifier
		% Recovered	% Recovered	% Recovered		
Method Blank	P120410-MB	100	109	107	70-130	
Lab Control Sample	P120410-LCS	96	108	109	70-130	
31-1F-032412	P1201320-001	96	111	102	70-130	
37-1F-032612	P1201320-002	99	109	104	70-130	
29-1F-032612	P1201320-003	100	109	100	70-130	
DUP-54	P1201320-004	100	108	105	70-130	
30-1F-032612	P1201320-005	100	112	110	70-130	
33-1B-032712	P1201320-006	97	112	110	70-130	
32-1F-032812	P1201320-007	99	109	102	70-130	
36-1F-032812	P1201320-008	99	110	107	70-130	
34-1F-032912	P1201320-009	98	108	106	70-130	
35-1F-032912	P1201320-010	98	107	103	70-130	

Surrogate percent recovery is verified and accepted based on the on-column result.

Reported results are shown in concentration units and as a result of the calculation, may vary slightly from the on-column percent recovery.

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 1

**Client:** Montana DEQ  
**Client Sample ID:** Lab Control Sample  
**Client Project ID:** Background Indoor Air Study

CAS Project ID: P1201320  
CAS Sample ID: P120410-LCS

Test Code: EPA TO-15 SIM  
Instrument ID: Tekmar AUTOCAN/Agilent 5973N/HP6890A/MS7  
Analyst: Karen Ryan  
Sampling Media: 6.0 L Summa Canister  
Test Notes:

Date Collected: NA  
Date Received: NA  
Date Analyzed: 4/10/12  
Volume(s) Analyzed: 0.125 Liter(s)

CAS #	Compound	Spike Amount µg/m <sup>3</sup>	Result µg/m <sup>3</sup>	% Recovery	CAS	Data Qualifier
					Acceptance Limits	
75-01-4	Vinyl Chloride	3.88	<b>2.91</b>	<b>75</b>	56-127	
107-06-2	1,2-Dichloroethane	4.00	<b>2.90</b>	<b>73</b>	51-140	
71-43-2	Benzene	3.96	<b>3.49</b>	<b>88</b>	56-125	
79-01-6	Trichloroethene	3.88	<b>2.84</b>	<b>73</b>	51-127	
79-00-5	1,1,2-Trichloroethane	3.92	<b>2.77</b>	<b>71</b>	49-137	
127-18-4	Tetrachloroethene	3.68	<b>2.82</b>	<b>77</b>	58-134	
100-41-4	Ethylbenzene	3.96	<b>3.25</b>	<b>82</b>	56-126	
79-34-5	1,1,2,2-Tetrachloroethane	3.84	<b>2.92</b>	<b>76</b>	53-148	
91-20-3	Naphthalene	3.44	<b>3.12</b>	<b>91</b>	25-149	

Laboratory Control Sample percent recovery is verified and accepted based on the on-column result.  
Reported results are shown in concentration units and as a result of the calculation, may vary slightly.

Massachusetts APH  
Hydrocarbon Ranges

ICAL Method: M8011412A.M

ICAL Date: 1/14/12

Instrument ID: MS08

areas

	0.5	1	5	25	50	100	0.5	1	5	25	50	100
<b>C5-C8 Aliphatics</b>												
Isopentane	60478	112606	488243	2480105	4710733	9222808	0.520	1.04	5.20	26.0	52.0	104
n-Hexane	59953	110631	507126	2577865	4931102	9678199	0.490	0.98	4.90	24.5	49.0	98
Cyclohexane	63203	115230	548910	2874726	5524402	10890911	0.500	1.00	5.00	25.0	50.0	100
2,3-Dimethylpentane	68648	127222	585183	2998017	5742104	11319597	0.520	1.04	5.20	26.0	52.0	104
n-Heptane	66143	120286	558074	2884817	5509801	10873046	0.490	0.98	4.90	24.5	49.0	98
n-Octane	76169	137829	632891	3286017	6205035	12092514	0.505	1.01	5.05	25.3	50.5	101
area sum:	<b>394594</b>	<b>723804</b>	<b>3320427</b>	<b>17101547</b>	<b>32623177</b>	<b>64077075</b>	<b>3.025</b>	<b>6.050</b>	<b>30.25</b>	<b>151.30</b>	<b>302.50</b>	<b>605.00</b>

C9-C12 Aliphatics

	0.5	1	5	25	50	100	0.5	1	5	25	50	100
2,3-Dimethylheptane	78780	147724	682518	3501487	6691428	13148840	0.485	0.97	4.85	24.3	48.5	97
n-Nonane	77464	140680	662221	3417128	6555165	12940625	0.485	0.97	4.85	24.3	48.5	97
n-Decane	80593	147724	705259	3651028	7021424	13870530	0.490	0.98	4.90	24.5	49.0	98
n-Undecane	88915	166782	776903	4026931	7749936	15415340	0.495	0.99	4.95	24.8	49.5	99
n-Dodecane	80867	152557	719618	3772331	7263077	14385881	0.480	0.96	4.80	24.0	48.0	96
area sum:	<b>476664</b>	<b>899213</b>	<b>4233774</b>	<b>22042429</b>	<b>42396042</b>	<b>83802799</b>	<b>2.935</b>	<b>5.870</b>	<b>29.35</b>	<b>146.90</b>	<b>293.50</b>	<b>587.00</b>

C9-C10 Aromatics

	0.5	1	5	25	50	100	0.5	1	5	25	50	100
Isopropylbenzene	10658	19169	87439	448287	849124	1695737	0.485	0.97	4.85	24.3	48.5	97
3-Ethyltoluene	11805	20564	97823	504510	977020	1973178	0.470	0.94	4.70	23.5	47.0	94
1,3,5-Trimethylbenzene	15759	29260	134073	700133	1353305	2726630	0.490	0.98	4.90	24.5	49.0	98
1,2,3-Trimethylbenzene	17195	31328	143473	757946	1493373	3063059	0.495	0.99	4.95	24.8	49.5	99
p-Isopropyltoluene	9490	18116	83401	440962	872799	1814314	0.460	0.92	4.60	23.0	46.0	92
area sum:	<b>64907</b>	<b>118437</b>	<b>546209</b>	<b>2851838</b>	<b>5545621</b>	<b>11272918</b>	<b>2.400</b>	<b>4.800</b>	<b>24.00</b>	<b>120.10</b>	<b>240.00</b>	<b>480.00</b>

Massachusetts APH  
Hydrocarbon Ranges

ICAL: M8011412A.M

ICAL Date: 1/14/12

Instrument ID: MS08

areas

<u>Internal Standards (TIC)</u>	0.5	1	5	25	50	100
Bromochloromethane (IS1)	1312793	1307309	1293586	1289283	1284194	1283817
1,4-Difluorobenzene (IS2)	2547079	2541159	2512498	2475255	2422119	2352159
Chlorobenzene-d5 (IS3)	2929189	2913374	2880258	2895080	2878314	2908396

Internal Standards (EIC)

Bromochloromethane (IS1)	249696	247899	245985	245585	244327	243561
1,4-Difluorobenzene (IS2)	1153657	1149947	1133233	1136223	1127203	1138730
Chlorobenzene-d5 (IS3)	1007702	1004638	991212	992925	982812	986950

Surrogates (TIC)

	0.5	1	5	25	50	100
1,2-Dichloroethane-d4	1162497	1154617	1105630	1126201	1103958	1095945
88 Toluene-d8	3508018	3490493	3455382	3466688	3437986	3457323
88 p-Bromofluorobenzene	2298220	2284797	2197239	2227396	2192159	2197999

C5-C8 Aliphatics

	0.5	1	5	25	50	100	RRF <sub>avg</sub>	%RSD
	2.8268	2.6009	2.4215	2.4870	2.3919	2.3252	2.509	7.24

C9-C12 Aliphatics

	0.5	1	5	25	50	100	RRF <sub>avg</sub>	%RSD
	4.0291	3.8120	3.6383	3.7780	3.6744	3.6163	3.758	4.09

C9-C10 Aromatics

	0.5	1	5	25	50	100	RRF <sub>avg</sub>	%RSD
	0.6709	0.6140	0.5740	0.5979	0.5878	0.5949	0.607	5.63

**Massachusetts APH**  
**Continuing Calibration Verification Check Sheet**

Data File Name: 04071202.D  
 Data File Path: J:\MS08\Data\2012\_04\06\  
 Operator: EM  
 Date Acquired: 4/7/12 3:06  
 Acq. Method File: TO15.M  
 Sample Name: 25ng TO-15/MAPH CCV STD  
 Misc Info: S25-04031203/S25-03161202  
 Instrument Name: MS08

Enter RRFs from current ICAL!

<b>Internal Standards</b>		<u>RT</u>	<u>Area</u>						
7)	1,4-Difluorobenzene (IS2)	15.89	1082412						
16)	Chlorobenzene-d5 (IS3)	21.69	941687						
<b>C5-C8 Aliphatics</b>		<u>RT</u>	<u>Area</u>	<u>RRF</u>	<u>ng</u>	<u>% D</u>	<u>LCL</u>	<u>UCL</u>	<u>Pass/Fai</u>
3)	Isopentane	8.19	2388606	2.538	153.0	1.14	-30	30	Pass
4)	n-Hexane	13.08	2507742						
9)	Cyclohexane	15.81	2792389						
10)	2,3-Dimethylpentane	16.17	2922324	Spike	ICAL				
11)	n-Heptane	17.35	2804136	Amt (ng)	RRF				
14)	n-Octane	20.69	3208272	151.30	2.509				
			<b>16623469</b>						
<b>C9-C12 Aliphatics</b>		<u>RT</u>	<u>Area</u>	<u>RRF</u>	<u>ng</u>	<u>% D</u>	<u>LCL</u>	<u>UCL</u>	<u>Pass/Fai</u>
18)	2,3-Dimethylheptane	22.31	3401295	3.834	149.9	2.02	-30	30	Pass
19)	n-Nonane	23.30	3317221						
25)	n-Decane	25.26	3533579						
28)	Butylcyclohexane	25.99	3853565	Spike	ICAL				
29)	n-Undecane	26.76	3619770	Amt (ng)	RRF				
30)	n-Dodecane	28.00	3488040	146.90	3.758				
			<b>21213470</b>						
<b>C9-C10 Aromatics</b>		<u>RT</u>	<u>Area</u>	<u>RRF</u>	<u>ng</u>	<u>% D</u>	<u>LCL</u>	<u>UCL</u>	<u>Pass/Fai</u>
22)	Isopropylbenzene	23.78	439879	0.601	118.9	-1.03	-30	30	Pass
23)	1-Methyl-3-ethylbenzene	24.52	482296						
24)	1,3,5-Trimethylbenzene	24.66	665021						
26)	p-Isopropyltoluene	25.68	416582	Spike	ICAL				
27)	1,2,3-Trimethylbenzene	25.68	714045	Amt (ng)	RRF				
			<b>2717823</b>	120.1	0.607				

*Em 4/7/12*



**Massachusetts APH**  
**Continuing Calibration Verification Check Sheet**

Data File Name: 04091202.D  
 Data File Path: J:\MS08\Data\2012\_04\09\  
 Operator: EM  
 Date Acquired: 4/9/12 9:14  
 Acq. Method File: TO15.M  
 Sample Name: 25ng TO-15/MAPH CCV STD  
 Misc Info: S25-04031203/S25-03161202  
 Instrument Name: MS08

Enter RRFs from current ICAL!

Internal Standards	RT	Area
7) 1,4-Difluorobenzene (IS2)	15.89	1084690
16) Chlorobenzene-d5 (IS3)	21.69	936982

C5-C8 Aliphatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
3) Isopentane	8.20	2405833	2.533	152.7	0.94	-30	30	Pass
4) n-Hexane	13.08	2517847						
9) Cyclohexane	15.81	2766222						
10) 2,3-Dimethylpentane	16.17	2918072	Spike	ICAL				
11) n-Heptane	17.35	2801114	Amt (ng)	RRF				
14) n-Octane	20.69	3216100	151.30	2.509				
		<b>16625188</b>						

C9-C12 Aliphatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
18) 2,3-Dimethylheptane	22.31	3390681	3.837	150.0	2.09	-30	30	Pass
19) n-Nonane	23.30	3309192						
25) n-Decane	25.26	3535601						
28) Butylcyclohexane	25.99	3847831	Spike	ICAL				
29) n-Undecane	26.76	3623338	Amt (ng)	RRF				
30) n-Dodecane	28.00	3416131	146.90	3.758				
		<b>21122774</b>						

C9-C10 Aromatics	RT	Area	RRF	ng	% D	LCL	UCL	Pass/Fai
22) Isopropylbenzene	23.78	435892	0.602	119.1	-0.84	-30	30	Pass
23) 1-Methyl-3-ethylbenzene	24.52	480661						
24) 1,3,5-Trimethylbenzene	24.66	664726						
26) p-Isopropyltoluene	25.68	414020	Spike	ICAL				
27) 1,2,3-Trimethylbenzene	25.68	714015	Amt (ng)	RRF				
		<b>2709314</b>	120.1	0.607				

*EM 4/10/12*

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Mon Dec 12 09:45:45 2011  
 Response Via : Initial Calibration

Calibration Files

0.1 =12091118.D 0.2 =12091119.D 0.5 =12091120.D 1.0 =12091121.D 5.0 =12091122.D 25 =12091123.D  
 50 =12091124.D 100 =12091125.D

Compound	0.1	0.2	0.5	1.0	5.0	25	50	100	Avg	%RSD
1) IR Bromochloromethane...				ISTD						
2) T Propene	2.487	1.965	1.952	1.692	2.194	2.303	2.185	2.036	2.102	11.62
3) T Dichlorodifluo...	3.545	2.805	2.940	2.748	2.432	2.361	2.278	2.199	2.664	16.73
4) T Chloromethane	3.290	2.811	2.875	2.512	2.176	2.286	1.977	1.664	2.449	21.62
5) T 1,2-Dichloro-1...	1.901	1.661	1.575	1.458	1.359	1.333	1.200	1.190	1.460	16.68
6) T Vinyl Chloride	2.906	2.590	2.631	2.374	2.145	2.129	1.950	1.939	2.333	15.07
7) T 1,3-Butadiene	2.145	1.788	1.790	1.718	1.581	1.689	1.621	1.616	1.743	10.33
8) T Bromomethane	1.687	1.328	1.349	1.212	1.094	1.142	1.111	1.092	1.252	16.21
9) T Chloroethane	1.671	1.259	1.333	1.193	1.101	1.091	1.046	1.044	1.217	17.29
9.10) T Ethanol	2.040	1.345	1.186	1.085	1.022	1.096	1.076	1.064	1.239	27.34
11) T Acetonitrile	4.365	2.973	2.723	2.546	2.480	2.562	2.482	2.468	2.825	22.85
12) T Acrolein	1.328	0.931	0.907	0.861	0.772	0.814	0.789	0.779	0.897	20.48
13) T Acetone	1.645	1.316	1.249	1.129	0.994	1.015	0.981	0.964	1.161	20.26
14) T Trichlorofluor...	2.776	2.385	2.496	2.272	2.097	2.135	2.042	2.011	2.277	11.59
15) T 2-Propanol (Is...	4.463	4.536	4.168	2.920	2.763	2.577	2.672	3.443	26.09	26.09
16) T Acrylonitrile	2.701	1.907	1.878	1.783	1.745	1.813	1.765	1.738	1.916	16.85
17) T 1,1-Dichloroet...	1.458	1.418	1.452	1.263	1.206	1.223	1.172	1.152	1.293	9.95
18) T 2-Methyl-2-Pro...	4.464	3.710	3.939	3.616	3.558	3.584			3.812	9.14
19) T Methylene Chlo...	2.347	2.260	2.345	2.220	2.150	2.296	2.230	2.191	2.255	20.88
20) T 3-Chloro-1-pro...	1.348	1.244	1.299	1.193	1.089	1.098	1.062	1.035	1.171	3.15
21) T Trichlorotrifl...	6.210	5.678	5.057	4.654	4.699	4.488	4.424	5.030	13.40	10.01
22) T Carbon Disulfide	2.436	2.237	2.206	2.142	1.951	1.953	1.883	1.835	2.080	10.00
23) T trans-1,2-Dich...	2.896	2.757	2.747	2.500	2.307	2.379	2.288	2.236	2.514	10.07
24) T 1,1-Dichloroet...	5.128	4.611	4.650	4.307	4.119	4.185	4.042	3.984	4.378	8.93
25) T Methyl tert-Bu...	0.430	0.368	0.423	0.399	0.393	0.410	0.402	0.398	0.403	4.70
26) T Vinyl Acetate	1.265	1.003	1.038	0.972	0.927	0.964	0.928	0.794	0.986	13.56
27) T 2-Butanone (MEK)	2.442	2.056	2.101	1.938	1.826	1.836	1.764	1.720	1.960	12.05
28) T cis-1,2-Dichlo...	1.446	1.223	1.323	1.262	1.171	1.188	1.153	1.148	1.239	8.26
29) T Diisopropyl Ether	0.611	0.554	0.573	0.545	0.517	0.535	0.520	0.516	0.546	6.02
30) T Ethyl Acetate	3.541	3.006	3.181	2.805	2.582	2.633	2.545	2.543	2.855	12.70
31) T n-Hexane										

Response Factor Report MS08

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M

Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

Title	2.679	2.408	2.409	2.204	1.983	2.054	1.974	1.937	2.206	12.18
32) T Chloroform	2.679	2.408	2.409	2.204	1.983	2.054	1.974	1.937	2.206	12.18
33) S 1,2-Dichloroet...	1.586	1.577	1.582	1.590	1.535	1.533	1.520	1.504	1.553	2.19
34) T Tetrahydrofura...	1.274	0.932	1.100	0.938	0.910	0.879	0.830	0.821	0.961	15.96
35) T Ethyl tert-But...	1.999	1.832	2.033	1.818	1.732	1.757	1.708	1.687	1.821	7.18
36) T 1,2-Dichloroet...	2.099	1.992	1.976	1.806	1.683	1.700	1.599	1.531	1.798	11.39
37) IR 1,4-Difluorobenzen...										
38) T 1,1,1-Trichloror...	0.531	0.465	0.486	0.453	0.419	0.434	0.413	0.403	0.451	9.46
39) T Isopropyl Acetate	0.227	0.205	0.220	0.206	0.194	0.201	0.192	0.188	0.204	6.71
40) T 1-Butanol		0.348	0.319	0.300	0.320	0.335	0.324	0.318	0.323	4.66
41) T Benzene	1.633	1.385	1.359	1.227	1.136	1.158	1.101	1.070	1.258	15.11
42) T Carbon Tetrach...	0.394	0.356	0.369	0.340	0.334	0.352	0.337	0.333	0.352	6.03
43) T Cyclohexane	0.666	0.576	0.588	0.539	0.492	0.500	0.479	0.469	0.539	12.59
44) T tert-Amyl Meth...	1.092	0.926	0.998	0.937	0.900	0.918	0.881	0.858	0.939	7.95
45) T 1,2-Dichloropr...	0.397	0.361	0.368	0.324	0.312	0.316	0.300	0.293	0.334	11.12
46) T Bromodichlorom...	0.458	0.387	0.392	0.378	0.362	0.378	0.362	0.353	0.384	8.57
47) T Trichloroethene	0.466	0.375	0.379	0.344	0.323	0.328	0.316	0.309	0.355	14.62
48) T 1,4-Dioxane	0.305	0.282	0.288	0.248	0.245	0.246	0.237	0.230	0.260	10.55
49) T 2,2,4-Trimethy...	1.964	1.629	1.686	1.546	1.447	1.461	1.387	1.337	1.557	12.98
50) T Methyl Methacr...	0.156	0.139	0.150	0.135	0.136	0.142	0.136	0.135	0.141	5.53
51) T n-Heptane	0.427	0.373	0.385	0.364	0.341	0.344	0.332	0.324	0.361	9.35
52) T cis-1,3-Dichlo...	0.571	0.502	0.521	0.491	0.482	0.500	0.481	0.471	0.502	6.33
53) T 4-Methyl-2-pen...	0.359	0.295	0.308	0.303	0.297	0.303	0.289	0.284	0.305	7.68
54) T trans-1,3-Dich...	0.476	0.413	0.443	0.421	0.429	0.453	0.437	0.427	0.437	4.59
55) T 1,1,2-Trichlor...	0.340	0.301	0.310	0.284	0.276	0.285	0.271	0.265	0.292	8.44
56) IR Chlorobenzene-d5 (...)										
57) S Toluene-d8 (SS2)	2.305	2.306	2.305	2.309	2.306	2.300	2.310	2.304	2.306	0.14
58) T Toluene	3.623	2.920	3.025	2.738	2.571	2.574	2.473	2.421	2.793	14.19
59) T 2-Hexanone	2.049	1.429	1.515	1.403	1.395	1.422	1.367	1.337	1.490	15.57
60) T Dibromochlorom...	0.696	0.618	0.636	0.636	0.612	0.648	0.629	0.625	0.638	4.14
61) T 1,2-Dibromoethane	0.831	0.730	0.718	0.658	0.641	0.662	0.638	0.628	0.688	9.97
62) T n-Butyl Acetate	2.338	1.850	1.857	1.721	1.704	1.765	1.737	1.737	1.839	11.40
63) T n-Octane	0.785	0.652	0.714	0.643	0.602	0.611	0.580	0.572	0.645	11.28
64) T Tetrachloroethene	1.090	0.904	0.904	0.826	0.767	0.779	0.756	0.755	0.848	13.63
65) T Chlorobenzene	2.375	1.868	1.886	1.744	1.610	1.637	1.574	1.548	1.780	15.30
66) T Ethylbenzene	3.846	3.171	3.188	2.956	2.803	2.848	2.743	2.680	3.030	12.51
67) T m- & p-Xylenes	3.081	2.499	2.611	2.392	2.216	2.273	2.170	2.112	2.419	13.08
68) T Bromoform	0.622	0.494	0.556	0.541	0.551	0.598	0.589	0.594	0.568	7.19
69) T Styrene	2.478	1.895	1.937	1.800	1.735	1.803	1.726	1.675	1.881	13.64
70) T o-Xylene	3.279	2.708	2.726	2.479	2.331	2.373	2.283	2.224	2.550	13.65

Method Path : J:\MS08\Methods\  
 Method File : R8120911A.M

Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

71)	T	n-Nonane	2.005	1.753	1.847	1.675	1.581	1.597	1.518	1.470	1.681	10.69
72)	T	1,1,2,2-Tetrac...	1.348	1.164	1.262	1.150	1.083	1.121	1.085	1.066	1.160	8.48
73)	S	Bromofluoroben...	0.774	0.771	0.779	0.781	0.775	0.778	0.782	0.780	0.777	0.50
74)	T	Cumene	4.522	3.495	3.524	3.169	2.962	3.002	2.884	2.831	3.299	16.98
75)	T	alpha-Pinene	1.782	1.636	1.734	1.601	1.537	1.578	1.518	1.498	1.611	6.35
76)	T	n-Propylbenzene	5.441	4.229	4.306	3.924	3.710	3.754	3.553	3.394	4.039	16.01
77)	T	3-Ethyltoluene	4.091	3.373	3.461	3.038	2.964	2.981	2.893	2.845	3.206	13.14
78)	T	4-Ethyltoluene	4.003	3.202	3.301	3.151	2.837	2.983	2.810	2.776	3.133	12.84
79)	T	1,3,5-Trimethy...	3.324	2.662	2.723	2.537	2.378	2.426	2.344	2.403	2.600	12.43
80)	T	alpha-Methylst...	1.742	1.386	1.449	1.371	1.390	1.402	1.354	1.399	1.436	8.80
81)	T	2-Ethyltoluene	4.143	3.362	3.430	3.169	2.975	3.048	2.919	2.985	3.254	12.43
82)	T	1,2,4-Trimethy...	3.356	2.663	2.791	2.551	2.433	2.513	2.454	2.548	2.664	11.37
83)	T	n-Decane	1.969	1.702	1.844	1.705	1.582	1.626	1.554	1.609	1.699	8.39
84)	T	Benzyl Chloride	2.373	1.648	1.772	1.688	1.799	2.058	2.054	2.190	1.948	13.35
85)	T	1,3-Dichlorobe...	2.498	1.658	1.649	1.455	1.377	1.409	1.374	1.436	1.607	23.49
86)	T	1,4-Dichlorobe...	2.655	1.785	1.618	1.492	1.380	1.428	1.388	1.449	1.649	25.99
87)	T	sec-Butylbenzene	4.527	3.773	3.816	3.471	3.264	3.328	3.214	3.176	3.571	12.80
88)	T	4-Isopropyltol...	4.426	3.674	3.683	3.416	3.254	3.356	3.253	3.366	3.554	10.98
889)	T	1,2,3-Trimethy...	3.266	2.650	2.852	2.588	2.471	2.573	2.506	2.619	2.691	9.63
890)	T	1,2-Dichlorobe...	2.173	1.589	1.545	1.435	1.332	1.387	1.370	1.415	1.531	17.89
891)	T	d-Limonene	1.124	0.998	1.106	1.069	1.064	1.072	1.040	1.065	1.067	3.58
92)	T	1,2-Dibromo-3-...	0.563	0.488	0.501	0.486	0.494	0.525	0.532	0.542	0.517	5.50
93)	T	n-Undecane	1.922	1.637	1.755	1.671	1.608	1.630	1.608	1.592	1.678	6.63
94)	T	1,2,4-Trichlor...		1.335	1.165	1.062	1.017	1.071	1.086	1.118	1.122	9.32
95)	T	Naphthalene		4.848	3.923	3.638	3.497	3.721	3.755	3.783	3.881	11.50
96)	T	n-Dodecane	1.866	1.599	1.669	1.586	1.545	1.575	1.553	1.509	1.613	6.99
97)	T	Hexachlorobuta...	0.841	0.749	0.730	0.673	0.635	0.664	0.663	0.676	0.704	9.50
98)	T	Cyclohexanone	1.172	0.979	1.021	0.969	0.952	0.962	0.931	0.903	0.986	8.37
99)	T	tert-Butylbenzene	3.486	2.799	2.880	2.658	2.449	2.533	2.453	2.528	2.723	12.71
100)	T	n-Butylbenzene	3.596	2.904	2.958	2.670	2.572	2.625	2.613	2.593	2.817	12.33

(#) = Out of Range

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_04\07\  
 Data File : 04071201.D  
 Acq On : 7 Apr 2012 2:23  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-04031203/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 07 07:01:39 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	IR Bromochloromethane (IS1)	1.000	1.000	0.0	82	-0.02
2	T Propene	2.102	2.305	-9.7	83	0.00
3	T Dichlorodifluoromethane (CF	2.664	2.312	13.2	81	0.00
4	T Chloromethane	2.449	2.172	11.3	78	0.00
5	T 1,2-Dichloro-1,1,2,2-tetra	1.460	1.212	17.0	75	0.00
6	T Vinyl Chloride	2.333	2.037	12.7	79	0.00
7	T 1,3-Butadiene	1.743	1.733	0.6	85	0.00
8	T Bromomethane	1.252	1.113	11.1	80	0.00
9	T Chloroethane	1.217	1.092	10.3	83	0.00
10	T Ethanol	1.239	1.182	4.6	89	-0.07
11	T Acetonitrile	2.825	2.653	6.1	85	-0.05
12	T Acrolein	0.897	0.829	7.6	84	-0.02
13	T Acetone	1.161	1.067	8.1	87	-0.05
14	T Trichlorofluoromethane	2.277	2.119	6.9	82	0.00
15	T 2-Propanol (Isopropanol)	3.443	2.761	19.8	82	-0.05
16	T Acrylonitrile	1.916	1.842	3.9	84	-0.03
17	T 1,1-Dichloroethene	1.293	1.185	8.4	80	0.00
18	T 2-Methyl-2-Propanol (tert-B	3.812	3.075	19.3	71	-0.03
19	T Methylene Chloride	1.421	1.250	12.0	81	-0.01
20	T 3-Chloro-1-propene (Allyl C	2.255	2.354	-4.4	85	-0.02
21	T Trichlorotrifluoroethane	1.171	1.055	9.9	79	0.00
22	T Carbon Disulfide	5.030	4.592	8.7	81	0.00
23	T trans-1,2-Dichloroethene	2.080	1.974	5.1	83	-0.01
24	T 1,1-Dichloroethane	2.514	2.417	3.9	84	-0.02
25	T Methyl tert-Butyl Ether	4.378	4.248	3.0	84	0.00
26	T Vinyl Acetate	0.403	0.414	-2.7	83	-0.03
27	T 2-Butanone (MEK)	0.986	0.977	0.9	84	-0.03
28	T cis-1,2-Dichloroethene	1.960	1.856	5.3	83	-0.01
29	T Diisopropyl Ether	1.239	1.196	3.5	83	-0.01
30	T Ethyl Acetate	0.546	0.542	0.7	84	-0.03
31	T n-Hexane	2.855	2.627	8.0	82	-0.01
32	T Chloroform	2.206	2.031	7.9	82	-0.03
33	S 1,2-Dichloroethane-d4 (SS1)	1.553	1.635	-5.3	88	-0.02
34	T Tetrahydrofuran (THF)	0.961	0.905	5.8	85	-0.01
35	T Ethyl tert-Butyl Ether	1.821	1.755	3.6	82	-0.01
36	T 1,2-Dichloroethane	1.798	1.708	5.0	83	-0.02
37	IR 1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	87	-0.01
38	T 1,1,1-Trichloroethane	0.451	0.412	8.6	82	-0.01

*Em* 4/7/12

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_04\07\  
 Data File : 04071201.D  
 Acq On : 7 Apr 2012 2:23  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-04031203/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 07 07:01:39 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
39 T	Isopropyl Acetate	0.204	0.194	4.9	83	-0.02
40 T	1-Butanol	0.323	0.356	-10.2	92	-0.05
41 T	Benzene	1.258	1.097	12.8	82	-0.02
42 T	Carbon Tetrachloride	0.352	0.326	7.4	80	-0.01
43 T	Cyclohexane	0.539	0.482	10.6	83	-0.02
44 T	tert-Amyl Methyl Ether	0.939	0.875	6.8	82	-0.02
45 T	1,2-Dichloropropane	0.334	0.303	9.3	83	-0.02
46 T	Bromodichloromethane	0.384	0.357	7.0	82	-0.01
47 T	Trichloroethene	0.355	0.302	14.9	80	-0.02
48 T	1,4-Dioxane	0.260	0.238	8.5	84	-0.02
49 T	2,2,4-Trimethylpentane (Iso	1.557	1.423	8.6	84	-0.01
50 T	Methyl Methacrylate	0.141	0.134	5.0	82	-0.03
51 T	n-Heptane	0.361	0.336	6.9	84	-0.02
52 T	cis-1,3-Dichloropropene	0.502	0.476	5.2	82	-0.01
53 T	4-Methyl-2-pentanone	0.305	0.298	2.3	85	-0.02
54 T	trans-1,3-Dichloropropene	0.437	0.429	1.8	82	-0.01
55 T	1,1,2-Trichloroethane	0.292	0.268	8.2	81	-0.02
56 IR	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	87	0.00
57 S	Toluene-d8 (SS2)	2.306	2.271	1.5	86	0.00
58 T	Toluene	2.793	2.427	13.1	82	-0.01
59 T	2-Hexanone	1.490	1.418	4.8	87	-0.02
60 T	Dibromochloromethane	0.638	0.594	6.9	80	-0.01
61 T	1,2-Dibromoethane	0.688	0.609	11.5	80	-0.01
62 T	n-Butyl Acetate	1.839	1.752	4.7	86	-0.02
63 T	n-Octane	0.645	0.620	3.9	88	-0.01
64 T	Tetrachloroethene	0.848	0.702	17.2	78	0.00
65 T	Chlorobenzene	1.780	1.527	14.2	81	0.00
66 T	Ethylbenzene	3.030	2.719	10.3	83	0.00
67 T	m- & p-Xylenes	2.419	2.179	9.9	83	-0.02
68 T	Bromoform	0.568	0.540	4.9	79	-0.01
69 T	Styrene	1.881	1.741	7.4	84	-0.01
70 T	o-Xylene	2.550	2.268	11.1	83	-0.01
71 T	n-Nonane	1.681	1.549	7.9	84	-0.01
72 T	1,1,2,2-Tetrachloroethane	1.160	1.053	9.2	82	-0.02
73 S	Bromofluorobenzene (SS3)	0.777	0.774	0.4	87	0.00
74 T	Cumene	3.299	2.975	9.8	86	-0.01
75 T	alpha-Pinene	1.611	1.486	7.8	82	-0.01
76 T	n-Propylbenzene	4.039	3.560	11.9	82	-0.01

*Em* 4/7/12

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_04\07\  
 Data File : 04071201.D  
 Acq On : 7 Apr 2012 2:23  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-04031203/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 07 07:01:39 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
77 T	3-Ethyltoluene	3.206	2.844	11.3	83	-0.01
78 T	4-Ethyltoluene	3.133	2.708	13.6	79	-0.02
79 T	1,3,5-Trimethylbenzene	2.600	2.285	12.1	82	-0.01
80 T	alpha-Methylstyrene	1.436	1.322	7.9	82	-0.01
81 T	2-Ethyltoluene	3.254	2.841	12.7	81	-0.01
82 T	1,2,4-Trimethylbenzene	2.664	2.354	11.6	81	-0.02
83 T	n-Decane	1.699	1.553	8.6	83	-0.02
84 T	Benzyl Chloride	1.948	2.052	-5.3	87	-0.02
85 T	1,3-Dichlorobenzene	1.607	1.281	20.3	79	-0.01
86 T	1,4-Dichlorobenzene	1.649	1.304	20.9	79	-0.02
87 T	sec-Butylbenzene	3.571	3.140	12.1	82	-0.01
88 T	4-Isopropyltoluene (p-Cymen	3.554	3.145	11.5	82	-0.01
89 T	1,2,3-Trimethylbenzene	2.691	2.401	10.8	81	-0.01
90 T	1,2-Dichlorobenzene	1.531	1.260	17.7	79	-0.01
91 T	d-Limonene	1.067	1.048	1.8	85	-0.01
92 T	1,2-Dibromo-3-Chloropropane	0.517	0.480	7.2	80	-0.01
93 T	n-Undecane	1.678	1.570	6.4	84	-0.01
94 T	1,2,4-Trichlorobenzene	1.122	0.972	13.4	79	-0.01
95 T	Naphthalene	3.881	3.485	10.2	81	-0.01
96 T	n-Dodecane	1.613	1.578	2.2	87	0.00
97 T	Hexachlorobutadiene	0.704	0.592	15.9	78	0.00
98 T	Cyclohexanone	0.986	0.996	-1.0	90	-0.02
99 T	tert-Butylbenzene	2.723	2.389	12.3	82	-0.02
100 T	n-Butylbenzene	2.817	2.452	13.0	81	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*EM 4/7/12*

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_04\09\  
 Data File : 04091201.D  
 Acq On : 9 Apr 2012 8:31  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-04031203/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 09 09:37:31 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	IR Bromochloromethane (IS1)	1.000	1.000	0.0	85	-0.02
2	T Propene	2.102	1.988	5.4	73	0.00
3	T Dichlorodifluoromethane (CF	2.664	2.317	13.0	83	0.00
4	T Chloromethane	2.449	2.268	7.4	84	0.00
5	T 1,2-Dichloro-1,1,2,2-tetra	1.460	1.239	15.1	79	0.00
6	T Vinyl Chloride	2.333	2.103	9.9	84	0.00
7	T 1,3-Butadiene	1.743	1.776	-1.9	89	0.00
8	T Bromomethane	1.252	1.094	12.6	81	0.00
9	T Chloroethane	1.217	1.068	12.2	83	0.00
10	T Ethanol	1.239	1.175	5.2	91	-0.07
11	T Acetonitrile	2.825	2.582	8.6	85	-0.05
12	T Acrolein	0.897	0.815	9.1	85	-0.02
13	T Acetone	1.161	1.051	9.5	88	-0.05
14	T Trichlorofluoromethane	2.277	2.083	8.5	83	0.00
15	T 2-Propanol (Isopropanol)	3.443	3.492	-1.4	107	-0.05
16	T Acrylonitrile	1.916	1.810	5.5	85	-0.03
17	T 1,1-Dichloroethene	1.293	1.183	8.5	82	0.00
18	T 2-Methyl-2-Propanol (tert-B	3.812	3.307	13.2	78	-0.03
19	T Methylene Chloride	1.421	1.225	13.8	81	-0.01
20	T 3-Chloro-1-propene (Allyl C	2.255	2.381	-5.6	88	-0.02
21	T Trichlorotrifluoroethane	1.171	1.041	11.1	80	0.00
22	T Carbon Disulfide	5.030	4.651	7.5	84	0.00
23	T trans-1,2-Dichloroethene	2.080	1.949	6.3	85	-0.02
24	T 1,1-Dichloroethane	2.514	2.391	4.9	85	-0.02
25	T Methyl tert-Butyl Ether	4.378	4.214	3.7	85	-0.01
26	T Vinyl Acetate	0.403	0.413	-2.5	85	-0.03
27	T 2-Butanone (MEK)	0.986	0.968	1.8	85	-0.03
28	T cis-1,2-Dichloroethene	1.960	1.838	6.2	85	-0.02
29	T Diisopropyl Ether	1.239	1.180	4.8	84	-0.02
30	T Ethyl Acetate	0.546	0.537	1.6	85	-0.03
31	T n-Hexane	2.855	2.627	8.0	84	-0.02
32	T Chloroform	2.206	2.007	9.0	83	-0.03
33	S 1,2-Dichloroethane-d4 (SS1)	1.553	1.644	-5.9	91	-0.02
34	T Tetrahydrofuran (THF)	0.961	0.893	7.1	86	-0.02
35	T Ethyl tert-Butyl Ether	1.821	1.753	3.7	84	-0.02
36	T 1,2-Dichloroethane	1.798	1.705	5.2	85	-0.02
37	IR 1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	88	-0.01
38	T 1,1,1-Trichloroethane	0.451	0.411	8.9	83	-0.01



Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_04\09\  
 Data File : 04091201.D  
 Acq On : 9 Apr 2012 8:31  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-04031203/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 09 09:37:31 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
39 T	Isopropyl Acetate	0.204	0.196	3.9	85	-0.03
40 T	1-Butanol	0.323	0.354	-9.6	93	-0.05
41 T	Benzene	1.258	1.101	12.5	84	-0.02
42 T	Carbon Tetrachloride	0.352	0.325	7.7	81	-0.02
43 T	Cyclohexane	0.539	0.479	11.1	84	-0.02
44 T	tert-Amyl Methyl Ether	0.939	0.877	6.6	84	-0.02
45 T	1,2-Dichloropropane	0.334	0.303	9.3	84	-0.02
46 T	Bromodichloromethane	0.384	0.358	6.8	83	-0.01
47 T	Trichloroethene	0.355	0.297	16.3	80	-0.02
48 T	1,4-Dioxane	0.260	0.237	8.8	85	-0.02
49 T	2,2,4-Trimethylpentane (Iso	1.557	1.432	8.0	86	-0.02
50 T	Methyl Methacrylate	0.141	0.134	5.0	83	-0.03
51 T	n-Heptane	0.361	0.338	6.4	86	-0.02
52 T	cis-1,3-Dichloropropene	0.502	0.480	4.4	84	-0.01
53 T	4-Methyl-2-pentanone	0.305	0.304	0.3	88	-0.02
54 T	trans-1,3-Dichloropropene	0.437	0.432	1.1	84	-0.01
55 T	1,1,2-Trichloroethane	0.292	0.265	9.2	82	-0.02
56 IR	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	88	0.00
57 S	Toluene-d8 (SS2)	2.306	2.307	-0.0	88	0.00
58 T	Toluene	2.793	2.439	12.7	84	-0.01
59 T	2-Hexanone	1.490	1.449	2.8	90	-0.02
60 T	Dibromochloromethane	0.638	0.597	6.4	81	-0.01
61 T	1,2-Dibromoethane	0.688	0.615	10.6	82	-0.01
62 T	n-Butyl Acetate	1.839	1.784	3.0	89	-0.02
63 T	n-Octane	0.645	0.621	3.7	90	-0.01
64 T	Tetrachloroethene	0.848	0.701	17.3	79	-0.01
65 T	Chlorobenzene	1.780	1.514	14.9	82	0.00
66 T	Ethylbenzene	3.030	2.768	8.6	86	-0.01
67 T	m- & p-Xylenes	2.419	2.181	9.8	85	-0.02
68 T	Bromoform	0.568	0.545	4.0	80	-0.01
69 T	Styrene	1.881	1.734	7.8	85	-0.01
70 T	o-Xylene	2.550	2.276	10.7	85	-0.02
71 T	n-Nonane	1.681	1.568	6.7	87	-0.01
72 T	1,1,2,2-Tetrachloroethane	1.160	1.055	9.1	83	-0.02
73 S	Bromofluorobenzene (SS3)	0.777	0.754	3.0	85	0.00
74 T	Cumene	3.299	2.963	10.2	87	-0.01
75 T	alpha-Pinene	1.611	1.495	7.2	84	-0.01
76 T	n-Propylbenzene	4.039	3.572	11.6	84	-0.01

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_04\09\  
 Data File : 04091201.D  
 Acq On : 9 Apr 2012 8:31  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-04031203/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 09 09:37:31 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
77 T	3-Ethyltoluene	3.206	2.767	13.7	82	-0.01
78 T	4-Ethyltoluene	3.133	2.782	11.2	82	-0.02
79 T	1,3,5-Trimethylbenzene	2.600	2.267	12.8	82	-0.01
80 T	alpha-Methylstyrene	1.436	1.314	8.5	83	-0.02
81 T	2-Ethyltoluene	3.254	2.834	12.9	82	-0.01
82 T	1,2,4-Trimethylbenzene	2.664	2.344	12.0	82	-0.02
83 T	n-Decane	1.699	1.566	7.8	85	-0.02
84 T	Benzyl Chloride	1.948	2.097	-7.6	90	-0.02
85 T	1,3-Dichlorobenzene	1.607	1.273	20.8	80	-0.02
86 T	1,4-Dichlorobenzene	1.649	1.273	22.8	79	-0.02
87 T	sec-Butylbenzene	3.571	3.138	12.1	83	-0.02
88 T	4-Isopropyltoluene (p-Cymen)	3.554	3.134	11.8	82	-0.01
89 T	1,2,3-Trimethylbenzene	2.691	2.416	10.2	83	-0.01
90 T	1,2-Dichlorobenzene	1.531	1.248	18.5	79	-0.02
91 T	d-Limonene	1.067	1.051	1.5	86	-0.02
92 T	1,2-Dibromo-3-Chloropropane	0.517	0.485	6.2	81	-0.01
93 T	n-Undecane	1.678	1.574	6.2	85	-0.01
94 T	1,2,4-Trichlorobenzene	1.122	0.928	17.3	76	-0.01
95 T	Naphthalene	3.881	3.358	13.5	80	-0.01
96 T	n-Dodecane	1.613	1.547	4.1	87	0.00
97 T	Hexachlorobutadiene	0.704	0.575	18.3	76	0.00
98 T	Cyclohexanone	0.986	1.009	-2.3	93	-0.02
99 T	tert-Butylbenzene	2.723	2.380	12.6	83	-0.02
100 T	n-Butylbenzene	2.817	2.457	12.8	83	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*Em 4/10/12*

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_04\10\  
 Data File : 04101201.D  
 Acq On : 10 Apr 2012 6:17  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-04031203/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 10 07:33:42 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
1 IR	Bromochloromethane (IS1)	1.000	1.000	0.0	82	-0.01
2 T	Propene	2.102	2.242	-6.7	79	0.00
3 T	Dichlorodifluoromethane (CF	2.664	2.319	13.0	80	0.00
4 T	Chloromethane	2.449	2.130	13.0	76	0.00
5 T	1,2-Dichloro-1,1,2,2-tetra	1.460	1.223	16.2	75	0.00
6 T	Vinyl Chloride	2.333	2.070	11.3	79	0.00
7 T	1,3-Butadiene	1.743	1.755	-0.7	85	0.00
8 T	Bromomethane	1.252	1.098	12.3	78	0.00
9 T	Chloroethane	1.217	1.080	11.3	81	0.00
10 T	Ethanol	1.239	1.168	5.7	87	-0.07
11 T	Acetonitrile	2.825	2.603	7.9	83	-0.05
12 T	Acrolein	0.897	0.817	8.9	82	-0.02
13 T	Acetone	1.161	1.045	10.0	84	-0.04
14 T	Trichlorofluoromethane	2.277	2.073	9.0	79	0.00
15 T	2-Propanol (Isopropanol)	3.443	2.874	16.5	85	-0.05
16 T	Acrylonitrile	1.916	1.809	5.6	81	-0.03
17 T	1,1-Dichloroethene	1.293	1.180	8.7	79	0.00
18 T	2-Methyl-2-Propanol (tert-B	3.812	3.165	17.0	72	-0.03
19 T	Methylene Chloride	1.421	1.244	12.5	80	0.00
20 T	3-Chloro-1-propene (Allyl C	2.255	2.321	-2.9	82	-0.01
21 T	Trichlorotrifluoroethane	1.171	1.047	10.6	78	0.00
22 T	Carbon Disulfide	5.030	4.594	8.7	80	0.00
23 T	trans-1,2-Dichloroethene	2.080	1.949	6.3	81	-0.01
24 T	1,1-Dichloroethane	2.514	2.370	5.7	81	-0.01
25 T	Methyl tert-Butyl Ether	4.378	4.163	4.9	81	0.00
26 T	Vinyl Acetate	0.403	0.407	-1.0	81	-0.03
27 T	2-Butanone (MEK)	0.986	0.963	2.3	81	-0.02
28 T	cis-1,2-Dichloroethene	1.960	1.818	7.2	81	-0.01
29 T	Diisopropyl Ether	1.239	1.180	4.8	81	0.00
30 T	Ethyl Acetate	0.546	0.533	2.4	81	-0.02
31 T	n-Hexane	2.855	2.640	7.5	82	0.00
32 T	Chloroform	2.206	1.998	9.4	79	-0.03
33 S	1,2-Dichloroethane-d4 (SS1)	1.553	1.632	-5.1	87	-0.01
34 T	Tetrahydrofuran (THF)	0.961	0.881	8.3	82	0.00
35 T	Ethyl tert-Butyl Ether	1.821	1.737	4.6	81	-0.01
36 T	1,2-Dichloroethane	1.798	1.688	6.1	81	-0.02
37 IR	1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	85	0.00
38 T	1,1,1-Trichloroethane	0.451	0.407	9.8	80	0.00

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_04\10\  
 Data File : 04101201.D  
 Acq On : 10 Apr 2012 6:17  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-04031203/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 10 07:33:42 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
39 T	Isopropyl Acetate	0.204	0.192	5.9	81	-0.02
40 T	1-Butanol	0.323	0.351	-8.7	89	-0.05
41 T	Benzene	1.258	1.099	12.6	81	-0.01
42 T	Carbon Tetrachloride	0.352	0.321	8.8	78	-0.01
43 T	Cyclohexane	0.539	0.476	11.7	81	-0.01
44 T	tert-Amyl Methyl Ether	0.939	0.871	7.2	81	-0.01
45 T	1,2-Dichloropropane	0.334	0.299	10.5	80	-0.01
46 T	Bromodichloromethane	0.384	0.354	7.8	80	0.00
47 T	Trichloroethene	0.355	0.299	15.8	78	-0.01
48 T	1,4-Dioxane	0.260	0.237	8.8	82	-0.02
49 T	2,2,4-Trimethylpentane (Iso	1.557	1.411	9.4	82	-0.01
50 T	Methyl Methacrylate	0.141	0.132	6.4	79	-0.02
51 T	n-Heptane	0.361	0.332	8.0	82	-0.01
52 T	cis-1,3-Dichloropropene	0.502	0.473	5.8	81	0.00
53 T	4-Methyl-2-pentanone	0.305	0.299	2.0	84	-0.02
54 T	trans-1,3-Dichloropropene	0.437	0.426	2.5	80	-0.01
55 T	1,1,2-Trichloroethane	0.292	0.261	10.6	78	-0.01
56 IR	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	85	0.00
57 S	Toluene-d8 (SS2)	2.306	2.284	1.0	85	0.00
58 T	Toluene	2.793	2.389	14.5	79	0.00
59 T	2-Hexanone	1.490	1.403	5.8	84	-0.02
60 T	Dibromochloromethane	0.638	0.586	8.2	77	0.00
61 T	1,2-Dibromoethane	0.688	0.604	12.2	78	-0.01
62 T	n-Butyl Acetate	1.839	1.731	5.9	84	-0.01
63 T	n-Octane	0.645	0.616	4.5	86	-0.01
64 T	Tetrachloroethene	0.848	0.697	17.8	76	0.00
65 T	Chlorobenzene	1.780	1.509	15.2	79	0.00
66 T	Ethylbenzene	3.030	2.669	11.9	80	0.00
67 T	m- & p-Xylenes	2.419	2.150	11.1	81	-0.02
68 T	Bromoform	0.568	0.538	5.3	77	-0.01
69 T	Styrene	1.881	1.716	8.8	81	-0.01
70 T	o-Xylene	2.550	2.247	11.9	81	-0.01
71 T	n-Nonane	1.681	1.523	9.4	81	-0.01
72 T	1,1,2,2-Tetrachloroethane	1.160	1.035	10.8	79	-0.02
73 S	Bromofluorobenzene (SS3)	0.777	0.772	0.6	85	0.00
74 T	Cumene	3.299	2.940	10.9	84	0.00
75 T	alpha-Pinene	1.611	1.473	8.6	80	0.00
76 T	n-Propylbenzene	4.039	3.520	12.8	80	0.00

*Em 4/10/12*

Evaluate Continuing Calibration Report

Data Path : J:\MS08\Data\2012\_04\10\  
 Data File : 04101201.D  
 Acq On : 10 Apr 2012 6:17  
 Operator : EM  
 Sample : 25ng TO-15 CCV STD  
 Misc : S25-04031203/S25-03191204  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 10 07:33:42 2012  
 Quant Method : J:\MS08\Methods\R8120911A.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Mon Dec 12 09:45:45 2011  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
77 T	3-Ethyltoluene	3.206	2.771	13.6	79	-0.01
78 T	4-Ethyltoluene	3.133	2.718	13.2	78	-0.01
79 T	1,3,5-Trimethylbenzene	2.600	2.246	13.6	79	-0.01
80 T	alpha-Methylstyrene	1.436	1.310	8.8	80	-0.01
81 T	2-Ethyltoluene	3.254	2.814	13.5	79	-0.01
82 T	1,2,4-Trimethylbenzene	2.664	2.306	13.4	78	-0.01
83 T	n-Decane	1.699	1.542	9.2	81	-0.01
84 T	Benzyl Chloride	1.948	2.028	-4.1	84	-0.02
85 T	1,3-Dichlorobenzene	1.607	1.269	21.0	77	-0.01
86 T	1,4-Dichlorobenzene	1.649	1.282	22.3	77	-0.02
87 T	sec-Butylbenzene	3.571	3.095	13.3	79	-0.01
88 T	4-Isopropyltoluene (p-Cymen)	3.554	3.092	13.0	79	-0.01
89 T	1,2,3-Trimethylbenzene	2.691	2.371	11.9	79	-0.01
90 T	1,2-Dichlorobenzene	1.531	1.242	18.9	77	-0.01
91 T	d-Limonene	1.067	1.026	3.8	82	-0.01
92 T	1,2-Dibromo-3-Chloropropane	0.517	0.478	7.5	78	-0.01
93 T	n-Undecane	1.678	1.544	8.0	81	0.00
94 T	1,2,4-Trichlorobenzene	1.122	0.949	15.4	76	-0.01
95 T	Naphthalene	3.881	3.417	12.0	78	-0.01
96 T	n-Dodecane	1.613	1.535	4.8	83	0.00
97 T	Hexachlorobutadiene	0.704	0.579	17.8	75	0.00
98 T	Cyclohexanone	0.986	0.977	0.9	87	-0.02
99 T	tert-Butylbenzene	2.723	2.358	13.4	80	-0.01
100 T	n-Butylbenzene	2.817	2.428	13.8	79	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

*EM 4/10/12*

Response Factor Report MS07

Method : J:\Ms07\METHODS\X7021712B.M (RTE Integrator)  
Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
Last Update : Fri Mar 02 08:26:43 2012  
Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	AVG	%RSD
1) I Bromochloromethan					ISTD						
2) T Dichlorodifluorom	3.784	3.527	3.467	3.822	3.279	2.986	2.875	3.155	2.920	3.313	10.83
3) T Chloromethane		1.470	0.956	1.190	0.904	0.778	0.569	0.897	0.827	0.949	28.81
4) T Vinyl Chloride	2.802	2.749	2.588	3.044	2.597	2.406	2.314	2.588	2.446	2.615	8.57
5) T Bromomethane	1.894	1.496	1.211	1.613	1.335	1.208	1.105	1.433	1.367	1.407	17.11
6) T Chloroethane	1.511	1.416	1.357	1.517	1.295	1.237	1.195	1.312	1.224	1.340	8.92
7) T Acetone				1.548	1.214	1.122	1.080	1.193	1.090	1.208	14.52
8) T Trichlorofluorome	3.051	2.887	2.831	3.145	2.692	2.565	2.447	2.691	2.536	2.761	8.59
9) T 1,1-Dichloroethen	1.591	1.440	1.366	1.547	1.322	1.229	1.214	1.385	1.295	1.376	9.52
10) T Methylene Chlorid		2.134	1.795	1.866	1.536	1.426	1.404	1.494	1.411	1.633	16.43
11) T Trichlorotrifluor	1.408	1.325	1.295	1.376	1.169	1.081	1.062	1.170	1.097	1.220	10.87
12) T trans-1,2-Dichlor	1.972	1.710	1.589	1.762	1.505	1.406	1.393	1.512	1.439	1.588	12.18
13) T 1,1-Dichloroethan	3.265	3.020	3.016	3.383	2.925	2.756	2.681	2.937	2.841	2.980	7.60
14) T Methyl tert-Butyl	5.168	4.740	4.762	5.045	4.335	4.137	4.166	4.690	4.471	4.613	7.90
15) T cis-1,2-Dichloroe	1.885	1.626	1.630	1.776	1.524	1.434	1.388	1.575	1.482	1.591	10.09
16) T Chloroform				3.041	3.257	2.586	2.390	2.524	2.374	2.636	14.03
17) S 1,2-Dichloroethan	2.000	2.002	2.032	2.018	1.989	1.973	1.960	1.938	1.932	1.983	1.75
18) T 1,2-Dichloroethan	2.582	2.460	2.375	2.604	2.222	2.090	1.988	2.222	2.094	2.293	9.73
19) T 1,1,1-Trichloroet	2.587	2.431	2.340	2.573	2.191	2.059	2.007	2.278	2.150	2.291	9.18
20) T Benzene				7.536	7.534	6.055	5.665	5.903	5.468	6.222	14.88
21) T Carbon Tetrachlor	2.280	1.842	1.769	1.950	1.676	1.579	1.549	1.771	1.692	1.790	12.41

22) I 1,4-Difluorobenze					ISTD						
23) T 1,2-Dichloropropa	0.417	0.397	0.373	0.409	0.354	0.331	0.319	0.358	0.336	0.366	9.68
24) T Bromodichlorometh	0.491	0.464	0.436	0.489	0.420	0.398	0.394	0.447	0.416	0.439	8.26
25) T Trichloroethene	0.418	0.379	0.349	0.390	0.324	0.305	0.287	0.326	0.307	0.343	12.96
26) T 1,4-Dioxane	0.324	0.290	0.270	0.278	0.242	0.231	0.227	0.258	0.243	0.263	11.99
27) T cis-1,3-Dichlorop	0.567	0.523	0.508	0.540	0.499	0.484	0.488	0.559	0.522	0.521	5.67
28) T trans-1,3-Dichlor	0.457	0.434	0.431	0.444	0.419	0.409	0.417	0.505	0.482	0.444	7.22

(#) = Out of Range ## Number of calibration levels exceeded format ##

Response Factor Report MS07

Method : J:\Ms07\METHODS\X7021712B.M (RTE Integrator)  
 Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 Last Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Calibration Files

10 =02171204.D 25 =02171205.D 75 =02171206.D 100 =02171207.D 500 =02171208.D  
 1000 =02171209.D 2500 =02171210.D 9999 =02171211.D 20K =02171212.D

Compound	10	25	75	100	500	1000	2500	9999	20K	Avg	%RSD
29) T 1,1,2-Trichloroet	0.315	0.296	0.286	0.305	0.263	0.248	0.242	0.274	0.257	0.276	9.32
30) S Toluene-d8 (SS2)	1.053	1.071	1.054	1.047	1.051	1.052	1.071	1.060	1.049	1.056	0.84
31) T Toluene	1.782	1.541	1.420	1.485	1.269	1.198	1.177	1.307	1.186	1.374	14.76
32) T 1,2-Dibromoethane	0.379	0.362	0.345	0.362	0.323	0.307	0.304	0.351	0.331	0.340	7.65
33) T Tetrachloroethene	0.423	0.394	0.371	0.400	0.341	0.320	0.311	0.347	0.326	0.359	11.03
-----ISTD-----											
34) I Chlorobenzene-d5											
35) T Chlorobenzene	4.581	4.172	4.017	4.308	3.688	3.469	3.311	3.705	3.345	3.844	11.68
36) T Ethylbenzene	7.604	7.082	6.707	7.235	6.297	6.070	5.915	6.682	5.878	6.608	9.30
37) T m,p-Xylene	5.861	5.347	5.132	5.536	4.930	4.777	4.719	5.268	4.558	5.125	8.24
38) T o-Xylene	6.505	5.826	5.526	5.889	5.188	4.995	4.950	5.594	4.932	5.490	9.67
39) T 1,1,2,2-Tetrachlo	2.655	2.674	2.710	2.646	2.437	2.371	2.298	2.761	2.498	2.561	6.40
40) S Bromofluorobenze	1.712	1.701	1.695	1.697	1.683	1.680	1.670	1.642	1.593	1.675	2.20
41) T 1,3-Dichlorobenze	3.304	3.261	3.222	3.204	2.835	2.719	2.608	3.138	2.846	3.015	8.71
42) T 1,4-Dichlorobenze	3.470	3.303	3.208	3.208	2.803	2.671	2.579	3.174	2.891	3.034	10.13
43) T 1,2-Dichlorobenze	3.271	3.123	3.112	3.026	2.662	2.536	2.418	3.053	2.772	2.886	10.31
44) T 1,2,4-Trichlorobe	2.184	2.118	2.169	1.753	1.641	1.608	1.575	2.139	1.978	1.907	13.66
45) T Naphthalene	6.857	6.355	6.704	5.143	5.107	5.172	5.297	8.126	7.328	6.232	17.82
46) T Hexachlorobutadie	1.486	1.423	1.403	1.217	1.055	1.013	0.975	1.227	1.145	1.216	15.40

(#) = Out of Range ### Number of calibration levels exceeded format ###  
 X7021712B.M Tue Apr 10 10:05:50 2012

Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_04\10\  
 Data File : 04101203.D  
 Acq On : 10 Apr 2012 9:21 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00163  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Apr 10 10:05:18 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
1 I	Bromochloromethane (IS1)	1.000	1.000	0.0	78	0.00
2 T	Dichlorodifluoromethane (CF)	3.313	2.565	22.6	61	0.00
3 T	Chloromethane	0.949	0.735	22.6	63	0.00
4 T	Vinyl Chloride	2.615	2.063	21.1	62	0.00
5 T	Bromomethane	1.407	1.209	14.1	71	0.00
6 T	Chloroethane	1.340	1.091	18.6	66	0.00
7 T	Acetone	1.208	1.070	11.4	69	0.00
8 T	Trichlorofluoromethane	2.761	2.217	19.7	64	0.00
9 T	1,1-Dichloroethene	1.376	1.163	15.5	69	0.00
10 T	Methylene Chloride	1.633	1.296	20.6	66	0.00
11 T	Trichlorotrifluoroethane	1.220	1.078	11.6	72	0.00
12 T	trans-1,2-Dichloroethene	1.588	1.296	18.4	67	0.00
13 T	1,1-Dichloroethane	2.980	2.472	17.0	66	0.00
14 T	Methyl tert-Butyl Ether	4.613	3.623	21.5	65	0.00
15 T	cis-1,2-Dichloroethene	1.591	1.314	17.4	67	0.00
16 T	Chloroform	2.636	2.183	17.2	66	0.00
17 S	1,2-Dichloroethane-d4 (SS1)	1.983	1.970	0.7	77	0.00
18 T	1,2-Dichloroethane	2.293	1.741	24.1	61	0.00
19 T	1,1,1-Trichloroethane	2.291	1.853	19.1	66	0.00
20 T	Benzene	6.222	5.270	15.3	68	0.00
21 T	Carbon Tetrachloride	1.790	1.442	19.4	67	0.00
22 I	1,4-Difluorobenzene (IS2)	1.000	1.000	0.0	88	0.00
23 T	1,2-Dichloropropane	0.366	0.260	29.0	65	0.00
24 T	Bromodichloromethane	0.439	0.306	30.3#	64	0.00
25 T	Trichloroethene	0.343	0.253	26.2	69	0.00
26 T	1,4-Dioxane	0.263	0.199	24.3	73	0.00
27 T	cis-1,3-Dichloropropene	0.521	0.396	24.0	70	0.00
28 T	trans-1,3-Dichloropropene	0.444	0.339	23.6	71	0.00
29 T	1,1,2-Trichloroethane	0.276	0.200	27.5	67	0.00
30 S	Toluene-d8 (SS2)	1.056	1.154	-9.3	97	0.00
31 T	Toluene	1.374	1.027	25.3	72	0.00
32 T	1,2-Dibromoethane	0.340	0.257	24.4	70	0.00
33 T	Tetrachloroethene	0.359	0.284	20.9	74	0.00
34 I	Chlorobenzene-d5 (IS3)	1.000	1.000	0.0	84	0.00
35 T	Chlorobenzene	3.844	3.248	15.5	74	0.00
36 T	Ethylbenzene	6.608	5.549	16.0	74	0.00
37 T	m,p-Xylene	5.125	4.350	15.1	74	0.00

KR 4/10/12



Evaluate Continuing Calibration Report

Data Path : J:\Ms07\DATA\2012\_04\10\  
 Data File : 04101203.D  
 Acq On : 10 Apr 2012 9:21 am  
 Operator : KR  
 Sample : 500pg TO-15 SIM CCV STD  
 Misc : STD00161/STD00163  
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Apr 10 10:05:18 2012  
 Quant Method : J:\Ms07\METHODS\X7021712B.M  
 Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)  
 QLast Update : Fri Mar 02 08:26:43 2012  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev (min)
38 T	o-Xylene	5.490	4.592	16.4	74	0.00
39 T	1,1,2,2-Tetrachloroethane	2.561	2.087	18.5	72	0.00
40 S	Bromofluorobenzene (SS3)	1.675	1.858	-10.9	93	0.00
41 T	1,3-Dichlorobenzene	3.015	2.719	9.8	80	0.00
42 T	1,4-Dichlorobenzene	3.034	2.710	10.7	81	0.00
43 T	1,2-Dichlorobenzene	2.886	2.620	9.2	83	0.00
44 T	1,2,4-Trichlorobenzene	1.907	1.807	5.2	92	0.00
45 T	Naphthalene	6.232	6.463	-3.7	106	0.00
46 T	Hexachlorobutadiene	1.216	1.083	10.9	86	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

# Appendix B

Data validation was performed for six analytical laboratory reports issued by Columbia Analytical Services identified as P1200947, P1201004, P1201043, P1201074, P1201121 and P1201247. Indoor air samples were analyzed for EPA Methods TO-15, TO-15 SIM and air-phase petroleum hydrocarbons (APH). Laboratory quality assurance/quality control (QA/QC) provided for each analysis was evaluated. All holding times were met for each analysis and method blank results were non-detect for all analyses. Laboratory control sample (LCS) spike recoveries were met for all analyses and surrogate spike recoveries were acceptable for the TO-15 methods. All laboratory duplicate results were acceptable. A small number of results required qualification due to field duplicate results. Co-elution of peaks was observed in two samples with high concentrations, however, this did not affect the quality of the data. Overall, the laboratory results demonstrated an acceptable level of accuracy (LCS and surrogate results) and precision (duplicate results). The few samples requiring qualification are all considered usable results.

**Columbia Analytical Services**  
**Service Request # P1200947**

Sample ID	Chemical	Reported	Qualified			
		Conc. (ug/m <sup>3</sup> )	Lab Flag	Value (ug/m <sup>3</sup> )	Flag	Note
03-1F-030512	n-Nonane	3.6		3.6	J	(1)
01-1F-030512	methylene chloride	0.83		0.83	J	(2)
	vinyl acetate	8.1	U	8.1	UJ	(3)
	methyl methacrylate	1.9		1.9	J	(4)
	n-butyl acetate	4.3		4.3	J	(4)
	vinyl chloride	0.048		0.048	J	(2)
	TCE	0.29		0.29	J	(4)
	naphthalene	0.36		0.36	J	(4)
	C9-C12 aliphatics	21		21	J	(4)
	C9-C10 aromatics	12		12	J	(2)
51-1F-030512	methylene chloride	0.86	U	0.86	UJ	(5)
	vinyl acetate	18		18	J	(6)
	methyl methacrylate	3.2		3.2	J	(4)
	n-butyl acetate	3		3	J	(4)
	vinyl chloride	0.043	U	0.043	UJ	(5)
	TCE	0.47		0.47	J	(4)
	naphthalene	0.21		0.21	J	(4)
	C9-C12 aliphatics	40		40	J	(4)
	C9-C10 aromatics	8.6	U	8.6	UJ	(5)

Notes:

- (1): Detected value is qualified as estimated and flagged with a J qualifier due to coelution with a non-target compound. The result may be biased high.
- (2): Value is qualified as estimated and flagged with a J qualifier as the chemical was detected in the parent sample but nondetect in the field duplicate. The result may be biased high.
- (3): Nondetect value is qualified as estimated and flagged with a UJ qualifier as the chemical was not detected in the parent sample but detected in the field duplicate. The result may be biased low.
- (4): Value is qualified as estimated and flagged with a J qualifier based on a high relative percent difference (RPD) calculated between parent and field duplicate samples.
- (5): Nondetect value is qualified as estimated and flagged with a UJ qualifier as the chemical was not detected in the field duplicate but detected in the parent sample. The result may be biased low.
- (6): Value is qualified as estimated and flagged with a J qualifier as the chemical was detected in the field duplicate but nondetect in the parent sample. The result may be biased high.

**Columbia Analytical Services**  
**Service Request # P1201043**

Sample ID	Chemical	Reported	Qualified			
		Conc. (ug/m <sup>3</sup> )	Lab Flag	Value (ug/m <sup>3</sup> )	Flag	Note
15-1F-031212	Benzene	0.78		0.78	J	(1)
	m,p-Xylenes	1.6		1.6	J	(1)
	Tetrachloroethene	0.098		0.098	J	(2)
	C9-C12 aliphatics	15	U	15	UJ	(3)

52-1F-031212	Benzene	0.88	U	0.88	UJ	(4)
	m,p-Xylenes	1.8	U	1.8	UJ	(4)
	Tetrachloroethene	0.14		0.14	J	(2)
	C9-C12 aliphatics	19		19	J	(5)

Notes:

- (1): Value is qualified as estimated and flagged with a J qualifier as the chemical was detected in the sample but nondetect in the laboratory duplicate. The result may be biased high.
- (2): Value is qualified as estimated and flagged with a J qualifier based on a high relative percent difference (RPD) calculated between parent and field duplicate samples.
- (3): Nondetect value is qualified as estimated and flagged with a UJ qualifier as the chemical was not detected in the parent sample but detected in the field duplicate. The result may be biased low.
- (4): Nondetect value is qualified as estimated and flagged with a UJ qualifier as the chemical was not detected in the field duplicate but detected in the parent sample. The result may be biased low.
- (5): Value is qualified as estimated and flagged with a J qualifier as the chemical was detected in the field duplicate but nondetect in the parent sample. The result may be biased high.

**Columbia Analytical Services**  
**Service Request # P1201121**

Sample ID	Chemical	Reported		Qualified		
		Conc. (ug/m <sup>3</sup> )	Lab Flag	Value (ug/m <sup>3</sup> )	Flag	Note
26-1F-031912	2-Propanol	11		11	J	(1)
19-1F-031412	Propene	2.4		2.4	J	(2)
	2-Propanol	12		12	J	(2)
	Tetrachloroethene	0.070		0.070	J	(2)
	C9-C12 aliphatics	17	U	17	UJ	(3)
53-1F-031412	Propene	0.72	U		UJ	(4)
	2-Propanol	1.4	U		UJ	(4)
	Tetrachloroethene	0.036	U		UJ	(4)
	C9-C12 aliphatics	20			J	(5)

Notes:

- (1): Detected value is qualified as estimated and flagged with a J qualifier due to coelution with a non-target compound. The result may be biased high.
- (2): Value is qualified as estimated and flagged with a J qualifier as the chemical was detected in the sample but nondetect in the laboratory duplicate. The result may be biased high.
- (3): Nondetect value is qualified as estimated and flagged with a UJ qualifier as the chemical was not detected in the parent sample but detected in the field duplicate. The result may be biased low.
- (4): Nondetect value is qualified as estimated and flagged with a UJ qualifier as the chemical was not detected in the field duplicate but detected in the parent sample. The result may be biased low.
- (5): Value is qualified as estimated and flagged with a J qualifier as the chemical was detected in the field duplicate but nondetect in the parent sample. The result may be biased high.

**Columbia Analytical Services**  
**Service Request # P1201319**

Sample ID	Chemical	Reported	Qualified			
		Conc. (ug/m <sup>3</sup> )	Lab Flag	Value (ug/m <sup>3</sup> )	Flag	Note
40-1F-032812	Trichlorotrifluoroethane	0.88	U	0.88	UJ	(1)
	Naphthalene	0.42		0.42	J	(2)
	1,2-Dichloroethane	0.36		0.36	J	(2)
	C9-C12 aliphatics	30		30	J	(2)
55-1F-032812	Trichlorotrifluoroethane	0.82		0.82	J	(3)
	Naphthalene	0.30		0.30	J	(2)
	1,2-Dichloroethane	0.24		0.24	J	(2)
	C9-C12 aliphatics	19		19	J	(2)

Notes:

- (1): Nondetect value is qualified as estimated and flagged with a UJ qualifier as the chemical was not detected in the parent sample but detected in the field duplicate. The result may be biased low.
- (2): Value is qualified as estimated and flagged with a J qualifier based on a high relative percent difference (RPD) calculated between parent and field duplicate samples.
- (3): Value is qualified as estimated and flagged with a J qualifier as the chemical was detected in the field duplicate but nondetect in the parent sample. The result may be biased high.

**Columbia Analytical Services**  
**Service Request # P1201320**

Sample ID	Chemical	Reported	Qualified			
		Conc. (ug/m <sup>3</sup> )	Lab Flag	Value (ug/m <sup>3</sup> )	Flag	Note
29-1F-032612	Propene	5.0		5.0	J	(1)
	2-Propanol	21		21	J	(1)
	Chloromethane	0.77		0.77	J	(2)
	Ethyl acetate	1.6		1.6	J	(2)
	n-Nonane	0.88		0.88	J	(2)
	d-Limonene	11		11	J	(1)
	Naphthalene	0.92		0.92	J	(1)
DUP-54	Propene	12		12	J	(1)
	2-Propanol	67		67	J	(1)
	Chloromethane	0.96	U	0.96	UJ	(3)
	Ethyl acetate	1.9	U	1.9	UJ	(3)
	n-Nonane	0.96	U	0.96	UJ	(3)
	d-Limonene	7.0		7.0	J	(1)
	Naphthalene	0.63		0.63	J	(1)

Notes:

- (1): Value is qualified as estimated and flagged with a J qualifier based on a high relative percent difference (RPD) calculated between parent and field duplicate samples.
- (2): Value is qualified as estimated and flagged with a J qualifier as the chemical was detected in the sample but nondetect in the laboratory duplicate. The result may be biased high.
- (3): Nondetect value is qualified as estimated and flagged with a UJ qualifier as the chemical was not detected in the field duplicate but detected in the parent sample. The result may be biased low.

# Appendix C

**1,1,1-Trichloroethane**



Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C2	50	0.67	11	1.134	0.84	2.178	1.476	0.119	6.404	42.98	1.301

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C2	50	0.71	0.729	0.768	0.773	0.84	0.93	0.946	1.04	1.72	7.178

**1,1,2-Trichloroethane**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.13	0.38	0.176	0.165	0.00184	0.0429	0.0222	3.089	11.54	0.244

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.14	0.149	0.15	0.15	0.165	0.188	0.19	0.2	0.25	0.341

**1,1,2-Trichlorotrifluoroethane**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**1,1-Dichloroethane**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**1,1-Dichloroethene**



Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.868	26.86	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.93	0.93	0.982	1.275	2.563

**1,2,4-Trichlorobenzene**

From File: 1,2,4-TCB.wst

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

## **1,2,4-Trimethylbenzene**

From File: P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\1,2,4-TMB\1,2,4-TMB data.wst

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
1,2,4-TMB	50	0.67	17	2.233	1.05	8.952	2.992	0.467	3.331	12.59	1.34

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
1,2,4-TMB	50	0.719	0.767	0.828	0.863	1.05	1.875	2.14	4.91	8.665	13.57

**1,2-Dibromoethane**

From File: P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\1,2-DBA\1,2-DBA data.wst

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**1,2-Dichloroethane**



Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\1,2-DCA\1,2-DCA data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

C0

Raw Statistics

Number of Valid Observations 50  
Number of Distinct Observations 36  
Minimum 0.053  
Maximum 2.1  
Mean of Raw Data 0.374  
Standard Deviation of Raw Data 0.446  
Kstar 1.084  
Mean of Log Transformed Data -1.482  
Standard Deviation of Log Transformed Data 0.966

Normal Distribution Test Results

Correlation Coefficient R 0.824  
Shapiro Wilk Test Statistic 0.687  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 5.239E-13  
Lilliefors Test Statistic 0.236  
Lilliefors Critical (0.95) Value 0.125

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.97  
A-D Test Statistic 1.955  
A-D Critical (0.95) Value 0.776  
K-S Test Statistic 0.186  
K-S Critical(0.95) Value 0.129

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.974  
Shapiro Wilk Test Statistic 0.931  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 0.00754  
Lilliefors Test Statistic 0.146  
Lilliefors Critical (0.95) Value 0.125

Data not Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C0	50	0.053	2.1	0.374	0.17	0.199	0.446	0.134	2.458	6.662	1.192

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C0	50	0.0675	0.0839	0.0942	0.11	0.17	0.478	0.572	0.82	1.155	2.051

**Nonparametric Background Statistics for Full Data Sets**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\1,2-DCA\1,2-DCA data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Number of Bootstrap Operations	2000

C0

**Some Non-Parametric Statistics**

Number of Valid Observations	50
Number of Distinct Observations	36
Minimum	0.053
Maximum	2.1
Second Largest	2
Mean	0.374
Geometric Mean	0.227
First Quartile	0.11
Median	0.17
Third Quartile	0.478
SD	0.446
Variance	0.199
Coefficient of Variation	1.192
Skewness	2.458
Mean of Log-Transformed data	-1.482
SD of Log-Transformed data	0.966

**Data do not follow a Discernable Distribution (0.05)**

**Non-Parametric Background Statistics**

90% Percentile	0.82
95% Percentile	1.155
99% Percentile	2.051

**95% UTL with 50% Coverage**

Order Statistic	30
Achieved CC	0.941
UTL	0.23

95% BCA Bootstrap UTL with 50% Coverage	0.285
95% Percentile Bootstrap UTL with 50% Coverage	0.35

95% UPL	1.56
95% Chebyshev UPL	2.339

Upper Limit Based upon IQR	1.029
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From File: C:\Documents and Settings\martichbj\Desktop\Data Stats\1,2-DCA\1,2-DCA data.wst

Summary Statistics for Raw Full Data Sets

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
AG	27	0.067	2.1	0.456	0.35	0.213	0.462	0.341	2.051	5.186	1.013
DG	23	0.053	2	0.279	0.13	0.174	0.417	0.083	3.544	14.06	1.496

Percentiles for Raw Full Data Sets

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
AG	27	0.088	0.104	0.12	0.125	0.35	0.6	0.676	1.04	1.17	1.866
DG	23	0.0554	0.0692	0.0862	0.0885	0.13	0.275	0.372	0.478	0.768	1.736

Lognormal Background Statistics for Full Data Sets

User Selected Options

From File C:\Documents and Settings\martichb\Desktop\Data Stats\1,2-DCA\1,2-DCA data.wst  
 Full Precision OFF  
 Confidence Coefficient 95%  
 Coverage 50%  
 Different or Future K Values 1  
 Number of Bootstrap Operations 2000

AG

Log-Transformed Statistics

Number of Valid Observations 27  
 Number of Distinct Observations 23  
 Minimum -2.703  
 Maximum 0.742  
 Second Largest 0.182  
 Mean -1.212  
 First Quartile -2.08  
 Median -1.05  
 Third Quartile -0.511  
 SD 0.943

Lognormal Distribution Test

Shapiro Wilk Test Statistic 0.949  
 5% Shapiro Wilk Critical Value 0.923

Data appear Lognormal at 5% Significance Level

Background Statistics Assuming Lognormal Distribution

90% Percentile (z) 0.996  
 95% Percentile (z) 1.402  
 99% Percentile (z) 2.666  
 95% UPL 1.53  
 Tolerance Factor K N/A  
 95% UTL with 50% Coverage N/A

Some Nonparametric Background Statistics

95% Chebyshev UPL 2.506  
 95% Bootstrap BCA UTL with 50% Coverage 0.4  
 95% Percentile Bootstrap UTL with 50% Coverage 0.5

Nonparametric Background Statistics for Full Data Sets

User Selected Options

From File C:\Documents and Settings\martichb\Desktop\Data Stats1,2-DCA\1,2-DCA data.wst  
 Full Precision OFF  
 Confidence Coefficient 95%  
 Coverage 50%  
 Number of Bootstrap Operations 2000

DG

Some Non-Parametric Statistics

Number of Valid Observations	23
Number of Distinct Observations	19
Minimum	0.053
Maximum	2
Second Largest	0.8
Mean	0.279
Geometric Mean	0.166
First Quartile	0.0885
Median	0.13
Third Quartile	0.275
SD	0.417
Variance	0.174
Coefficient of Variation	1.496
Skewness	3.544
Mean of Log-Transformed data	-1.798
SD of Log-Transformed data	0.914

Data do not follow a Discernable Distribution (0.05)

Non-Parametric Background Statistics

90% Percentile	0.478
95% Percentile	0.768
99% Percentile	1.736

95% UTL with 50% Coverage

Order Statistic	15
Achieved CC	0.953
UTL	0.17

95% BCA Bootstrap UTL with 50% Coverage	0.17
95% Percentile Bootstrap UTL with 50% Coverage	0.19

95% UPL	1.76
95% Chebyshev UPL	2.136

Upper Limit Based upon IQR	0.555
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Wilcoxon-Mann-Whitney Site vs Background Comparison Test for Full Data Sets without NDs

User Selected Options

From File C:\Documents and Settings\martichbj\Desktop\Data Stats\1,2-DCA\1,2-DCA data.wst  
Full Precision OFF  
Confidence Coefficient 95%  
Substantial Difference 0.000  
Selected Null Hypothesis Site or AOC Mean/Median Less Than or Equal to Background Mean/Median (Form 1)  
Alternative Hypothesis Site or AOC Mean/Median Greater Than Background Mean/Median

Area of Concern Data: AG

Background Data: DG

Raw Statistics

	Site	Background
Number of Valid Observations	27	23
Number of Distinct Observations	23	19
Minimum	0.067	0.053
Maximum	2.1	2
Mean	0.456	0.279
Median	0.35	0.13
SD	0.462	0.417
SE of Mean	0.0889	0.087

Wilcoxon-Mann-Whitney (WMW) Test

H0: Mean/Median of Site or AOC  $\leq$  Mean/Median of Background

Site Rank Sum W-Stat: 805  
WMW Test U-Stat: 2.258  
WMW Critical Value (0.050): 1.645  
P-Value: 0.012

Conclusion with Alpha = 0.05

Reject H0, Conclude Site > Background

P-Value < alpha (0.05)

**1,2-Dichlorobenzene**



Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**1,2-Dichloropropane**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**1,3,5-Trimethylbenzene**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	5.5	1.171	0.865	0.757	0.87	0.126	3.297	12.6	0.743

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.77	0.78	0.865	0.97	1.08	1.98	2.91	4.373

**1,3-Butadiene**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.92	0.835	0.147	0.383	0.126	4.848	26.73	0.417

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.835	0.93	0.932	0.991	1.275	2.563

**1,3-Dichlorobenzene**



Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**1,4-Dichlorobenzene**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	36	2.143	0.835	37.62	6.133	0.126	4.982	24.45	2.862

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.835	0.93	0.946	1.05	2.615	31.59

**1,4-Dioxane**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**2-Butanone**

From File: 2-butanone.wst

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	6.7	32	9.908	8.5	21.39	4.625	1.26	3.391	12.7	0.467

Percentiles for Raw Full Dataset

Percentiles for Raw F

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	7.1	7.29	7.6	7.7	8.5	9.625	9.84	13.2	17.65	29.06

**2-Hexanone**



From File: P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\2-Hexanone\2-hexanone data.wsl

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.933	0.84	0.15	0.388	0.119	4.636	24.85	0.416

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.725	0.739	0.768	0.77	0.84	0.93	0.932	0.982	1.455	2.563

## **2-Propanol**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\2-Propanol\2-propanol data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

C0

Raw Statistics

Number of Valid Observations 50  
Number of Distinct Observations 43  
Minimum 3.1  
Maximum 170  
Mean of Raw Data 35.94  
Standard Deviation of Raw Data 33.66  
Kstar 1.204  
Mean of Log Transformed Data 3.137  
Standard Deviation of Log Transformed Data 1.02

Normal Distribution Test Results

Correlation Coefficient R 0.909  
Shapiro Wilk Test Statistic 0.833  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 2.6966E-7  
Lilliefors Test Statistic 0.165  
Lilliefors Critical (0.95) Value 0.125

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.996  
A-D Test Statistic 0.304  
A-D Critical (0.95) Value 0.773  
K-S Test Statistic 0.0766  
K-S Critical(0.95) Value 0.128

Data appear Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.988  
Shapiro Wilk Test Statistic 0.96  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 0.156  
Lilliefors Test Statistic 0.089  
Lilliefors Critical (0.95) Value 0.125

Data appear Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C0	50	3.1	170	35.94	27	1133	33.66	28.17	1.813	4.341	0.937

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C0	50	4.27	5.56	7.86	11.25	27	46	56.2	76.5	94.85	145.5

**Lognormal Background Statistics for Full Data Sets**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\2-Propanol\2-propanol data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1
Number of Bootstrap Operations	2000

C0

**Log-Transformed Statistics**

Number of Valid Observations	50
Number of Distinct Observations	43
Minimum	1.131
Maximum	5.136
Second Largest	4.787
Mean	3.137
First Quartile	2.42
Median	3.296
Third Quartile	3.829
SD	1.02

**Lognormal Distribution Test**

Shapiro Wilk Test Statistic	0.96
5% Shapiro Wilk Critical Value	0.947

**Data appear Lognormal at 5% Significance Level**

**Background Statistics Assuming Lognormal Distribution**

90% Percentile (z)	85.19
95% Percentile (z)	123.4
99% Percentile (z)	247.4
95% UPL	129.7
Tolerance Factor K	0.236
95% UTL with 50% Coverage	29.32

**Some Nonparametric Background Statistics**

95% Chebyshev UPL	184.1
95% Bootstrap BCA UTL with 50% Coverage	36
95% Percentile Bootstrap UTL with 50% Coverage	39

**4-Ethyltoluene**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	5.4	1.163	0.865	0.737	0.858	0.133	3.262	12.25	0.738

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.77	0.78	0.865	0.97	1.04	1.97	2.91	4.322

**4-Methyl-2-pentanone**



Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.3	1.113	0.875	0.368	0.607	0.156	2.301	4.978	0.545

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.724	0.749	0.778	0.783	0.875	1	1.16	2.1	2.255	3.251

**Acetone**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Acetone\Acetone data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

C0

Raw Statistics

Number of Valid Observations 50  
Number of Distinct Observations 40  
Minimum 16  
Maximum 130  
Mean of Raw Data 51.58  
Standard Deviation of Raw Data 26.84  
Kstar 3.92  
Mean of Log Transformed Data 3.818  
Standard Deviation of Log Transformed Data 0.508

Normal Distribution Test Results

Correlation Coefficient R 0.954  
Shapiro Wilk Test Statistic 0.902  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 3.1846E-4  
Lilliefors Test Statistic 0.115  
Lilliefors Critical (0.95) Value 0.125

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.992  
A-D Test Statistic 0.214  
A-D Critical (0.95) Value 0.754  
K-S Test Statistic 0.0603  
K-S Critical(0.95) Value 0.126

Data appear Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.997  
Shapiro Wilk Test Statistic 0.979  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 0.695  
Lilliefors Test Statistic 0.0585  
Lilliefors Critical (0.95) Value 0.125

Data appear Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C0	50	16	130	51.58	46	720.1	26.84	23.72	1.181	1.425	0.52

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C0	50	19.45	23.9	29	30.5	46	65	71.2	83.4	102.8	130

Lognormal Background Statistics for Full Data Sets

User Selected Options

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Acetone\Acetone data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1
Number of Bootstrap Operations	2000

C0

Log-Transformed Statistics

Number of Valid Observations	50
Number of Distinct Observations	40
Minimum	2.773
Maximum	4.868
Second Largest	4.868
Mean	3.818
First Quartile	3.417
Median	3.828
Third Quartile	4.174
SD	0.508

Lognormal Distribution Test

Shapiro Wilk Test Statistic	0.979
5% Shapiro Wilk Critical Value	0.947

Data appear Lognormal at 5% Significance Level

Background Statistics Assuming Lognormal Distribution

90% Percentile (z)	87.27
95% Percentile (z)	105
99% Percentile (z)	148.4
95% UPL	107.6
Tolerance Factor K	0.236
95% UTL with 50% Coverage	51.31

Some Nonparametric Background Statistics

95% Chebyshev UPL	169.7
95% Bootstrap BCA UTL with 50% Coverage	53
95% Percentile Bootstrap UTL with 50% Coverage	53

**Acetonitrile**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.971	0.835	0.255	0.505	0.119	3.807	14.67	0.52

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.715	0.739	0.768	0.77	0.835	0.93	0.932	1	1.81	3.2

**Acrolein**



Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	2.7	13	3.894	3.5	2.902	1.703	0.593	3.861	17.48	0.437

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	2.845	2.9	3.08	3.1	3.5	3.9	4	4.8	6.835	10.8

**Acrylonitrile**

From File: P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Acrylonitrile\Acrylonitrile data.wst

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

## **Allyl Chloride**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**alpha-Pinene**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\alpha-Pinene\alpha-Pinene data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

Raw

Raw Statistics

Number of Valid Observations 48  
Number of Missing Values 2  
Number of Distinct Observations 34  
Minimum 0.83  
Maximum 52  
Mean of Raw Data 6.765  
Standard Deviation of Raw Data 9.23  
Kstar 1.003  
Mean of Log Transformed Data 1.368  
Standard Deviation of Log Transformed Data 0.993

Normal Distribution Test Results

Correlation Coefficient R 0.77  
Shapiro Wilk Test Statistic 0.618  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 1.399E-14  
Lilliefors Test Statistic 0.3  
Lilliefors Critical (0.95) Value 0.128

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.946  
A-D Test Statistic 1.711  
A-D Critical (0.95) Value 0.777  
K-S Test Statistic 0.199  
K-S Critical(0.95) Value 0.131

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.981  
Shapiro Wilk Test Statistic 0.952  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 0.0802  
Lilliefors Test Statistic 0.118  
Lilliefors Critical (0.95) Value 0.128

Data appear Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
LnROS_Pinene	50	0.429	52	6.535	3.75	83.02	9.112	3.089	3.284	12.92	1.394

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
LnROS_Pinene	50	0.985	1.1	1.5	1.65	3.75	5.5	9.48	13.1	24.65	41.22



Lognormal Background Statistics for Data Sets with Non-Detects

User Selected Options

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\alpha-Pinene\alpha-Pinene data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1
Number of Bootstrap Operations	2000

Pinene

Log-Transformed Statistics

Total Number of Data	50
Number of Non-Detect Data	2
Number of Detected Data	48
Minimum Detected	-0.186
Maximum Detected	3.951
Percent Non-Detects	4.00%
Minimum Non-detect	-0.0943
Maximum Non-detect	1.163
Mean of Detected data	1.368
SD of Detected data	0.993

Lognormal Distribution Test with Detected Values Only

Shapiro Wilk Test Statistic	0.952
5% Shapiro Wilk Critical Value	0.947

Data appear Lognormal at 5% Significance Level

Background Statistics Assuming Lognormal Distribution

DL/2 Substitution Method

Mean (Log Scale)	1.307
SD (Log Scale)	1.026
Tolerance Factor K	0.236
95% UTL 50% Coverage	4.706
95% UPL	20.99
90% Percentile (z)	13.76
95% Percentile (z)	19.98
99% Percentile (z)	40.2

Note: DL/2 is not a recommended method.

Log ROS Method

Mean in Log Scale	1.306
SD in Log Scale	1.028
Mean in Original Scale	6.535
SD in Original Scale	9.112
95% UTL 50% Coverage	4.705
95% BCA UTL with 50% Coverage	4.9
95% Bootstrap (%) UTL with 50% Coverage	4.95
95% UPL (t)	21.06
90% Percentile (z)	13.79



**Bromodichloromethane**

From File: P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\BDCM\BDCM data.wsl

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.95	0.84	0.179	0.423	0.119	3.989	17.75	0.445

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.768	0.77	0.84	0.93	0.946	1	1.72	2.71

**Benzene**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Benzene\Benzene data.wst  
 Full Precision OFF  
 Confidence Coefficient 0.95

Raw

Raw Statistics

Number of Valid Observations: 49  
 Number of Missing Values: 1  
 Number of Distinct Observations: 36  
     Minimum: 0.35  
     Maximum: 20  
 Mean of Raw Data: 2.785  
 Standard Deviation of Raw Data: 4.24  
     Kstar: 0.78  
 Mean of Log Transformed Data: 0.299  
 Standard Deviation of Log Transformed Data: 1.105

Normal Distribution Test Results

Correlation Coefficient R: 0.773  
 Shapiro Wilk Test Statistic: 0.609  
 Shapiro Wilk Critical (0.95) Value: 0.947  
 Approximate Shapiro Wilk P Value: 2.998E-15  
 Lilliefors Test Statistic: 0.317  
 Lilliefors Critical (0.95) Value: 0.127

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R: 0.965  
 A-D Test Statistic: 3.66  
 A-D Critical (0.95) Value: 0.789  
 K-S Test Statistic: 0.212  
 K-S Critical(0.95) Value: 0.131

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R: 0.947  
 Shapiro Wilk Test Statistic: 0.88  
 Shapiro Wilk Critical (0.95) Value: 0.947  
 Approximate Shapiro Wilk P Value: 4.0091E-5  
 Lilliefors Test Statistic: 0.146  
 Lilliefors Critical (0.95) Value: 0.127

Data not Lognormal at (0.05) Significance Level

From File: C:\Documents and Settings\martichbj\Desktop\Data Stats\Benzene\Benzene data.wst

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
Benzene	50	0.14	20	2.732	0.895	17.75	4.213	0.645	2.551	6.457	1.542

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
Benzene	50	0.38	0.46	0.508	0.52	0.895	2.5	3.28	8.36	12.1	17.55

Nonparametric Background Statistics for Data Sets with Non-Detects

User Selected Options

From File C:\Documents and Settings\martichbj\Desktop\Data Stats\Benzene\Benzene data.wst  
 Full Precision OFF  
 Confidence Coefficient 95%  
 Coverage 50%  
 Different or Future K Values 1

Benzene

Total Number of Data 50  
 Number of Non-Detect Data 1  
 Number of Detected Data 49  
 Minimum Detected 0.35  
 Maximum Detected 20  
 Percent Non-Detects 2.00%  
 Minimum Non-detect 0.14  
 Maximum Non-detect 0.14  
 Mean of Detected Data 2.785  
 SD of Detected Data 4.24  
 Mean of Log-Transformed Detected Data 0.299  
 SD of Log-Transformed Detected Data 1.105

Data do not follow a Discernable Distribution (0.05)

Nonparametric Background Statistics

Tolerance Factor K 0.236  
 95% UTL with 50% Coverage  
 Order Statistic 30  
 Achieved CC 0.941  
 UTL 1.3  
 Largest Non-detect at Order 1  
 95% UPL  
 95% UPL 13.9

Kaplan-Meier (KM) Method

Mean 2.736  
 SD 4.168  
 Standard Error of Mean 0.596  
 95% UTL 50% Coverage 3.72  
 95% KM Chebyshev UPL 21.09  
 95% KM UPL (t) 9.794  
 90% KM Percentile (z) 8.078  
 95% KM Percentile (z) 9.593  
 99% KM Percentile (z) 12.43



Summary Statistics for Raw Full Data Sets

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
AG	27	0.46	20	4.513	1.9	26.22	5.12	1.779	1.673	2.233	1.135
DG	23	0.14	1.4	0.642	0.52	0.0967	0.311	0.208	1.274	1.601	0.484

Percentiles for Raw Full Data Sets

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
AG	27	0.546	0.778	0.956	1.1	1.9	5.95	8.02	11.8	14.4	18.7
DG	23	0.353	0.38	0.464	0.475	0.52	0.745	0.79	1.054	1.37	1.4

Nonparametric Background Statistics for Full Data Sets

User Selected Options

From File C:\Documents and Settings\martichbj\Desktop\Data Stats\Benzene\Benzene data.wst  
 Full Precision OFF  
 Confidence Coefficient 95%  
 Coverage 50%  
 Number of Bootstrap Operations 2000

AG

Some Non-Parametric Statistics

Number of Valid Observations	27
Number of Distinct Observations	23
Minimum	0.46
Maximum	20
Second Largest	15
Mean	4.513
Geometric Mean	2.559
First Quartile	1.1
Median	1.9
Third Quartile	5.95
SD	5.12
Variance	26.22
Coefficient of Variation	1.135
Skewness	1.673
Mean of Log-Transformed data	0.94
SD of Log-Transformed data	1.088

Data appear Lognormal at 5% Significance Level

Non-Parametric Background Statistics

90% Percentile	11.8
95% Percentile	14.4
99% Percentile	18.7

95% UTL with 50% Coverage

Order Statistic	17
Achieved CC	0.939
UTL	3.1

95% BCA Bootstrap UTL with 50% Coverage 3.1

95% Percentile Bootstrap UTL with 50% Coverage 4

95% UPL 18

95% Chebyshev UPL 27.24

Upper Limit Based upon IQR 13.23

Nonparametric Background Statistics for Data Sets with Non-Detects

User Selected Options

From File C:\Documents and Settings\martichbj\Desktop\Data Stats\Benzene\Benzene data.wst  
 Full Precision OFF  
 Confidence Coefficient 95%  
 Coverage 50%  
 Different or Future K Values 1

DG

Total Number of Data 23  
 Number of Non-Detect Data 1  
 Number of Detected Data 22  
 Minimum Detected 0.35  
 Maximum Detected 1.4  
 Percent Non-Detects 4.35%  
 Minimum Non-detect 0.14  
 Maximum Non-detect 0.14  
 Mean of Detected Data 0.665  
 SD of Detected Data 0.298  
 Mean of Log-Transformed Detected Data -0.487  
 SD of Log-Transformed Detected Data 0.389

Data appear Lognormal at 5% Significance Level

Nonparametric Background Statistics

Tolerance Factor K N/A  
 95% UTL with 50% Coverage  
 Order Statistic 15  
 Achieved CC 0.953  
 UTL 0.64  
 Largest Non-detect at Order 1

95% UPL  
 95% UPL 1.4

Kaplan-Meier (KM) Method

Mean 0.651  
 SD 0.292  
 Standard Error of Mean 0.0623  
 95% UTL 50% Coverage N/A  
 95% KM Chebyshev UPL 1.951  
 95% KM UPL (t) 1.163  
 90% KM Percentile (z) 1.025  
 95% KM Percentile (z) 1.131  
 99% KM Percentile (z) 1.33

Nonparametric Background Statistics for Data Sets with Non-Detects

User Selected Options

From File C:\Documents and Settings\martichbj\Desktop\Data Stats\Benzene\Benzene data.wst  
 Full Precision OFF  
 Confidence Coefficient 95%  
 Coverage 50%  
 Different or Future K Values 1

DG

Total Number of Data 23  
 Number of Non-Detect Data 1  
 Number of Detected Data 22  
 Minimum Detected 0.35  
 Maximum Detected 1.4  
 Percent Non-Detects 4.35%  
 Minimum Non-detect 0.14  
 Maximum Non-detect 0.14  
 Mean of Detected Data 0.665  
 SD of Detected Data 0.298  
 Mean of Log-Transformed Detected Data -0.487  
 SD of Log-Transformed Detected Data 0.389

Data appear Lognormal at 5% Significance Level

Nonparametric Background Statistics

Tolerance Factor K N/A  
 95% UTL with 50% Coverage  
 Order Statistic 15  
 Achieved CC 0.953  
 UTL 0.64  
 Largest Non-detect at Order 1  
 95% UPL  
 95% UPL 1.4

Kaplan-Meier (KM) Method

Mean 0.651  
 SD 0.292  
 Standard Error of Mean 0.0623  
 95% UTL 50% Coverage N/A  
 95% KM Chebyshev UPL 1.951  
 95% KM UPL (t) 1.163  
 90% KM Percentile (z) 1.025  
 95% KM Percentile (z) 1.131  
 99% KM Percentile (z) 1.33

Wilcoxon-Mann-Whitney Site vs Background Comparison Test for Full Data Sets without NDs

User Selected Options

From File C:\Documents and Settings\martichbj\Desktop\Data Stats\Benzene\Benzene data.wst  
Full Precision OFF  
Confidence Coefficient 95%  
Substantial Difference 0.000  
Selected Null Hypothesis Site or AOC Mean/Median Less Than or Equal to Background Mean/Median (Form 1)  
Alternative Hypothesis Site or AOC Mean/Median Greater Than Background Mean/Median

Area of Concern Data: AG

Background Data: DG

Raw Statistics

	Site	Background
Number of Valid Observations	27	23
Number of Distinct Observations	23	18
Minimum	0.46	0.14
Maximum	20	1.4
Mean	4.513	0.642
Median	1.9	0.52
SD	5.12	0.311
SE of Mean	0.985	0.0648

Wilcoxon-Mann-Whitney (WMW) Test

H0: Mean/Median of Site or AOC  $\leq$  Mean/Median of Background

Site Rank Sum W-Stat 938.5  
WMW Test U-Stat 4.857  
WMW Critical Value (0.050) 1.645  
P-Value 5.9716E-7

Conclusion with Alpha = 0.05

Reject H0, Conclude Site > Background

P-Value < alpha (0.05)

## **Benzyl Chloride**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**Bromoform**



From File: P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Bromoform\Bromoform data.wst

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**Bromomethane**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**C<sub>5</sub> - C<sub>8</sub> Aliphatics**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\C5-C8 Aliphatics\C5-C8 Aliphatics data.
Full Precision	OFF
Confidence Coefficient	0.95

raw

Raw Statistics

Number of Valid Observations	45
Number of Missing Values	5
Number of Distinct Observations	33
Minimum	31
Maximum	500
Mean of Raw Data	131.1
Standard Deviation of Raw Data	119.1
Kstar	1.673
Mean of Log Transformed Data	4.569
Standard Deviation of Log Transformed Data	0.759

Normal Distribution Test Results

Correlation Coefficient R	0.863
Shapiro Wilk Test Statistic	0.741
Shapiro Wilk Critical (0.95) Value	0.945
Approximate Shapiro Wilk P Value	9.436E-10
Lilliefors Test Statistic	0.226
Lilliefors Critical (0.95) Value	0.132

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R	0.964
A-D Test Statistic	1.462
A-D Critical (0.95) Value	0.764
K-S Test Statistic	0.149
K-S Critical(0.95) Value	0.134

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R	0.978
Shapiro Wilk Test Statistic	0.937
Shapiro Wilk Critical (0.95) Value	0.945
Approximate Shapiro Wilk P Value	0.0238
Lilliefors Test Statistic	0.105
Lilliefors Critical (0.95) Value	0.132

Data not Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	31	500	123.3	81	13435	115.9	59.3	1.973	3.43	0.94

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	33	34.9	41	46	81	160	182	224	414	480.4

**Nonparametric Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\C5-C8 Aliphatics\C5-C8 Aliphatics data.
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1

**C5-C8**

Total Number of Data	50
Number of Non-Detect Data	5
Number of Detected Data	45
Minimum Detected	31
Maximum Detected	500
Percent Non-Detects	10.00%
Minimum Non-detect	33
Maximum Non-detect	130
Mean of Detected Data	131.1
SD of Detected Data	119.1
Mean of Log-Transformed Detected Data	4.569
SD of Log-Transformed Detected Data	0.759

**Data do not follow a Discernable Distribution (0.05)**

**Nonparametric Background Statistics**

Tolerance Factor K	0.236
<b>95% UTL with 50% Coverage</b>	
Order Statistic	30
Achieved CC	0.941
UTL	94
<b>Warning: Largest Non-detect at Order</b>	<b>36</b>

**95% UPL**

95% UPL	454.5
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**Kaplan-Meier (KM) Method**

Mean	121.8
SD	115.3
Standard Error of Mean	16.49
95% UTL 50% Coverage	149
95% KM Chebyshev UPL	629.2
95% KM UPL (t)	317
90% KM Percentile (z)	269.5
95% KM Percentile (z)	311.4
99% KM Percentile (z)	389.9

Summary Statistics for Raw Full Data Sets

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
AG	27	34	500	163.9	110	19335	139.1	88.95	1.374	0.722	0.849
DG	23	31	200	75.74	50	2690	51.86	26.69	1.265	0.508	0.685

Percentiles for Raw Full Data Sets

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
AG	27	42.5	46.6	62.6	76	110	200	208	402	457	489.6
DG	23	32.1	33	35	37	50	93.5	116.8	162	179	195.6



Nonparametric Background Statistics for Data Sets with Non-Detects

User Selected Options

From File C:\Documents and Settings\martichbj\Desktop\Data Stats\C5-C8 Aliphatics\C5-C8 Aliphatics data.wst  
 Full Precision OFF  
 Confidence Coefficient 95%  
 Coverage 50%  
 Different or Future K Values 1

AG

Total Number of Data 27  
 Number of Non-Detect Data 1  
 Number of Detected Data 26  
 Minimum Detected 41  
 Maximum Detected 500  
 Percent Non-Detects 3.70%  
 Minimum Non-detect 34  
 Maximum Non-detect 34  
 Mean of Detected Data 168.8  
 SD of Detected Data 139.3  
 Mean of Log-Transformed Detected Data 4.845  
 SD of Log-Transformed Detected Data 0.75

Data appear Lognormal at 5% Significance Level

Nonparametric Background Statistics

Tolerance Factor K N/A  
 95% UTL with 50% Coverage  
 Order Statistic 17  
 Achieved CC 0.939  
 UTL 120  
 Largest Non-detect at Order 1  
 95% UPL  
 95% UPL 484

Kaplan-Meier (KM) Method

Mean 164.1  
 SD 136.2  
 Standard Error of Mean 26.73  
 95% UTL 50% Coverage N/A  
 95% KM Chebyshev UPL 768.7  
 95% KM UPL (t) 400.7  
 90% KM Percentile (z) 338.7  
 95% KM Percentile (z) 388.2  
 99% KM Percentile (z) 481

DG

Total Number of Data	23
Number of Non-Detect Data	4
Number of Detected Data	19
Minimum Detected	31
Maximum Detected	200
Percent Non-Detects	17.39%
Minimum Non-detect	33
Maximum Non-detect	130
Mean of Detected Data	79.42
SD of Detected Data	53.14
Mean of Log-Transformed Detected Data	4.191
SD of Log-Transformed Detected Data	0.604

Data Follow Appr. Gamma Distribution at 5% Significance Level

**Nonparametric Background Statistics**

Tolerance Factor K	N/A
95% UTL with 50% Coverage	
Order Statistic	15
Achieved CC	0.953
UTL	71
Warning: Largest Non-detect at Order	19

95% UPL

95% UPL	196
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**Kaplan-Meier (KM) Method**

Mean	71.98
SD	50.05
Standard Error of Mean	10.77
95% UTL 50% Coverage	N/A
95% KM Chebyshev UPL	294.8
95% KM UPL (t)	159.8
90% KM Percentile (z)	136.1
95% KM Percentile (z)	154.3
99% KM Percentile (z)	188.4

Wilcoxon-Mann-Whitney Site vs Background Comparison Test for Data Sets with Non-Detects

User Selected Options

From File C:\Documents and Settings\martichbj\Desktop\Data Stats\C5-C8 Aliphatics\C5-C8 Aliphatics data.wst  
 Full Precision OFF  
 Confidence Coefficient 95%  
 Substantial Difference (S) 0.000  
 Selected Null Hypothesis Site or AOC Mean/Median Less Than or Equal to Background Mean/Median (Form 1)  
 Alternative Hypothesis Site or AOC Mean/Median Greater Than Background Mean/Median

Area of Concern Data: AG

Background Data: DG

Raw Statistics

	Site	Background
Number of Valid Data	27	23
Number of Non-Detect Data	1	4
Number of Detect Data	26	19
Minimum Non-Detect	34	33
Maximum Non-Detect	34	130
Percent Non detects	3.70%	17.39%
Minimum Detected	41	31
Maximum Detected	500	200
Mean of Detected Data	168.8	79.42
Median of Detected Data	110	66
SD of Detected Data	139.3	53.14

Wilcoxon-Mann-Whitney Site vs Background Test

All observations  $\leq 130$  (Max DL) are ranked the same

Wilcoxon-Mann-Whitney (WMW) Test

H0: Mean/Median of Site or AOC  $\leq$  Mean/Median of Background

Site Rank Sum W-Stat	773.5
WMW Test U-Stat	1.645
WMW Critical Value (0.050)	1.645
P-Value	0.05

Conclusion with Alpha = 0.05

Do Not Reject H0, Conclude Site  $\leq$  Background

P-Value  $\geq$  alpha (0.05)

## **C<sub>9</sub> – C<sub>10</sub> Aromatics**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\C9-C10 Aromatics\C9-C10 Aromatics da
Full Precision	OFF
Confidence Coefficient	0.95

raw

Raw Statistics

Number of Valid Observations	20
Number of Missing Values	30
Number of Distinct Observations	14
Minimum	8.6
Maximum	70
Mean of Raw Data	19.9
Standard Deviation of Raw Data	16.36
Kstar	2.004
Mean of Log Transformed Data	2.76
Standard Deviation of Log Transformed Data	0.647

Normal Distribution Test Results

Correlation Coefficient R	0.847
Shapiro Wilk Test Statistic	0.725
Shapiro Wilk Critical (0.95) Value	0.905
Approximate Shapiro Wilk P Value	2.9039E-5
Lilliefors Test Statistic	0.268
Lilliefors Critical (0.95) Value	0.198

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R	0.958
A-D Test Statistic	1.431
A-D Critical (0.95) Value	0.751
K-S Test Statistic	0.247
K-S Critical(0.95) Value	0.196

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R	0.93
Shapiro Wilk Test Statistic	0.854
Shapiro Wilk Critical (0.95) Value	0.905
Approximate Shapiro Wilk P Value	0.00591
Lilliefors Test Statistic	0.218
Lilliefors Critical (0.95) Value	0.198

Data not Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C9-C10	50	6.7	70	13.84	9.2	142.6	11.94	2.001	3.055	10.46	0.863

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C9-C10	50	7.19	7.67	8.14	8.4	9.2	12	15	23.9	37.85	58.73

**Nonparametric Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\C9-C10 Aromatics\C9-C10 Aromatics da
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1

**C9-C10**

Total Number of Data	50
Number of Non-Detect Data	30
Number of Detected Data	20
Minimum Detected	8.6
Maximum Detected	70
Percent Non-Detects	60.00%
Minimum Non-detect	6.7
Maximum Non-detect	32
Mean of Detected Data	19.9
SD of Detected Data	16.36
Mean of Log-Transformed Detected Data	2.76
SD of Log-Transformed Detected Data	0.647

**Data do not follow a Discernable Distribution (0.05)**

**Nonparametric Background Statistics**

Tolerance Factor K	0.236
<b>95% UTL with 50% Coverage</b>	
Order Statistic	30
Achieved CC	0.941
UTL	9.5
<b>Warning: Largest Non-detect at Order</b>	<b>46</b>

<b>95% UPL</b>	
95% UPL	43.7

**Kaplan-Meier (KM) Method**

Mean	13.18
SD	11.5
Standard Error of Mean	1.67
95% UTL 50% Coverage	15.89
95% KM Chebyshev UPL	63.79
95% KM UPL (t)	32.65
90% KM Percentile (z)	27.91
95% KM Percentile (z)	32.09
99% KM Percentile (z)	39.93

## **C<sub>9</sub> – C<sub>12</sub> Aliphatics**



Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\C9-C12 Aliphatics\C9-C12 Aliphatics.ws  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 42  
Number of Missing Values 8  
Number of Distinct Observations 33  
Minimum 17  
Maximum 850  
Mean of Raw Data 87.86  
Standard Deviation of Raw Data 151.6  
Kstar 1.014  
Mean of Log Transformed Data 3.943  
Standard Deviation of Log Transformed Data 0.867

Normal Distribution Test Results

Correlation Coefficient R 0.644  
Shapiro Wilk Test Statistic 0.436  
Shapiro Wilk Critical (0.95) Value 0.942  
Approximate Shapiro Wilk P Value 2.220E-16  
Lilliefors Test Statistic 0.373  
Lilliefors Critical (0.95) Value 0.137

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.848  
A-D Test Statistic 3.232  
A-D Critical (0.95) Value 0.776  
K-S Test Statistic 0.228  
K-S Critical(0.95) Value 0.14

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.939  
Shapiro Wilk Test Statistic 0.843  
Shapiro Wilk Critical (0.95) Value 0.942  
Approximate Shapiro Wilk P Value 2.7374E-4  
Lilliefors Test Statistic 0.127  
Lilliefors Critical (0.95) Value 0.137

Data not Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C9-C12	50	14	850	76.58	33	19910	141.1	23.72	4.518	21.71	1.843

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C9-C12	50	15.45	16.9	21	22.25	33	72.75	83	100	195.5	722.6

**Nonparametric Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\C9-C12 Aliphatics\C9-C12 Aliphatics.ws
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1

**C9-C12**

Total Number of Data	50
Number of Non-Detect Data	8
Number of Detected Data	42
Minimum Detected	17
Maximum Detected	850
Percent Non-Detects	16.00%
Minimum Non-detect	14
Maximum Non-detect	29
Mean of Detected Data	87.86
SD of Detected Data	151.6
Mean of Log-Transformed Detected Data	3.943
SD of Log-Transformed Detected Data	0.867

**Data do not follow a Discernable Distribution (0.05)**

**Nonparametric Background Statistics**

Tolerance Factor K	0.236
<b>95% UTL with 50% Coverage</b>	
Order Statistic	30
Achieved CC	0.941
UTL	44
Largest Non-detect at Order	20

**95% UPL**

95% UPL	375.5
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**Kaplan-Meier (KM) Method**

Mean	76.59
SD	139.7
Standard Error of Mean	19.99
95% UTL 50% Coverage	109.5
95% KM Chebyshev UPL	691.4
95% KM UPL (t)	313.1
90% KM Percentile (z)	255.6
95% KM Percentile (z)	306.3
99% KM Percentile (z)	401.5

Summary Statistics for Raw Full Data Sets

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
AG	27	14	590	68.96	39	12374	111.2	31.13	4.294	20.06	1.613
DG	23	15	850	85.57	32	29560	171.9	22.24	4.355	19.91	2.009

Percentiles for Raw Full Data Sets

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
AG	27	16.3	17	21	22	39	68	73	94.6	170	488.6
DG	23	15.2	17	21.4	24	32	86.5	97	100	181	704.8

Nonparametric Background Statistics for Data Sets with Non-Detects

User Selected Options

From File C:\Documents and Settings\martichbj\Desktop\Data Stats\C9-C12 Aliphatics\C9-C12 Aliphatics.wst  
 Full Precision OFF  
 Confidence Coefficient 95%  
 Coverage 50%  
 Different or Future K Values 1

AG

Total Number of Data 27  
 Number of Non-Detect Data 4  
 Number of Detected Data 23  
 Minimum Detected 18  
 Maximum Detected 590  
 Percent Non-Detects 14.81%  
 Minimum Non-detect 14  
 Maximum Non-detect 17  
 Mean of Detected Data 78.17  
 SD of Detected Data 118.4  
 Mean of Log-Transformed Detected Data 3.919  
 SD of Log-Transformed Detected Data 0.812

Data do not follow a Discernable Distribution (0.05)

Nonparametric Background Statistics

Tolerance Factor K N/A  
 95% UTL with 50% Coverage  
 Order Statistic 17  
 Achieved CC 0.939  
 UTL 56  
 Largest Non-detect at Order 4  
 95% UPL  
 95% UPL 434

Kaplan-Meier (KM) Method

Mean 69.26  
 SD 109  
 Standard Error of Mean 21.45  
 95% UTL 50% Coverage N/A  
 95% KM Chebyshev UPL 553.2  
 95% KM UPL (t) 258.6  
 90% KM Percentile (z) 209  
 95% KM Percentile (z) 248.6  
 99% KM Percentile (z) 322.9

DG

Total Number of Data	23
Number of Non-Detect Data	4
Number of Detected Data	19
Minimum Detected	17
Maximum Detected	850
Percent Non-Detects	17.39%
Minimum Non-detect	15
Maximum Non-detect	29
Mean of Detected Data	99.58
SD of Detected Data	186.9
Mean of Log-Transformed Detected Data	3.972
SD of Log-Transformed Detected Data	0.952

Data do not follow a Discernable Distribution (0.05)

**Nonparametric Background Statistics**

Tolerance Factor K	N/A
95% UTL with 50% Coverage	
Order Statistic	15
Achieved CC	0.953
UTL	43
Largest Non-detect at Order	10
95% UPL	
95% UPL	718

**Kaplan-Meier (KM) Method**

Mean	85.4
SD	168.2
Standard Error of Mean	36.03
95% UTL 50% Coverage	N/A
95% KM Chebyshev UPL	834.3
95% KM UPL (t)	380.4
90% KM Percentile (z)	301
95% KM Percentile (z)	362.1
99% KM Percentile (z)	476.7

Wilcoxon-Mann-Whitney Site vs Background Comparison Test for Data Sets with Non-Detects

User Selected Options:

From File C:\Documents and Settings\martichbj\Desktop\Data Stats\C9-C12 Aliphatics\C9-C12 Aliphatics.wst  
Full Precision OFF  
Confidence Coefficient 95%  
Substantial Difference (S) 0.000  
Selected Null Hypothesis Site or AOC Mean/Median Less Than or Equal to Background Mean/Median (Form 1)  
Alternative Hypothesis Site or AOC Mean/Median Greater Than Background Mean/Median

Area of Concern Data: AG

Background Data: DG

Raw Statistics

	Site	Background
Number of Valid Data	27	23
Number of Non-Detect Data	4	4
Number of Detect Data	23	19
Minimum Non-Detect	14	15
Maximum Non-Detect	17	29
Percent Non detects	14.81%	17.39%
Minimum Detected	18	17
Maximum Detected	590	850
Mean of Detected Data	78.17	99.58
Median of Detected Data	46	39
SD of Detected Data	118.4	186.9

Wilcoxon-Mann-Whitney Site vs Background Test

All observations  $\leq 29$  (Max DL) are ranked the same

Wilcoxon-Mann-Whitney (WMW) Test

H0: Mean/Median of Site or AOC  $\leq$  Mean/Median of Background

Site Rank Sum W-Stat 683.5  
WMW Test U-Stat -0.107  
WMW Critical Value (0.050) 1.645  
P-Value 0.543

Conclusion with Alpha = 0.05

Do Not Reject H0, Conclude Site  $\leq$  Background

P-Value  $\geq$  alpha (0.05)

## **Carbon Disulfide**



Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	6.7	32	9.168	8.25	14.72	3.837	1.26	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	7.1	7.29	7.6	7.7	8.25	9.275	9.3	9.82	12.75	25.63

## **Carbon Tetrachloride**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.939	0.85	0.152	0.39	0.119	4.541	23.99	0.415

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.715	0.739	0.768	0.773	0.85	0.93	0.946	1	1.5	2.563

**cis-1,2-Dichloroethene**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**cis-1,3-Dichloropropene**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**Trichlorofluoromethane (CFC 11)**



Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\CFC 11\CFC11 data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 45  
Number of Missing Values 5  
Number of Distinct Observations 18  
Minimum 0.84  
Maximum 5.3  
Mean of Raw Data 1.583  
Standard Deviation of Raw Data 0.916  
Kstar 4.549  
Mean of Log Transformed Data 0.353  
Standard Deviation of Log Transformed Data 0.423

Normal Distribution Test Results

Correlation Coefficient R 0.808  
Shapiro Wilk Test Statistic 0.668  
Shapiro Wilk Critical (0.95) Value 0.945  
Approximate Shapiro Wilk P Value 4.770E-12  
Lilliefors Test Statistic 0.288  
Lilliefors Critical (0.95) Value 0.132

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.902  
A-D Test Statistic 4.251  
A-D Critical (0.95) Value 0.753  
K-S Test Statistic 0.272  
K-S Critical(0.95) Value 0.132

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.896  
Shapiro Wilk Test Statistic 0.804  
Shapiro Wilk Critical (0.95) Value 0.945  
Approximate Shapiro Wilk P Value 1.2809E-7  
Lilliefors Test Statistic 0.251  
Lilliefors Critical (0.95) Value 0.132

Data not Lognormal at (0.05) Significance Level

From File: P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\CFC 11\CFC11 data.wst

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
CFC11	50	0.84	5.3	1.596	1.2	0.823	0.907	0.222	2.309	5.632	0.568

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
CFC11	50	0.989	1	1.1	1.1	1.2	1.775	1.92	2.66	3.6	4.614

**Nonparametric Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\CFC 11\CFC11 data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1

CFC11

Total Number of Data	50
Number of Non-Detect Data	5
Number of Detected Data	45
Minimum Detected	0.84
Maximum Detected	5.3
Percent Non-Detects	10.00%
Minimum Non-detect	0.98
Maximum Non-detect	3.2
Mean of Detected Data	1.583
SD of Detected Data	0.916
Mean of Log-Transformed Detected Data	0.353
SD of Log-Transformed Detected Data	0.423

Data do not follow a Discernable Distribution (0.05)

**Nonparametric Background Statistics**

Tolerance Factor K	0.236
<b>95% UTL with 50% Coverage</b>	
Order Statistic	30
Achieved CC	0.941
UTL	1.3
<b>Warning: Largest Non-detect at Order</b>	<b>46</b>
<b>95% UPL</b>	
95% UPL	3.735

**Kaplan-Meier (KM) Method**

Mean	1.531
SD	0.878
Standard Error of Mean	0.126
95% UTL 50% Coverage	1.738
95% KM Chebyshev UPL	5.395
95% KM UPL (t)	3.017
90% KM Percentile (z)	2.656
95% KM Percentile (z)	2.975
99% KM Percentile (z)	3.573

**1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114)**

From File: P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\CFC 114\CFC-114 data.wst

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**Dichlorodifluoromethane (CFC 12)**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\CFC 12\CFC12 data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 49  
Number of Missing Values 1  
Number of Distinct Observations 10  
Minimum 1.4  
Maximum 2.7  
Mean of Raw Data 2.008  
Standard Deviation of Raw Data 0.213  
Kstar 84.12  
Mean of Log Transformed Data 0.692  
Standard Deviation of Log Transformed Data 0.108

Normal Distribution Test Results

Correlation Coefficient R 0.952  
Shapiro Wilk Test Statistic 0.931  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 0.00854  
Lilliefors Test Statistic 0.17  
Lilliefors Critical (0.95) Value 0.127

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.953  
A-D Test Statistic 1.517  
A-D Critical (0.95) Value 0.748  
K-S Test Statistic 0.168  
K-S Critical(0.95) Value 0.126

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.951  
Shapiro Wilk Test Statistic 0.929  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 0.00693  
Lilliefors Test Statistic 0.17  
Lilliefors Critical (0.95) Value 0.127

Data not Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
raw	49	1.4	2.7	2.008	2	0.0453	0.213	0.148	0.186	2.468	0.106

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
raw	49	1.7	1.8	1.8	1.9	2	2.1	2.1	2.2	2.2	2.604



**Nonparametric Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\CFC 12\CFC12 data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1

CFC12

Total Number of Data	50
Number of Non-Detect Data	1
Number of Detected Data	49
Minimum Detected	1.4
Maximum Detected	2.7
Percent Non-Detects	2.00%
Minimum Non-detect	3.2
Maximum Non-detect	3.2
Mean of Detected Data	2.008
SD of Detected Data	0.213
Mean of Log-Transformed Detected Data	0.692
SD of Log-Transformed Detected Data	0.108

**Data do not follow a Discernable Distribution (0.05)**

**Nonparametric Background Statistics**

Tolerance Factor K	0.236
<b>95% UTL with 50% Coverage</b>	
Order Statistic	30
Achieved CC	0.941
UTL	2.1
<b>Warning: Largest Non-detect at Order</b>	<b>50</b>

**95% UPL**

95% UPL	2.59
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**Kaplan-Meier (KM) Method**

Mean	2.008
SD	0.211
Standard Error of Mean	0.0304
95% UTL 50% Coverage	2.058
95% KM Chebyshev UPL	2.936
95% KM UPL (t)	2.365
90% KM Percentile (z)	2.278
95% KM Percentile (z)	2.355
99% KM Percentile (z)	2.498

## **Chlorobenzene**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**Chloroethane**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**Chloroform**

**Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects**

**User Selected Options**

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Chloroform\Chloroform data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

**Raw Statistics**

Number of Valid Observations 18  
Number of Missing Values 32  
Number of Distinct Observations 14  
Minimum 1.1  
Maximum 5.2  
Mean of Raw Data 2.406  
Standard Deviation of Raw Data 1.252  
Kstar 3.885  
Mean of Log Transformed Data 0.766  
Standard Deviation of Log Transformed Data 0.476

**Normal Distribution Test Results**

Correlation Coefficient R 0.926  
Shapiro Wilk Test Statistic 0.85  
Shapiro Wilk Critical (0.95) Value 0.897  
Approximate Shapiro Wilk P Value 0.00785  
Lilliefors Test Statistic 0.208  
Lilliefors Critical (0.95) Value 0.209

**Data not Normal at (0.05) Significance Level**

**Gamma Distribution Test Results**

Correlation Coefficient R 0.971  
A-D Test Statistic 0.498  
A-D Critical (0.95) Value 0.743  
K-S Test Statistic 0.162  
K-S Critical(0.95) Value 0.204

**Data appear Gamma Distributed at (0.05) Significance Level**

**Lognormal Distribution Test Results**

Correlation Coefficient R 0.978  
Shapiro Wilk Test Statistic 0.943  
Shapiro Wilk Critical (0.95) Value 0.897  
Approximate Shapiro Wilk P Value 0.37  
Lilliefors Test Statistic 0.131  
Lilliefors Critical (0.95) Value 0.209

**Data appear Lognormal at (0.05) Significance Level**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
CFM	50	0.67	5.2	1.428	0.93	1.12	1.058	0.237	2.253	5.062	0.741

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
CFM	50	0.715	0.767	0.78	0.815	0.93	1.725	2.02	2.73	3.55	5.151



**Lognormal Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Chloroform\Chloroform data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1
Number of Bootstrap Operations	2000

**CFM**

**Log-Transformed Statistics**

Total Number of Data	50
Number of Non-Detect Data	32
Number of Detected Data	18
Minimum Detected	0.0953
Maximum Detected	1.649
Percent Non-Detects	64.00%
Minimum Non-detect	-0.4
Maximum Non-detect	0.642
Mean of Detected data	0.766
SD of Detected data	0.476

**Lognormal Distribution Test with Detected Values Only**

Shapiro Wilk Test Statistic	0.943
5% Shapiro Wilk Critical Value	0.897

**Data appear Lognormal at 5% Significance Level**

**Background Statistics Assuming Lognormal Distribution**

**DL/2 Substitution Method**

Mean (Log Scale)	-0.263
SD (Log Scale)	0.841
Tolerance Factor K	0.236
95% UTL 50% Coverage	0.937
95% UPL	3.191
90% Percentile (z)	2.257
95% Percentile (z)	3.064
99% Percentile (z)	5.434

**Note: DL/2 is not a recommended method.**

**Log ROS Method**

Mean in Log Scale	-0.154
SD in Log Scale	0.78
Mean in Original Scale	1.204
SD in Original Scale	1.176
95% UTL 50% Coverage	1.03
95% BCA UTL with 50% Coverage	0.77
95% Bootstrap (%) UTL with 50% Coverage	0.77
95% UPL (t)	3.212
90% Percentile (z)	2.33



**Chloromethane**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.923	0.835	0.146	0.383	0.119	4.865	26.87	0.415

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.715	0.739	0.768	0.77	0.835	0.93	0.932	1	1.275	2.563

**Cumene**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**Cyclohexane**

From File: P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Cyclohexane\Cyclohexane data.wst

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	1.3	6.8	2.236	1.8	1.528	1.236	0.297	2.514	6.083	0.553

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	1.4	1.49	1.58	1.6	1.8	2.225	2.44	3.81	5.055	6.555



**Dibromochloromethane**

From File: P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\DBCMDBCM data.wsl

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.928	0.835	0.15	0.387	0.119	4.691	25.39	0.417

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.768	0.77	0.835	0.93	0.932	1	1.41	2.563

**1,2-Dibromo-3-chloropropane**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**d-Limonene**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\d-Limonene\d-Limonene data.wst
Full Precision	OFF
Confidence Coefficient	0.95

C0

Raw Statistics

Number of Valid Observations	50
Number of Distinct Observations	33
Minimum	3.2
Maximum	230
Mean of Raw Data	28.97
Standard Deviation of Raw Data	40.7
Kstar	1.147
Mean of Log Transformed Data	2.897
Standard Deviation of Log Transformed Data	0.891

Normal Distribution Test Results

Correlation Coefficient R	0.72
Shapiro Wilk Test Statistic	0.545
Shapiro Wilk Critical (0.95) Value	0.947
Approximate Shapiro Wilk P Value	0
Lilliefors Test Statistic	0.27
Lilliefors Critical (0.95) Value	0.125

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R	0.9
A-D Test Statistic	2.296
A-D Critical (0.95) Value	0.775
K-S Test Statistic	0.176
K-S Critical(0.95) Value	0.128

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R	0.977
Shapiro Wilk Test Statistic	0.954
Shapiro Wilk Critical (0.95) Value	0.947
Approximate Shapiro Wilk P Value	0.0897
Lilliefors Test Statistic	0.116
Lilliefors Critical (0.95) Value	0.125

Data appear Lognormal at (0.05) Significance Level

From File: P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\d-Limonene\d-Limonene data.wst

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C0	50	3.2	230	28.97	16	1656	40.7	10.38	3.645	14.5	1.405

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C0	50	4.23	7.15	10.8	11	16	27.5	34	46.4	94.8	200.6

**Lognormal Background Statistics for Full Data Sets**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\d-Limonene\d-Limonene data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1
Number of Bootstrap Operations	2000

C0

**Log-Transformed Statistics**

Number of Valid Observations	50
Number of Distinct Observations	33
Minimum	1.163
Maximum	5.438
Second Largest	5.136
Mean	2.897
First Quartile	2.398
Median	2.773
Third Quartile	3.314
SD	0.891

**Lognormal Distribution Test**

Shapiro Wilk Test Statistic	0.954
5% Shapiro Wilk Critical Value	0.947

**Data appear Lognormal at 5% Significance Level**

**Background Statistics Assuming Lognormal Distribution**

90% Percentile (z)	56.8
95% Percentile (z)	78.52
99% Percentile (z)	144.1
95% UPL	81.98
Tolerance Factor K	0.236
95% UTL with 50% Coverage	22.37

**Some Nonparametric Background Statistics**

95% Chebyshev UPL	208.1
95% Bootstrap BCA UTL with 50% Coverage	21
95% Percentile Bootstrap UTL with 50% Coverage	22.5



**Ethanol**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Ethanol\Ethanol data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

C0

Raw Statistics

Number of Valid Observations 50  
Number of Distinct Observations 40  
Minimum 69  
Maximum 5900  
Mean of Raw Data 986.3  
Standard Deviation of Raw Data 1078  
Kstar 1.092  
Mean of Log Transformed Data 6.399  
Standard Deviation of Log Transformed Data 1.046

Normal Distribution Test Results

Correlation Coefficient R 0.855  
Shapiro Wilk Test Statistic 0.75  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 1.096E-10  
Lilliefors Test Statistic 0.197  
Lilliefors Critical (0.95) Value 0.125

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.982  
A-D Test Statistic 0.406  
A-D Critical (0.95) Value 0.776  
K-S Test Statistic 0.0889  
K-S Critical(0.95) Value 0.129

Data appear Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.994  
Shapiro Wilk Test Statistic 0.977  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 0.638  
Lilliefors Test Statistic 0.0737  
Lilliefors Critical (0.95) Value 0.125

Data appear Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C0	50	69	5900	986.3	625	1161081	1078	659.7	2.532	8.44	1.093

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C0	50	120	130	198	287.5	625	1375	1420	2020	2930	4822

**Lognormal Background Statistics for Full Data Sets**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Ethanol\Ethanol data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1
Number of Bootstrap Operations	2000

C0

**Log-Transformed Statistics**

Number of Valid Observations	50
Number of Distinct Observations	40
Minimum	4.234
Maximum	8.683
Second Largest	8.216
Mean	6.399
First Quartile	5.66
Median	6.437
Third Quartile	7.226
SD	1.046

**Lognormal Distribution Test**

Shapiro Wilk Test Statistic	0.977
5% Shapiro Wilk Critical Value	0.947

Data appear Lognormal at 5% Significance Level

**Background Statistics Assuming Lognormal Distribution**

90% Percentile (z)	2296
95% Percentile (z)	3356
99% Percentile (z)	6844
95% UPL	3530
Tolerance Factor K	0.236
95% UTL with 50% Coverage	769.2

**Some Nonparametric Background Statistics**

95% Chebyshev UPL	5730
95% Bootstrap BCA UTL with 50% Coverage	780
95% Percentile Bootstrap UTL with 50% Coverage	860

**Ethyl Acetate**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Ethyl Acetate\Ehtyl acetate data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 45  
Number of Missing Values 5  
Number of Distinct Observations 33  
Minimum 1.6  
Maximum 40  
Mean of Raw Data 11.43  
Standard Deviation of Raw Data 9.902  
Kstar 1.477  
Mean of Log Transformed Data 2.084  
Standard Deviation of Log Transformed Data 0.86

Normal Distribution Test Results

Correlation Coefficient R 0.914  
Shapiro Wilk Test Statistic 0.826  
Shapiro Wilk Critical (0.95) Value 0.945  
Approximate Shapiro Wilk P Value 8.5417E-7  
Lilliefors Test Statistic 0.18  
Lilliefors Critical (0.95) Value 0.132

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.987  
A-D Test Statistic 0.929  
A-D Critical (0.95) Value 0.767  
K-S Test Statistic 0.138  
K-S Critical(0.95) Value 0.134

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.986  
Shapiro Wilk Test Statistic 0.953  
Shapiro Wilk Critical (0.95) Value 0.945  
Approximate Shapiro Wilk P Value 0.11  
Lilliefors Test Statistic 0.12  
Lilliefors Critical (0.95) Value 0.132

Data appear Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
EA	50	1.6	40	10.59	6.3	94.73	9.733	6.153	1.47	1.559	0.919

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
EA	50	1.8	2.27	3.5	3.725	6.3	14.5	16	25	31.3	38.53

**Lognormal Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Ethyl Acetate\Ethyl acetate data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1
Number of Bootstrap Operations	2000

EA

**Log-Transformed Statistics**

Total Number of Data	50
Number of Non-Detect Data	5
Number of Detected Data	45
Minimum Detected	0.47
Maximum Detected	3.689
Percent Non-Detects	10.00%
Minimum Non-detect	0.47
Maximum Non-detect	1.841
Mean of Detected data	2.084
SD of Detected data	0.86

**Lognormal Distribution Test with Detected Values Only**

Shapiro Wilk Test Statistic	0.953
5% Shapiro Wilk Critical Value	0.945

Data appear Lognormal at 5% Significance Level

**Background Statistics Assuming Lognormal Distribution**

**DL/2 Substitution Method**

Mean (Log Scale)	1.905
SD (Log Scale)	0.993
Tolerance Factor K	0.236
95% UTL 50% Coverage	8.492
95% UPL	36.11
90% Percentile (z)	23.99
95% Percentile (z)	34.42
99% Percentile (z)	67.73

Note: DL/2 is not a recommended method.

**Log ROS Method**

Mean in Log Scale	1.927
SD in Log Scale	0.957
Mean in Original Scale	10.47
SD in Original Scale	9.823
95% UTL 50% Coverage	8.606
95% BCA UTL with 50% Coverage	10.25
95% Bootstrap (%) UTL with 50% Coverage	10.25
95% UPL (t)	34.69
90% Percentile (z)	23.4





**Ethylbenzene**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Ethylbenzene\Ethylbenzene data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 47  
Number of Missing Values 3  
Number of Distinct Observations 38  
Minimum 0.18  
Maximum 11  
Mean of Raw Data 1.68  
Standard Deviation of Raw Data 2.085  
Kstar 1.002  
Mean of Log Transformed Data -0.025  
Standard Deviation of Log Transformed Data 1.025

Normal Distribution Test Results

Correlation Coefficient R 0.819  
Shapiro Wilk Test Statistic 0.689  
Shapiro Wilk Critical (0.95) Value 0.946  
Approximate Shapiro Wilk P Value 5.455E-12  
Lilliefors Test Statistic 0.239  
Lilliefors Critical (0.95) Value 0.129

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.972  
A-D Test Statistic 1.366  
A-D Critical (0.95) Value 0.777  
K-S Test Statistic 0.159  
K-S Critical(0.95) Value 0.133

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.988  
Shapiro Wilk Test Statistic 0.962  
Shapiro Wilk Critical (0.95) Value 0.946  
Approximate Shapiro Wilk P Value 0.211  
Lilliefors Test Statistic 0.093  
Lilliefors Critical (0.95) Value 0.129

Data appear Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
EB	50	0.17	11	1.589	0.775	4.212	2.052	0.778	2.682	8.706	1.291

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
EB	50	0.18	0.209	0.326	0.408	0.775	2.125	2.42	3.96	5.965	8.599

**Lognormal Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Ethylbenzene\Ethylbenzene data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1
Number of Bootstrap Operations	2000

EB

**Log-Transformed Statistics**

Total Number of Data	50
Number of Non-Detect Data	4
Number of Detected Data	46
Minimum Detected	-1.715
Maximum Detected	2.398
Percent Non-Detects	8.00%
Minimum Non-detect	-1.772
Maximum Non-detect	-0.151
Mean of Detected data	-0.0222
SD of Detected data	1.036

**Lognormal Distribution Test with Detected Values Only**

Shapiro Wilk Test Statistic	0.961
5% Shapiro Wilk Critical Value	0.945

**Data appear Lognormal at 5% Significance Level**

**Background Statistics Assuming Lognormal Distribution**

**DL/2 Substitution Method**

Mean (Log Scale)	-0.183
SD (Log Scale)	1.152
Tolerance Factor K	0.236
95% UTL 50% Coverage	1.093
95% UPL	5.86
90% Percentile (z)	3.647
95% Percentile (z)	5.542
99% Percentile (z)	12.15

**Note: DL/2 is not a recommended method.**

**Log ROS Method**

Mean in Log Scale	-0.185
SD in Log Scale	1.156
Mean in Original Scale	1.575
SD in Original Scale	2.06
95% UTL 50% Coverage	1.091
95% BCA UTL with 50% Coverage	1
95% Bootstrap (%) UTL with 50% Coverage	1.045
95% UPL (t)	5.887
90% Percentile (z)	3.657



Summary Statistics for Raw Full Data Sets

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
AG	27	0.31	11	2.416	1.4	6.109	2.472	1.364	1.97	4.507	1.023
DG	23	0.17	2.5	0.619	0.48	0.339	0.582	0.356	2.003	4.114	0.94

Percentiles for Raw Full Data Sets

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
AG	27	0.409	0.46	0.766	0.79	1.4	2.9	3.7	5.92	6.1	9.726
DG	23	0.18	0.18	0.204	0.215	0.48	0.665	0.704	1.46	1.59	2.302

Nonparametric Background Statistics for Full Data Sets

User Selected Options

From File C:\Documents and Settings\martichb\Desktop\Data Stats\Ethylbenzene\Ethylbenzene data.wst  
 Full Precision OFF  
 Confidence Coefficient 95%  
 Coverage 50%  
 Number of Bootstrap Operations 2000

AG

Some Non-Parametric Statistics

Number of Valid Observations	27
Number of Distinct Observations	22
Minimum	0.31
Maximum	11
Second Largest	6.1
Mean	2.416
Geometric Mean	1.564
First Quartile	0.79
Median	1.4
Third Quartile	2.9
SD	2.472
Variance	6.109
Coefficient of Variation	1.023
Skewness	1.97
Mean of Log-Transformed data	0.447
SD of Log-Transformed data	0.952

Data appear Gamma Distributed at 5% Significance Level

Non-Parametric Background Statistics

90% Percentile	5.92
95% Percentile	6.1
99% Percentile	9.726

95% UTL with 50% Coverage

Order Statistic	17
Achieved CC	0.939
UTL	2.3

95% BCA Bootstrap UTL with 50% Coverage 2.3

95% Percentile Bootstrap UTL with 50% Coverage 2.4

95% UPL 9.04

95% Chebyshev UPL 13.39

Upper Limit Based upon IQR 6.065



Nonparametric Background Statistics for Data Sets with Non-Detects

User Selected Options

From File C:\Documents and Settings\martichbj\Desktop\Data Stats\Ethylbenzene\Ethylbenzene data.wst  
 Full Precision OFF  
 Confidence Coefficient 95%  
 Coverage 50%  
 Different or Future K Values 1

DG

Total Number of Data 23  
 Number of Non-Detect Data 3  
 Number of Detected Data 20  
 Minimum Detected 0.18  
 Maximum Detected 2.5  
 Percent Non-Detects 13.04%  
 Minimum Non-detect 0.17  
 Maximum Non-detect 0.18  
 Mean of Detected Data 0.686  
 SD of Detected Data 0.597  
 Mean of Log-Transformed Detected Data -0.662  
 SD of Log-Transformed Detected Data 0.748

Data appear Gamma Distributed at 5% Significance Level

Nonparametric Background Statistics

Tolerance Factor K N/A  
 95% UTL with 50% Coverage  
 Order Statistic 15  
 Achieved CC 0.953  
 UTL 0.53  
 Largest Non-detect at Order 3  
 95% UPL  
 95% UPL 2.32

Kaplan-Meier (KM) Method

Mean 0.62  
 SD 0.569  
 Standard Error of Mean 0.122  
 95% UTL 50% Coverage N/A  
 95% KM Chebyshev UPL 3.152  
 95% KM UPL (t) 1.617  
 90% KM Percentile (z) 1.348  
 95% KM Percentile (z) 1.555  
 99% KM Percentile (z) 1.943

Wilcoxon-Mann-Whitney Site vs Background Comparison Test for Full Data Sets without NDs

User Selected Options:

From File C:\Documents and Settings\martichbj\Desktop\Data Stats\Ethylbenzene\Ethylbenzene data.wst  
 Full Precision OFF  
 Confidence Coefficient 95%  
 Substantial Difference 0.000  
 Selected Null Hypothesis Site or AOC Mean/Median Less Than or Equal to Background Mean/Median (Form 1)  
 Alternative Hypothesis Site or AOC Mean/Median Greater Than Background Mean/Median

Area of Concern Data: AG

Background Data: DG

Raw Statistics

	Site	Background
Number of Valid Observations	27	23
Number of Distinct Observations	22	20
Minimum	0.31	0.17
Maximum	11	2.5
Mean	2.416	0.619
Median	1.4	0.48
SD	2.472	0.582
SE of Mean	0.476	0.121

Wilcoxon-Mann-Whitney (WMW) Test

H0: Mean/Median of Site or AOC <= Mean/Median of Background

Site Rank Sum W-Stat	903.5
WMW Test U-Stat	4.175
WMW Critical Value (0.050)	1.645
P-Value	1.4880E-5

Conclusion with Alpha = 0.05

Reject H0, Conclude Site > Background

P-Value < alpha (0.05)

## **Hexachlorobutadiene**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**m,p-Xylenes**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\m&p-Xylenes\m&p-Xylenes data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 33  
Number of Missing Values 17  
Number of Distinct Observations 29  
Minimum 1.5  
Maximum 50  
Mean of Raw Data 8.745  
Standard Deviation of Raw Data 10.4  
Kstar 1.049  
Mean of Log Transformed Data 1.665  
Standard Deviation of Log Transformed Data 0.979

Normal Distribution Test Results

Correlation Coefficient R 0.83  
Shapiro Wilk Test Statistic 0.704  
Shapiro Wilk Critical (0.95) Value 0.931  
Approximate Shapiro Wilk P Value 7.0119E-8  
Lilliefors Test Statistic 0.243  
Lilliefors Critical (0.95) Value 0.154

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.977  
A-D Test Statistic 1.191  
A-D Critical (0.95) Value 0.773  
K-S Test Statistic 0.173  
K-S Critical(0.95) Value 0.157

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.974  
Shapiro Wilk Test Statistic 0.933  
Shapiro Wilk Critical (0.95) Value 0.931  
Approximate Shapiro Wilk P Value 0.0532  
Lilliefors Test Statistic 0.124  
Lilliefors Critical (0.95) Value 0.154

Data appear Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
m&p-x	50	1.5	50	6.454	2.65	81.45	9.025	1.705	3.045	11.06	1.398

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
m&p-x	50	1.5	1.5	1.6	1.7	2.65	6.6	8.92	16.2	24.2	39.71

**Lognormal Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\m&p-Xylenes\m&p-Xylenes data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1
Number of Bootstrap Operations	2000

m&p-x

**Log-Transformed Statistics**

Total Number of Data	50
Number of Non-Detect Data	17
Number of Detected Data	33
Minimum Detected	0.405
Maximum Detected	3.912
Percent Non-Detects	34.00%
Minimum Non-detect	0.405
Maximum Non-detect	1.841
Mean of Detected data	1.665
SD of Detected data	0.979

**Lognormal Distribution Test with Detected Values Only**

Shapiro Wilk Test Statistic	0.933
5% Shapiro Wilk Critical Value	0.931

Data appear Lognormal at 5% Significance Level

**Background Statistics Assuming Lognormal Distribution**

**DL/2 Substitution Method**

Mean (Log Scale)	1.073
SD (Log Scale)	1.167
Tolerance Factor K	0.236
95% UTL 50% Coverage	3.85
95% UPL	21.09
90% Percentile (z)	13.05
95% Percentile (z)	19.94
99% Percentile (z)	44.17

Note: DL/2 is not a recommended method.

**Log ROS Method**

Mean in Log Scale	0.953
SD in Log Scale	1.315
Mean in Original Scale	6.025
SD in Original Scale	9.237
95% UTL 50% Coverage	3.538
95% BCA UTL with 50% Coverage	3.35
95% Bootstrap (%) UTL with 50% Coverage	3.8
95% UPL (t)	24.03
90% Percentile (z)	13.98



95% Percentile (z)	22.55
99% Percentile (z)	55.22
<b>Kaplan Meier (KM) Method</b>	
Mean	6.306
SD	8.989
SE of Mean	1.291
95% UTL 50% Coverage	8.426
95% KM Chebyshev UPL	45.88
95% KM UPL (t)	21.53
95% KM Percentile (z)	21.09

## **Methyl Methacrylate**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	1.3	6.3	1.836	1.7	0.568	0.754	0.297	4.804	26.38	0.411

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	1.4	1.49	1.5	1.5	1.7	1.9	1.9	2	2.495	5.075

## **Methylene Chloride**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Methylene chloride\Methylene chloride c  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 19  
Number of Missing Values 31  
Number of Distinct Observations 19  
Minimum 0.83  
Maximum 170  
Mean of Raw Data 19.43  
Standard Deviation of Raw Data 43.31  
Kstar 0.382  
Mean of Log Transformed Data 1.379  
Standard Deviation of Log Transformed Data 1.656

Normal Distribution Test Results

Correlation Coefficient R 0.69  
Shapiro Wilk Test Statistic 0.497  
Shapiro Wilk Critical (0.95) Value 0.901  
Approximate Shapiro Wilk P Value 3.1139E-8  
Lilliefors Test Statistic 0.364  
Lilliefors Critical (0.95) Value 0.203

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.96  
A-D Test Statistic 2.128  
A-D Critical (0.95) Value 0.82  
K-S Test Statistic 0.301  
K-S Critical(0.95) Value 0.212

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.923  
Shapiro Wilk Test Statistic 0.843  
Shapiro Wilk Critical (0.95) Value 0.901  
Approximate Shapiro Wilk P Value 0.00457  
Lilliefors Test Statistic 0.242  
Lilliefors Critical (0.95) Value 0.203

Data not Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
MC	50	0.67	170	7.948	0.93	771.5	27.78	0.237	5.009	26.35	3.495

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
MC	50	0.71	0.739	0.778	0.783	0.93	1.45	2.12	8.95	28.55	135.7

**Nonparametric Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Methylene chloride\Methylene chloride c
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1

MC

Total Number of Data	50
Number of Non-Detect Data	31
Number of Detected Data	19
Minimum Detected	0.83
Maximum Detected	170
Percent Non-Detects	62.00%
Minimum Non-detect	0.67
Maximum Non-detect	3.2
Mean of Detected Data	19.43
SD of Detected Data	43.31
Mean of Log-Transformed Detected Data	1.379
SD of Log-Transformed Detected Data	1.656

Data do not follow a Discernable Distribution (0.05)

**Nonparametric Background Statistics**

Tolerance Factor K	0.236
<b>95% UTL with 50% Coverage</b>	
Order Statistic	30
Achieved CC	0.941
UTL	0.98
<b>Warning: Largest Non-detect at Order</b>	<b>43</b>

**95% UPL**

95% UPL	65.9
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**Kaplan-Meier (KM) Method**

Mean	7.901
SD	27.51
Standard Error of Mean	3.997
95% UTL 50% Coverage	14.39
95% KM Chebyshev UPL	129
95% KM UPL (t)	54.48
90% KM Percentile (z)	43.15
95% KM Percentile (z)	53.15
99% KM Percentile (z)	71.89

**Methyl tert-Butyl Ether**



From File: P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\MTBE\MTBE data.wst

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**Naphthalene**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Naphthalene\Naphthalene data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 40  
Number of Missing Values 10  
Number of Distinct Observations 33  
Minimum 0.18  
Maximum 130  
Mean of Raw Data 3.725  
Standard Deviation of Raw Data 20.48  
Kstar 0.32  
Mean of Log Transformed Data -0.757  
Standard Deviation of Log Transformed Data 1.074

Normal Distribution Test Results

Correlation Coefficient R 0.371  
Shapiro Wilk Test Statistic 0.172  
Shapiro Wilk Critical (0.95) Value 0.94  
Approximate Shapiro Wilk P Value 0  
Lilliefors Test Statistic 0.51  
Lilliefors Critical (0.95) Value 0.14

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.699  
A-D Test Statistic 10.98  
A-D Critical (0.95) Value 0.851  
K-S Test Statistic 0.419  
K-S Critical(0.95) Value 0.151

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.78  
Shapiro Wilk Test Statistic 0.646  
Shapiro Wilk Critical (0.95) Value 0.94  
Approximate Shapiro Wilk P Value 4.516E-11  
Lilliefors Test Statistic 0.206  
Lilliefors Critical (0.95) Value 0.14

Data not Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
Naph	50	0.14	130	3.017	0.355	335.9	18.33	0.23	7.067	49.96	6.075

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
Naph	50	0.155	0.17	0.18	0.21	0.355	0.51	0.572	0.821	1.32	67.23

Nonparametric Background Statistics for Data Sets with Non-Detects

User Selected Options

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Naphthalene\Naphthalene data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1

Naph

Total Number of Data	50
Number of Non-Detect Data	10
Number of Detected Data	40
Minimum Detected	0.18
Maximum Detected	130
Percent Non-Detects	20.00%
Minimum Non-detect	0.14
Maximum Non-detect	0.38
Mean of Detected Data	3.725
SD of Detected Data	20.48
Mean of Log-Transformed Detected Data	-0.757
SD of Log-Transformed Detected Data	1.074

Data do not follow a Discernable Distribution (0.05)

Nonparametric Background Statistics

Tolerance Factor K	0.236
<b>95% UTL with 50% Coverage</b>	
Order Statistic	30
Achieved CC	0.941
UTL	0.39
Largest Non-detect at Order	29

<b>95% UPL</b>	
95% UPL	1.68

**Kaplan-Meier (KM) Method**

Mean	3.017
SD	18.14
Standard Error of Mean	2.599
95% UTL 50% Coverage	7.297
95% KM Chebyshev UPL	82.89
95% KM UPL (t)	33.74
90% KM Percentile (z)	26.27
95% KM Percentile (z)	32.86
99% KM Percentile (z)	45.23

Summary Statistics for Raw Full Data Sets

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
AG	27	0.14	1.9	0.528	0.39	0.176	0.419	0.237	1.941	3.976	0.793
DG	23	0.15	130	5.939	0.24	731.4	27.04	0.133	4.796	23	4.554

Percentiles for Raw Full Data Sets

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
AG	27	0.17	0.176	0.236	0.265	0.39	0.61	0.722	0.992	1.38	1.796
DG	23	0.151	0.164	0.18	0.18	0.24	0.41	0.432	0.556	0.634	101.5

Nonparametric Background Statistics for Data Sets with Non-Detects

User Selected Options:

From File C:\Documents and Settings\martichbj\Desktop\Data Stats\Naphthalene\Naphthalene data.wst  
 Full Precision OFF  
 Confidence Coefficient 95%  
 Coverage 50%  
 Different or Future K Values 1

AG

Total Number of Data: 27  
 Number of Non-Detect Data: 4  
 Number of Detected Data: 23  
 Minimum Detected: 0.21  
 Maximum Detected: 1.9  
 Percent Non-Detects: 14.81%  
 Minimum Non-detect: 0.14  
 Maximum Non-detect: 0.18  
 Mean of Detected Data: 0.591  
 SD of Detected Data: 0.424  
 Mean of Log-Transformed Detected Data: -0.713  
 SD of Log-Transformed Detected Data: 0.596

Data appear Gamma Distributed at 5% Significance Level

Nonparametric Background Statistics

Tolerance Factor K: N/A  
 95% UTL with 50% Coverage  
 Order Statistic: 17  
 Achieved CC: 0.939  
 UTL: 0.51  
 Largest Non-detect at Order: 4

95% UPL  
 95% UPL: 1.74

Kaplan-Meier (KM) Method

Mean: 0.535  
 SD: 0.406  
 Standard Error of Mean: 0.0798  
 95% UTL 50% Coverage: N/A  
 95% KM Chebyshev UPL: 2.335  
 95% KM UPL (t): 1.239  
 90% KM Percentile (z): 1.055  
 95% KM Percentile (z): 1.202  
 99% KM Percentile (z): 1.478

DG

Total Number of Data	23
Number of Non-Detect Data	6
Number of Detected Data	17
Minimum Detected	0.18
Maximum Detected	130
Percent Non-Detects	26.09%
Minimum Non-detect	0.15
Maximum Non-detect	0.38
Mean of Detected Data	7.964
SD of Detected Data	31.45
Mean of Log-Transformed Detected Data	-0.818
SD of Log-Transformed Detected Data	1.522

Data do not follow a Discernable Distribution (0.05)

**Nonparametric Background Statistics**

Tolerance Factor K	N/A
95% UTL with 50% Coverage	
Order Statistic	15
Achieved CC	0.953
UTL	0.37
Warning: Largest Non-detect at Order	16

**95% UPL**

95% UPL	104.1
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**Kaplan-Meier (KM) Method**

Mean	5.935
SD	26.45
Standard Error of Mean	5.685
95% UTL 50% Coverage	N/A
95% KM Chebyshev UPL	123.7
95% KM UPL (t)	52.33
90% KM Percentile (z)	39.83
95% KM Percentile (z)	49.44
99% KM Percentile (z)	67.47



**Wilcoxon-Mann-Whitney Site vs Background Comparison Test for Data Sets with Non-Detects**

**User Selected Options**

From File C:\Documents and Settings\martichbj\Desktop\Data Stats\Naphthalene\Naphthalene data.wst  
 Full Precision OFF  
 Confidence Coefficient 95%  
 Substantial Difference (S) 0.000  
 Selected Null Hypothesis Site or AOC Mean/Median Less Than or Equal to Background Mean/Median (Form 1)  
 Alternative Hypothesis Site or AOC Mean/Median Greater Than Background Mean/Median

Area of Concern Data: AG

Background Data: DG

**Raw Statistics**

	Site	Background
Number of Valid Data	27	23
Number of Non-Detect Data	4	6
Number of Detect Data	23	17
Minimum Non-Detect	0.14	0.15
Maximum Non-Detect	0.18	0.38
Percent Non detects	14.81%	26.09%
Minimum Detected	0.21	0.18
Maximum Detected	1.9	130
Mean of Detected Data	0.591	7.964
Median of Detected Data	0.5	0.36
SD of Detected Data	0.424	31.45

**Wilcoxon-Mann-Whitney Site vs Background Test**

All observations  $\leq 0.38$  (Max DL) are ranked the same

**Wilcoxon-Mann-Whitney (WMW) Test**

**H0: Mean/Median of Site or AOC  $\leq$  Mean/Median of Background**

Site Rank Sum W-Stat	768
WMW Test U-Stat	1.538
WMW Critical Value (0.050)	1.645
P-Value	0.0621

**Conclusion with Alpha = 0.05**

**Do Not Reject H0, Conclude Site  $\leq$  Background**

**P-Value  $\geq$  alpha (0.05)**

**n-Butyl Acetate**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\In-Butyl Acetate\In-Butyl acetate data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 31  
Number of Missing Values 19  
Number of Distinct Observations 23  
Minimum 0.83  
Maximum 14  
Mean of Raw Data 2.882  
Standard Deviation of Raw Data 2.681  
Kstar 1.724  
Mean of Log Transformed Data 0.77  
Standard Deviation of Log Transformed Data 0.731

Normal Distribution Test Results

Correlation Coefficient R 0.833  
Shapiro Wilk Test Statistic 0.715  
Shapiro Wilk Critical (0.95) Value 0.929  
Approximate Shapiro Wilk P Value 3.1207E-7  
Lilliefors Test Statistic 0.222  
Lilliefors Critical (0.95) Value 0.159

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.952  
A-D Test Statistic 1.088  
A-D Critical (0.95) Value 0.76  
K-S Test Statistic 0.194  
K-S Critical(0.95) Value 0.16

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.969  
Shapiro Wilk Test Statistic 0.931  
Shapiro Wilk Critical (0.95) Value 0.929  
Approximate Shapiro Wilk P Value 0.0552  
Lilliefors Test Statistic 0.175  
Lilliefors Critical (0.95) Value 0.159

Data appear Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
BA	50	0.69	14	2.173	1.3	5.364	2.316	0.712	3.201	13.45	1.066

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
BA	50	0.744	0.779	0.838	0.853	1.3	2.475	3.34	4.37	5.675	10.72

**Lognormal Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\n-Butyl Acetate\n-Butyl acetate data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1
Number of Bootstrap Operations	2000

BA

**Log-Transformed Statistics**

Total Number of Data	50
Number of Non-Detect Data	19
Number of Detected Data	31
Minimum Detected	-0.186
Maximum Detected	2.639
Percent Non-Detects	38.00%
Minimum Non-detect	-0.371
Maximum Non-detect	1.163
Mean of Detected data	0.77
SD of Detected data	0.731

**Lognormal Distribution Test with Detected Values Only**

Shapiro Wilk Test Statistic	0.931
5% Shapiro Wilk Critical Value	0.929

**Data appear Lognormal at 5% Significance Level**

**Background Statistics Assuming Lognormal Distribution**

**DL/2 Substitution Method**

Mean (Log Scale)	0.188
SD (Log Scale)	0.97
Tolerance Factor K	0.236
95% UTL 50% Coverage	1.517
95% UPL	6.239
90% Percentile (z)	4.184
95% Percentile (z)	5.952
99% Percentile (z)	11.53

**Note: DL/2 is not a recommended method.**

**Log ROS Method**

Mean in Log Scale	0.176
SD in Log Scale	0.969
Mean in Original Scale	1.965
SD in Original Scale	2.41
95% UTL 50% Coverage	1.499
95% BCA UTL with 50% Coverage	1.3
95% Bootstrap (%) UTL with 50% Coverage	1.3
95% UPL (t)	6.153
90% Percentile (z)	4.129



**n-Heptane**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\n-Heptane\n-Heptane data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 31  
Number of Missing Values 19  
Number of Distinct Observations 25  
Minimum 0.71  
Maximum 20  
Mean of Raw Data 3.293  
Standard Deviation of Raw Data 3.608  
Kstar 1.469  
Mean of Log Transformed Data 0.848  
Standard Deviation of Log Transformed Data 0.794

Normal Distribution Test Results

Correlation Coefficient R 0.774  
Shapiro Wilk Test Statistic 0.629  
Shapiro Wilk Critical (0.95) Value 0.929  
Approximate Shapiro Wilk P Value 7.3984E-9  
Lilliefors Test Statistic 0.237  
Lilliefors Critical (0.95) Value 0.159

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.908  
A-D Test Statistic 0.802  
A-D Critical (0.95) Value 0.763  
K-S Test Statistic 0.135  
K-S Critical(0.95) Value 0.16

Data follow Appr. Gamma Distribution at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.977  
Shapiro Wilk Test Statistic 0.952  
Shapiro Wilk Critical (0.95) Value 0.929  
Approximate Shapiro Wilk P Value 0.212  
Lilliefors Test Statistic 0.125  
Lilliefors Critical (0.95) Value 0.159

Data appear Lognormal at (0.05) Significance Level



From File: P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\In-Heptane\In-Heptane data.wst

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
Hep	50	0.69	20	2.412	1.1	9.372	3.061	0.511	4.181	22.31	1.269

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
Hep	50	0.739	0.769	0.838	0.87	1.1	3.175	3.32	4.98	5.965	13.92

**Lognormal Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\n-Heptane\n-Heptane data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1
Number of Bootstrap Operations	2000

Hep

**Log-Transformed Statistics**

Total Number of Data	50
Number of Non-Detect Data	19
Number of Detected Data	31
Minimum Detected	-0.342
Maximum Detected	2.996
Percent Non-Detects	38.00%
Minimum Non-detect	-0.371
Maximum Non-detect	1.163
Mean of Detected data	0.848
SD of Detected data	0.794

**Lognormal Distribution Test with Detected Values Only**

Shapiro Wilk Test Statistic	0.952
5% Shapiro Wilk Critical Value	0.929

**Data appear Lognormal at 5% Significance Level**

**Background Statistics Assuming Lognormal Distribution**

**DL/2 Substitution Method**

Mean (Log Scale)	0.225
SD (Log Scale)	1.035
Tolerance Factor K	0.236
95% UTL 50% Coverage	1.599
95% UPL	7.223
90% Percentile (z)	4.718
95% Percentile (z)	6.87
99% Percentile (z)	13.91

**Note: DL/2 is not a recommended method.**

**Log ROS Method**

Mean in Log Scale	0.207
SD in Log Scale	1.046
Mean in Original Scale	2.211
SD in Original Scale	3.15
95% UTL 50% Coverage	1.574
95% BCA UTL with 50% Coverage	1.2
95% Bootstrap (%) UTL with 50% Coverage	1.5
95% UPL (t)	7.225
90% Percentile (z)	4.697



**n-Hexane**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\In-Hexane\In-Hexane data.wst  
 Full Precision OFF  
 Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations	31
Number of Missing Values	19
Number of Distinct Observations	23
Minimum	0.8
Maximum	19
Mean of Raw Data	5.387
Standard Deviation of Raw Data	5.715
Kstar	1.007
Mean of Log Transformed Data	1.16
Standard Deviation of Log Transformed Data	1.032

Normal Distribution Test Results

Correlation Coefficient R	0.879
Shapiro Wilk Test Statistic	0.76
Shapiro Wilk Critical (0.95) Value	0.929
Approximate Shapiro Wilk P Value	2.7710E-6
Lilliefors Test Statistic	0.292
Lilliefors Critical (0.95) Value	0.159

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R	0.965
A-D Test Statistic	1.632
A-D Critical (0.95) Value	0.773
K-S Test Statistic	0.263
K-S Critical(0.95) Value	0.162

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R	0.958
Shapiro Wilk Test Statistic	0.895
Shapiro Wilk Critical (0.95) Value	0.929
Approximate Shapiro Wilk P Value	0.00543
Lilliefors Test Statistic	0.221
Lilliefors Critical (0.95) Value	0.159

Data not Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
HX	50	0.75	19	3.745	1.5	24.61	4.961	0.993	1.998	3.036	1.325

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
HX	50	0.775	0.799	0.878	0.91	1.5	3.2	6.14	11.3	15.55	19

Nonparametric Background Statistics for Data Sets with Non-Detects

User Selected Options:

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\n-Hexane\n-Hexane data.wst  
 Full Precision OFF  
 Confidence Coefficient 95%  
 Coverage 50%  
 Different or Future K Values 1

HX

Total Number of Data 50  
 Number of Non-Detect Data 19  
 Number of Detected Data 31  
 Minimum Detected 0.8  
 Maximum Detected 19  
 Percent Non-Detects 38.00%  
 Minimum Non-detect 0.75  
 Maximum Non-detect 3.2  
 Mean of Detected Data 5.387  
 SD of Detected Data 5.715  
 Mean of Log-Transformed Detected Data 1.16  
 SD of Log-Transformed Detected Data 1.032

Data do not follow a Discernable Distribution (0.05)

Nonparametric Background Statistics

Tolerance Factor K 0.236  
 95% UTL with 50% Coverage  
 Order Statistic 30  
 Achieved CC 0.941  
 UTL 1.9  
 Warning: Largest Non-detect at Order 37

95% UPL  
 95% UPL 17.35

Kaplan-Meier (KM) Method

Mean 3.658  
 SD 4.948  
 Standard Error of Mean 0.711  
 95% UTL 50% Coverage 4.826  
 95% KM Chebyshev UPL 25.44  
 95% KM UPL (t) 12.04  
 90% KM Percentile (z) 9.999  
 95% KM Percentile (z) 11.8  
 99% KM Percentile (z) 15.17

**n-Nonane**



Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\In-Nonane\In-Nonane data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 19  
Number of Missing Values 31  
Number of Distinct Observations 16  
Minimum 0.79  
Maximum 39  
Mean of Raw Data 4.098  
Standard Deviation of Raw Data 8.532  
Kstar 0.82  
Mean of Log Transformed Data 0.786  
Standard Deviation of Log Transformed Data 0.882

Normal Distribution Test Results

Correlation Coefficient R 0.574  
Shapiro Wilk Test Statistic 0.36  
Shapiro Wilk Critical (0.95) Value 0.901  
Approximate Shapiro Wilk P Value 8.313E-10  
Lilliefors Test Statistic 0.41  
Lilliefors Critical (0.95) Value 0.203

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.789  
A-D Test Statistic 2.277  
A-D Critical (0.95) Value 0.772  
K-S Test Statistic 0.261  
K-S Critical(0.95) Value 0.205

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.902  
Shapiro Wilk Test Statistic 0.832  
Shapiro Wilk Critical (0.95) Value 0.901  
Approximate Shapiro Wilk P Value 0.00242  
Lilliefors Test Statistic 0.151  
Lilliefors Critical (0.95) Value 0.203

Data not Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
Non	50	0.67	39	2.156	0.93	29.24	5.408	0.245	6.718	46.52	2.508

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
Non	50	0.71	0.748	0.778	0.795	0.93	1.8	2.1	3.02	3.985	22.29

**Nonparametric Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\n-Nonane\n-Nonane data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1

**Non**

Total Number of Data	50
Number of Non-Detect Data	31
Number of Detected Data	19
Minimum Detected	0.79
Maximum Detected	39
Percent Non-Detects	62.00%
Minimum Non-detect	0.67
Maximum Non-detect	3.2
Mean of Detected Data	4.098
SD of Detected Data	8.532
Mean of Log-Transformed Detected Data	0.786
SD of Log-Transformed Detected Data	0.882

**Data do not follow a Discernable Distribution (0.05)**

**Nonparametric Background Statistics**

Tolerance Factor K	0.236
<b>95% UTL with 50% Coverage</b>	
Order Statistic	30
Achieved CC	0.941
UTL	1
<b>Warning: Largest Non-detect at Order 46</b>	
<b>95% UPL</b>	
95% UPL	4.57

**Kaplan-Meier (KM) Method**

Mean	2.057
SD	5.364
Standard Error of Mean	0.779
95% UTL 50% Coverage	3.322
95% KM Chebyshev UPL	25.67
95% KM UPL (t)	11.14
90% KM Percentile (z)	8.931
95% KM Percentile (z)	10.88
99% KM Percentile (z)	14.54

**n-Octane**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\In-Octane\In-Octane data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 20  
Number of Missing Values 30  
Number of Distinct Observations 13  
Minimum 0.82  
Maximum 12  
Mean of Raw Data 1.865  
Standard Deviation of Raw Data 2.437  
Kstar 1.658  
Mean of Log Transformed Data 0.339  
Standard Deviation of Log Transformed Data 0.605

Normal Distribution Test Results

Correlation Coefficient R 0.607  
Shapiro Wilk Test Statistic 0.398  
Shapiro Wilk Critical (0.95) Value 0.905  
Approximate Shapiro Wilk P Value 1.0374E-9  
Lilliefors Test Statistic 0.379  
Lilliefors Critical (0.95) Value 0.198

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.768  
A-D Test Statistic 2.666  
A-D Critical (0.95) Value 0.753  
K-S Test Statistic 0.267  
K-S Critical(0.95) Value 0.196

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.839  
Shapiro Wilk Test Statistic 0.723  
Shapiro Wilk Critical (0.95) Value 0.905  
Approximate Shapiro Wilk P Value 2.6574E-5  
Lilliefors Test Statistic 0.206  
Lilliefors Critical (0.95) Value 0.198

Data not Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
Oct	50	0.67	12	1.329	0.925	2.637	1.624	0.222	6.06	39.88	1.222

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
Oct	50	0.71	0.748	0.78	0.823	0.925	1.175	1.32	2.03	2.3	7.688

**Nonparametric Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\In-Octane\In-Octane data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1

Oct

Total Number of Data	50
Number of Non-Detect Data	30
Number of Detected Data	20
Minimum Detected	0.82
Maximum Detected	12
Percent Non-Detects	60.00%
Minimum Non-detect	0.67
Maximum Non-detect	3.2
Mean of Detected Data	1.865
SD of Detected Data	2.437
Mean of Log-Transformed Detected Data	0.339
SD of Log-Transformed Detected Data	0.605

Data do not follow a Discernable Distribution (0.05)

**Nonparametric Background Statistics**

Tolerance Factor K	0.236
<b>95% UTL with 50% Coverage</b>	
Order Statistic	30
Achieved CC	0.941
UTL	0.98
<b>Warning: Largest Non-detect at Order</b>	<b>49</b>

**95% UPL**

95% UPL	2.705
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**Kaplan-Meier (KM) Method**

Mean	1.247
SD	1.586
Standard Error of Mean	0.23
95% UTL 50% Coverage	1.621
95% KM Chebyshev UPL	8.229
95% KM UPL (t)	3.933
90% KM Percentile (z)	3.28
95% KM Percentile (z)	3.856
99% KM Percentile (z)	4.937

**n-Propylbenzene**



From File: n-Propylbenzene.wst

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	1.002	0.85	0.24	0.49	0.119	3.184	10.86	0.489

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.768	0.773	0.85	0.93	0.972	1.51	1.81	3.004

**o-Xylene**

**Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects**

**User Selected Options**

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\o-Xylenelo-Xylene data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

**Raw Statistics**

Number of Valid Observations 21  
Number of Missing Values 29  
Number of Distinct Observations 19  
Minimum 0.9  
Maximum 16  
Mean of Raw Data 3.929  
Standard Deviation of Raw Data 3.628  
Kstar 1.543  
Mean of Log Transformed Data 1.059  
Standard Deviation of Log Transformed Data 0.775

**Normal Distribution Test Results**

Correlation Coefficient R 0.862  
Shapiro Wilk Test Statistic 0.755  
Shapiro Wilk Critical (0.95) Value 0.908  
Approximate Shapiro Wilk P Value 6.9809E-5  
Lilliefors Test Statistic 0.25  
Lilliefors Critical (0.95) Value 0.193

**Data not Normal at (0.05) Significance Level**

**Gamma Distribution Test Results**

Correlation Coefficient R 0.971  
A-D Test Statistic 0.662  
A-D Critical (0.95) Value 0.756  
K-S Test Statistic 0.158  
K-S Critical(0.95) Value 0.192

**Data appear Gamma Distributed at (0.05) Significance Level**

**Lognormal Distribution Test Results**

Correlation Coefficient R 0.981  
Shapiro Wilk Test Statistic 0.957  
Shapiro Wilk Critical (0.95) Value 0.908  
Approximate Shapiro Wilk P Value 0.46  
Lilliefors Test Statistic 0.11  
Lilliefors Critical (0.95) Value 0.193

**Data appear Lognormal at (0.05) Significance Level**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
o-X	50	0.67	16	2.194	0.935	7.719	2.778	0.319	3.219	12.42	1.267

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
o-X	50	0.719	0.749	0.78	0.823	0.935	2.375	2.88	5.19	7.61	12.47

**Lognormal Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\o-Xylene\o-Xylene data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1
Number of Bootstrap Operations	2000

o-X

**Log-Transformed Statistics**

Total Number of Data	50
Number of Non-Detect Data	29
Number of Detected Data	21
Minimum Detected	-0.105
Maximum Detected	2.773
Percent Non-Detects	58.00%
Minimum Non-detect	-0.4
Maximum Non-detect	1.163
Mean of Detected data	1.059
SD of Detected data	0.775

**Lognormal Distribution Test with Detected Values Only**

Shapiro Wilk Test Statistic	0.957
5% Shapiro Wilk Critical Value	0.908

**Data appear Lognormal at 5% Significance Level**

**Background Statistics Assuming Lognormal Distribution**

**DL/2 Substitution Method**

Mean (Log Scale)	-0.0282
SD (Log Scale)	1.08
Tolerance Factor K	0.236
95% UTL 50% Coverage	1.254
95% UPL	6.052
90% Percentile (z)	3.88
95% Percentile (z)	5.744
99% Percentile (z)	11.99

**Note: DL/2 is not a recommended method.**

**Log ROS Method**

Mean in Log Scale	-0.17
SD in Log Scale	1.184
Mean in Original Scale	1.858
SD in Original Scale	2.923
95% UTL 50% Coverage	1.116
95% BCA UTL with 50% Coverage	1.05
95% Bootstrap (%) UTL with 50% Coverage	1.2
95% UPL (t)	6.266
90% Percentile (z)	3.848



## **Tetrachloroethene**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\PCE\PCE data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 42  
Number of Missing Values 8  
Number of Distinct Observations 37  
Minimum 0.045  
Maximum 3.9  
Mean of Raw Data 0.611  
Standard Deviation of Raw Data 1.019  
Kstar 0.564  
Mean of Log Transformed Data -1.543  
Standard Deviation of Log Transformed Data 1.344

Normal Distribution Test Results

Correlation Coefficient R 0.769  
Shapiro Wilk Test Statistic 0.585  
Shapiro Wilk Critical (0.95) Value 0.942  
Approximate Shapiro Wilk P Value 4.848E-13  
Lilliefors Test Statistic 0.378  
Lilliefors Critical (0.95) Value 0.137

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.95  
A-D Test Statistic 4.339  
A-D Critical (0.95) Value 0.805  
K-S Test Statistic 0.279  
K-S Critical(0.95) Value 0.143

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.929  
Shapiro Wilk Test Statistic 0.809  
Shapiro Wilk Critical (0.95) Value 0.942  
Approximate Shapiro Wilk P Value 1.0399E-5  
Lilliefors Test Statistic 0.18  
Lilliefors Critical (0.95) Value 0.137

Data not Lognormal at (0.05) Significance Level



Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
PCE	50	0.034	3.9	0.52	0.099	0.914	0.956	0.0793	2.265	4.035	1.837

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
PCE	50	0.0445	0.0459	0.0506	0.061	0.099	0.238	0.348	2.31	2.765	3.557

**Nonparametric Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\18469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\PCE\PCE data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1

**PCE**

Total Number of Data	50
Number of Non-Detect Data	8
Number of Detected Data	42
Minimum Detected	0.045
Maximum Detected	3.9
Percent Non-Detects	16.00%
Minimum Non-detect	0.034
Maximum Non-detect	0.051
Mean of Detected Data	0.611
SD of Detected Data	1.019
Mean of Log-Transformed Detected Data	-1.543
SD of Log-Transformed Detected Data	1.344

**Data do not follow a Discernable Distribution (0.05)**

**Nonparametric Background Statistics**

Tolerance Factor K	0.236
<b>95% UTL with 50% Coverage</b>	
Order Statistic	30
Achieved CC	0.941
UTL	0.14
Largest Non-detect at Order	11

**95% UPL**

95% UPL	3.035
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**Kaplan-Meier (KM) Method**

Mean	0.521
SD	0.946
Standard Error of Mean	0.135
95% UTL 50% Coverage	0.744
95% KM Chebyshev UPL	4.686
95% KM UPL (t)	2.123
90% KM Percentile (z)	1.733
95% KM Percentile (z)	2.077
99% KM Percentile (z)	2.722

**Propene**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Propene\Propene data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 48  
Number of Missing Values 2  
Number of Distinct Observations 41  
Minimum 0.94  
Maximum 850  
Mean of Raw Data 26.16  
Standard Deviation of Raw Data 122.1  
Kstar 0.413  
Mean of Log Transformed Data 1.734  
Standard Deviation of Log Transformed Data 1.214

Normal Distribution Test Results

Correlation Coefficient R 0.399  
Shapiro Wilk Test Statistic 0.198  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 0  
Lilliefors Test Statistic 0.454  
Lilliefors Critical (0.95) Value 0.128

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.7  
A-D Test Statistic 7.167  
A-D Critical (0.95) Value 0.832  
K-S Test Statistic 0.322  
K-S Critical(0.95) Value 0.137

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.937  
Shapiro Wilk Test Statistic 0.895  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 2.3501E-4  
Lilliefors Test Statistic 0.122  
Lilliefors Critical (0.95) Value 0.128

Data not Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
PPN	50	0.75	850	25.14	5.95	14327	119.7	5.486	6.955	48.84	4.761

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
PPN	50	0.945	1.29	1.86	2.175	5.95	8.65	11	18.1	35.3	471.7

**Nonparametric Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Propene\Propene data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1

**PPN**

Total Number of Data	50
Number of Non-Detect Data	2
Number of Detected Data	48
Minimum Detected	0.94
Maximum Detected	850
Percent Non-Detects	4.00%
Minimum Non-detect	0.75
Maximum Non-detect	0.77
Mean of Detected Data	26.16
SD of Detected Data	122.1
Mean of Log-Transformed Detected Data	1.734
SD of Log-Transformed Detected Data	1.214

**Data do not follow a Discernable Distribution (0.05)**

**Nonparametric Background Statistics**

Tolerance Factor K	0.236
<b>95% UTL with 50% Coverage</b>	
Order Statistic	30
Achieved CC	0.941
UTL	6.5
Largest Non-detect at Order	2

<b>95% UPL</b>	
95% UPL	60.95

**Kaplan-Meier (KM) Method**

Mean	25.15
SD	118.5
Standard Error of Mean	16.93
95% UTL 50% Coverage	53.1
95% KM Chebyshev UPL	546.8
95% KM UPL (t)	225.8
90% KM Percentile (z)	177
95% KM Percentile (z)	220
99% KM Percentile (z)	300.8

**Styrene**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Styrene\Styrene data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 17  
Number of Missing Values 33  
Number of Distinct Observations 14  
Minimum 0.85  
Maximum 2.9  
Mean of Raw Data 1.549  
Standard Deviation of Raw Data 0.659  
Kstar 5.19  
Mean of Log Transformed Data 0.355  
Standard Deviation of Log Transformed Data 0.414

Normal Distribution Test Results

Correlation Coefficient R 0.942  
Shapiro Wilk Test Statistic 0.874  
Shapiro Wilk Critical (0.95) Value 0.892  
Approximate Shapiro Wilk P Value 0.0302  
Lilliefors Test Statistic 0.235  
Lilliefors Critical (0.95) Value 0.215

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.965  
A-D Test Statistic 0.752  
A-D Critical (0.95) Value 0.741  
K-S Test Statistic 0.202  
K-S Critical(0.95) Value 0.21

Data appear Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.961  
Shapiro Wilk Test Statistic 0.904  
Shapiro Wilk Critical (0.95) Value 0.892  
Approximate Shapiro Wilk P Value 0.104  
Lilliefors Test Statistic 0.177  
Lilliefors Critical (0.95) Value 0.215

Data appear Lognormal at (0.05) Significance Level



Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
Sty	50	0.67	3.2	1.168	0.925	0.357	0.597	0.185	1.92	2.98	0.511

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
Sty	50	0.735	0.759	0.806	0.823	0.925	1.2	1.34	2.21	2.355	3.053

**Lognormal Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Styrene\Styrene data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1
Number of Bootstrap Operations	2000

Sty

**Log-Transformed Statistics**

Total Number of Data	50
Number of Non-Detect Data	33
Number of Detected Data	17
Minimum Detected	-0.163
Maximum Detected	1.065
Percent Non-Detects	66.00%
Minimum Non-detect	-0.4
Maximum Non-detect	1.163
Mean of Detected data	0.355
SD of Detected data	0.414

**Lognormal Distribution Test with Detected Values Only**

Shapiro Wilk Test Statistic	0.904
5% Shapiro Wilk Critical Value	0.892

**Data appear Lognormal at 5% Significance Level**

**Background Statistics Assuming Lognormal Distribution**

**DL/2 Substitution Method**

Mean (Log Scale)	-0.394
SD (Log Scale)	0.64
Tolerance Factor K	0.236
95% UTL 50% Coverage	0.784
95% UPL	1.993
90% Percentile (z)	1.532
95% Percentile (z)	1.933
99% Percentile (z)	2.99

**Note: DL/2 is not a recommended method.**

**Log ROS Method**

Mean in Log Scale	-0.398
SD in Log Scale	0.616
Mean in Original Scale	0.833
SD in Original Scale	0.646
95% UTL 50% Coverage	0.777
95% BCA UTL with 50% Coverage	0.602
95% Bootstrap (%) UTL with 50% Coverage	0.641
95% UPL (t)	1.905
90% Percentile (z)	1.479



## **Trichloroethene**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\TCENTCE data.wst
Full Precision	OFF
Confidence Coefficient	0.95

raw

Raw Statistics

Number of Valid Observations	22
Number of Missing Values	28
Number of Distinct Observations	20
Minimum	0.042
Maximum	5.2
Mean of Raw Data	0.61
Standard Deviation of Raw Data	1.16
Kstar	0.525
Mean of Log Transformed Data	-1.58
Standard Deviation of Log Transformed Data	1.415

Normal Distribution Test Results

Correlation Coefficient R	0.717
Shapiro Wilk Test Statistic	0.536
Shapiro Wilk Critical (0.95) Value	0.911
Approximate Shapiro Wilk P Value	3.4514E-8
Lilliefors Test Statistic	0.336
Lilliefors Critical (0.95) Value	0.189

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R	0.951
A-D Test Statistic	1.555
A-D Critical (0.95) Value	0.798
K-S Test Statistic	0.243
K-S Critical(0.95) Value	0.195

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R	0.953
Shapiro Wilk Test Statistic	0.9
Shapiro Wilk Critical (0.95) Value	0.911
Approximate Shapiro Wilk P Value	0.0272
Lilliefors Test Statistic	0.203
Lilliefors Critical (0.95) Value	0.189

Data not Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
TCE	50	0.034	5.2	0.294	0.048	0.656	0.81	0.0148	5.026	28.51	2.754

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
TCE	50	0.0355	0.0379	0.041	0.042	0.048	0.0955	0.224	0.582	1.3	3.681

**Nonparametric Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\TCEITCE data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1

TCE

Total Number of Data	50
Number of Non-Detect Data	28
Number of Detected Data	22
Minimum Detected	0.042
Maximum Detected	5.2
Percent Non-Detects	56.00%
Minimum Non-detect	0.034
Maximum Non-detect	0.096
Mean of Detected Data	0.61
SD of Detected Data	1.16
Mean of Log-Transformed Detected Data	-1.58
SD of Log-Transformed Detected Data	1.415

Data do not follow a Discernable Distribution (0.05)

**Nonparametric Background Statistics**

Tolerance Factor K	0.236
<b>95% UTL with 50% Coverage</b>	
Order Statistic	30
Achieved CC	0.941
UTL	0.057
<b>Warning: Largest Non-detect at Order</b>	<b>38</b>

**95% UPL**

95% UPL	1.66
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**Kaplan-Meier (KM) Method**

Mean	0.292
SD	0.803
Standard Error of Mean	0.116
95% UTL 50% Coverage	0.481
95% KM Chebyshev UPL	3.826
95% KM UPL (t)	1.651
90% KM Percentile (z)	1.321
95% KM Percentile (z)	1.612
99% KM Percentile (z)	2.159

**trans-1,2-Dichloroethene**



Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

**trans-1,3-Dichloropropene**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.67	3.2	0.917	0.825	0.147	0.384	0.126	4.87	26.88	0.418

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.71	0.729	0.76	0.77	0.825	0.928	0.93	0.982	1.275	2.563

## **Tetrahydrofuran**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\THF\THF data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 15  
Number of Missing Values 35  
Number of Distinct Observations 15  
Minimum 0.8  
Maximum 4.8  
Mean of Raw Data 2.63  
Standard Deviation of Raw Data 1.392  
Kstar 2.668  
Mean of Log Transformed Data 0.807  
Standard Deviation of Log Transformed Data 0.622

Normal Distribution Test Results

Correlation Coefficient R 0.975  
Shapiro Wilk Test Statistic 0.928  
Shapiro Wilk Critical (0.95) Value 0.881  
Approximate Shapiro Wilk P Value 0.348  
Lilliefors Test Statistic 0.111  
Lilliefors Critical (0.95) Value 0.229

Data appear Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.959  
A-D Test Statistic 0.387  
A-D Critical (0.95) Value 0.743  
K-S Test Statistic 0.138  
K-S Critical(0.95) Value 0.223

Data appear Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.965  
Shapiro Wilk Test Statistic 0.909  
Shapiro Wilk Critical (0.95) Value 0.881  
Approximate Shapiro Wilk P Value 0.184  
Lilliefors Test Statistic 0.17  
Lilliefors Critical (0.95) Value 0.229

Data appear Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
THF	50	0.67	4.8	1.447	0.9	1.299	1.14	0.17	1.801	2.122	0.788

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
THF	50	0.74	0.76	0.78	0.803	0.9	1.55	2.02	3.24	4.185	4.702

**Lognormal Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\THF\THF data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1
Number of Bootstrap Operations	2000

THF

**Log-Transformed Statistics**

Total Number of Data	50
Number of Non-Detect Data	35
Number of Detected Data	15
Minimum Detected	-0.223
Maximum Detected	1.569
Percent Non-Detects	70.00%
Minimum Non-detect	-0.4
Maximum Non-detect	1.163
Mean of Detected data	0.807
SD of Detected data	0.622

**Lognormal Distribution Test with Detected Values Only**

Shapiro Wilk Test Statistic	0.909
5% Shapiro Wilk Critical Value	0.881

Data appear Lognormal at 5% Significance Level

**Background Statistics Assuming Lognormal Distribution**

**DL/2 Substitution Method**

Mean (Log Scale)	-0.323
SD (Log Scale)	0.85
Tolerance Factor K	0.236
95% UTL 50% Coverage	0.884
95% UPL	3.054
90% Percentile (z)	2.152
95% Percentile (z)	2.931
99% Percentile (z)	5.233

Note: DL/2 is not a recommended method.

**Log ROS Method**

Mean in Log Scale	-0.442
SD in Log Scale	0.938
Mean in Original Scale	1.069
SD in Original Scale	1.278
95% UTL 50% Coverage	0.802
95% BCA UTL with 50% Coverage	0.536
95% Bootstrap (%) UTL with 50% Coverage	0.536
95% UPL (t)	3.148
90% Percentile (z)	2.139





**Toluene**

Goodness-of-Fit Test Statistics for Full Data Sets without Non-Detects

User Selected Options

From File P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Toluene\Toluene data.wst  
Full Precision OFF  
Confidence Coefficient 0.95

raw

Raw Statistics

Number of Valid Observations 49  
Number of Missing Values 1  
Number of Distinct Observations 41  
Minimum 1.5  
Maximum 130  
Mean of Raw Data 18.69  
Standard Deviation of Raw Data 24.52  
Kstar 0.938  
Mean of Log Transformed Data 2.34  
Standard Deviation of Log Transformed Data 1.06

Normal Distribution Test Results

Correlation Coefficient R 0.806  
Shapiro Wilk Test Statistic 0.669  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 2.730E-13  
Lilliefors Test Statistic 0.249  
Lilliefors Critical (0.95) Value 0.127

Data not Normal at (0.05) Significance Level

Gamma Distribution Test Results

Correlation Coefficient R 0.973  
A-D Test Statistic 1.703  
A-D Critical (0.95) Value 0.78  
K-S Test Statistic 0.181  
K-S Critical(0.95) Value 0.13

Data not Gamma Distributed at (0.05) Significance Level

Lognormal Distribution Test Results

Correlation Coefficient R 0.987  
Shapiro Wilk Test Statistic 0.963  
Shapiro Wilk Critical (0.95) Value 0.947  
Approximate Shapiro Wilk P Value 0.215  
Lilliefors Test Statistic 0.114  
Lilliefors Critical (0.95) Value 0.127

Data appear Lognormal at (0.05) Significance Level

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
Tol	50	0.87	130	18.33	8.25	595.5	24.4	7.339	2.767	9.141	1.331

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
Tol	50	1.88	3.14	4.48	5	8.25	24.5	29.2	49.2	59.85	109.9

**Lognormal Background Statistics for Data Sets with Non-Detects**

**User Selected Options**

From File	P:\8469 (DEQ)\Indoor Air Study\Technical Evaluation\Data Stats\Toluene\Toluene data.wst
Full Precision	OFF
Confidence Coefficient	95%
Coverage	50%
Different or Future K Values	1
Number of Bootstrap Operations	2000

Tol

**Log-Transformed Statistics**

Total Number of Data	50
Number of Non-Detect Data	1
Number of Detected Data	49
Minimum Detected	0.405
Maximum Detected	4.868
Percent Non-Detects	2.00%
Minimum Non-detect	-0.139
Maximum Non-detect	-0.139
Mean of Detected data	2.34
SD of Detected data	1.06

**Lognormal Distribution Test with Detected Values Only**

Shapiro Wilk Test Statistic	0.963
5% Shapiro Wilk Critical Value	0.947

**Data appear Lognormal at 5% Significance Level**

**Background Statistics Assuming Lognormal Distribution**

**DL/2 Substitution Method**

Mean (Log Scale)	2.277
SD (Log Scale)	1.141
Tolerance Factor K	0.236
95% UTL 50% Coverage	12.76
95% UPL	67.25
90% Percentile (z)	42.05
95% Percentile (z)	63.64
99% Percentile (z)	138.5

**Note: DL/2 is not a recommended method.**

**Log ROS Method**

Mean in Log Scale	2.285
SD in Log Scale	1.119
Mean in Original Scale	18.33
SD in Original Scale	24.41
95% UTL 50% Coverage	12.8
95% BCA UTL with 50% Coverage	10.4
95% Bootstrap (%) UTL with 50% Coverage	11.5
95% UPL (t)	65.34
90% Percentile (z)	41.23



## **Vinyl Chloride**

Summary Statistics for Raw Full Dataset

Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	0.034	0.096	0.0442	0.042	1.1604E-4	0.0108	0.00593	3.145	11.95	0.244

Percentiles for Raw Full Dataset

Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	0.0355	0.0369	0.038	0.038	0.042	0.046	0.047	0.049	0.0636	0.0852

## **Vinyl Acetate**



Summary Statistics for Raw Full Dataset

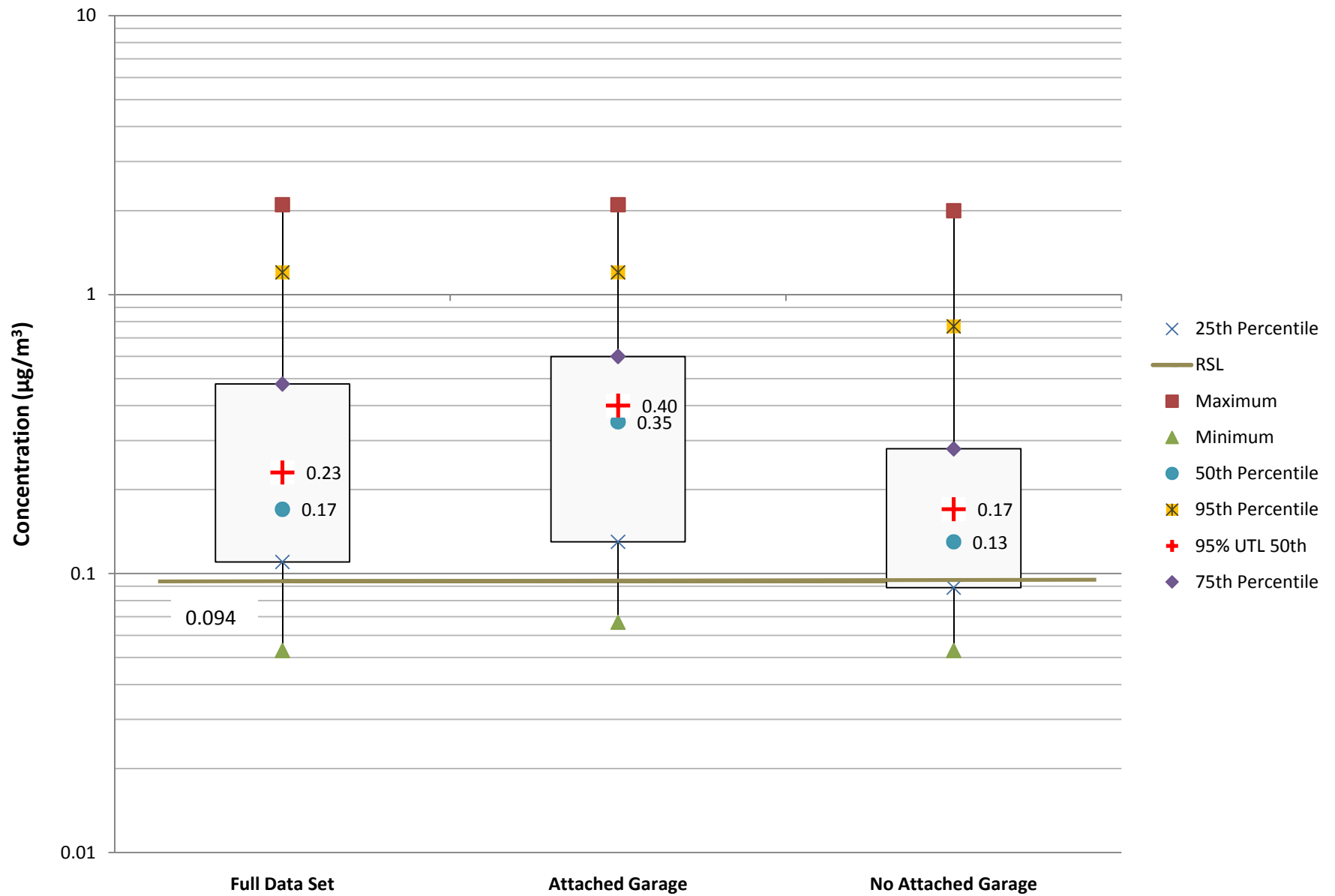
Variable	NumObs	Minimum	Maximum	Mean	Median	Variance	SD	MAD/0.675	Skewness	Kurtosis	CV
C1	50	6.7	32	9.3	8.35	15.12	3.888	1.186	4.617	24.69	0.418

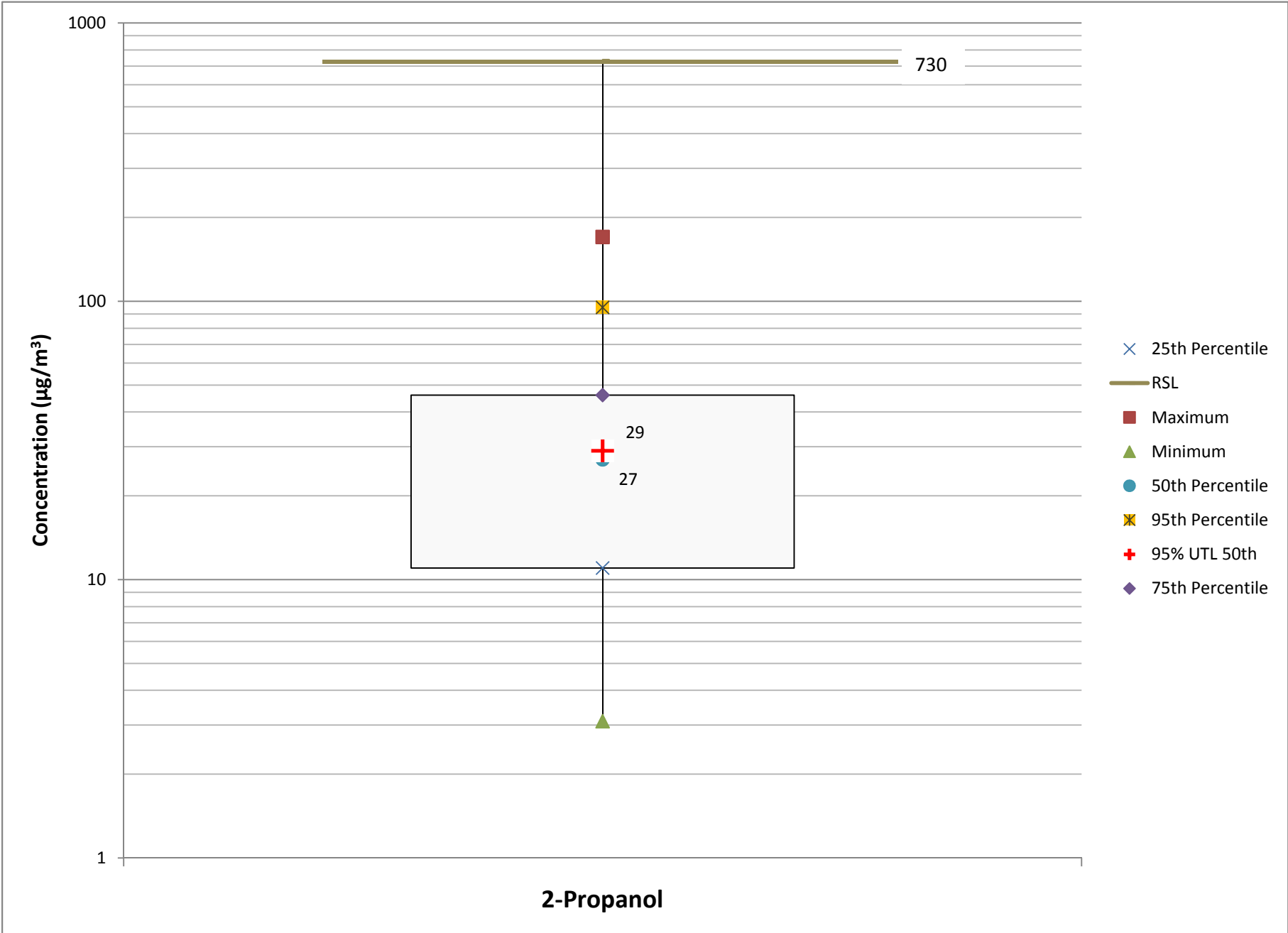
Percentiles for Raw Full Dataset

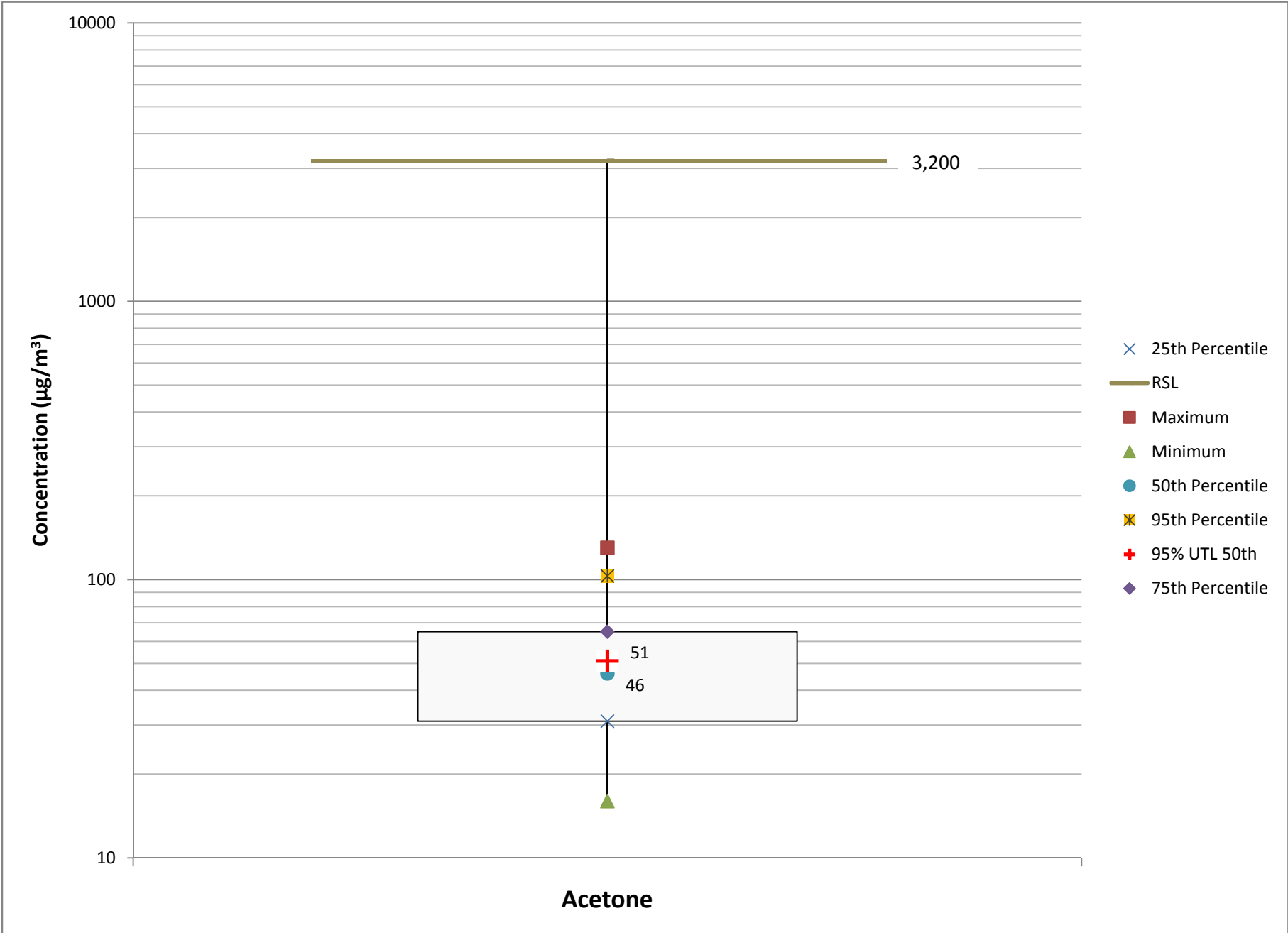
Variable	NumObs	5%ile	10%ile	20%ile	25%ile(Q1)	50%ile(Q2)	75%ile(Q3)	80%ile	90%ile	95%ile	99%ile
C1	50	7.1	7.29	7.68	7.7	8.35	9.3	9.32	10	14.55	25.63

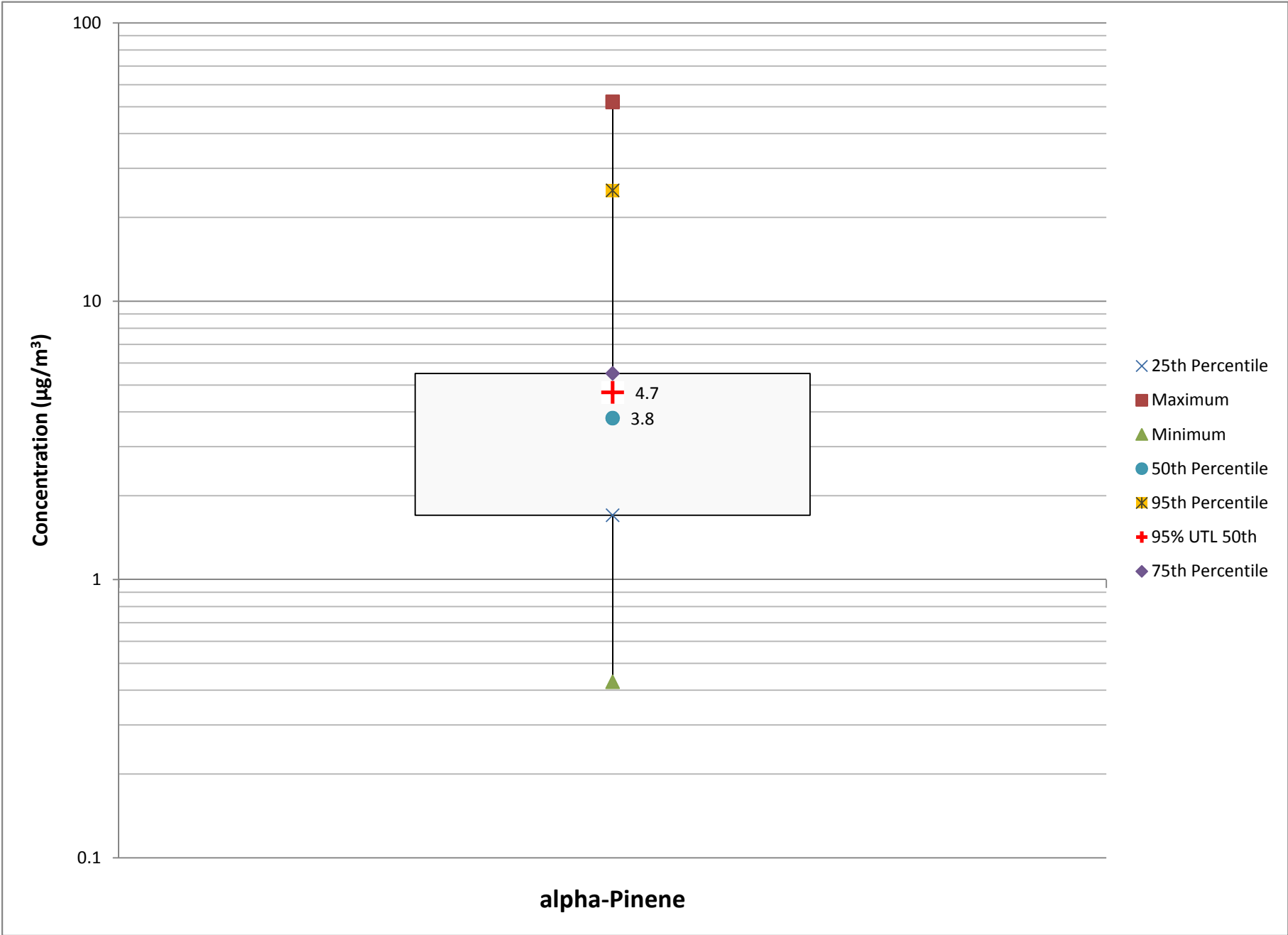
# Appendix D

# 1,2-Dichloroethane

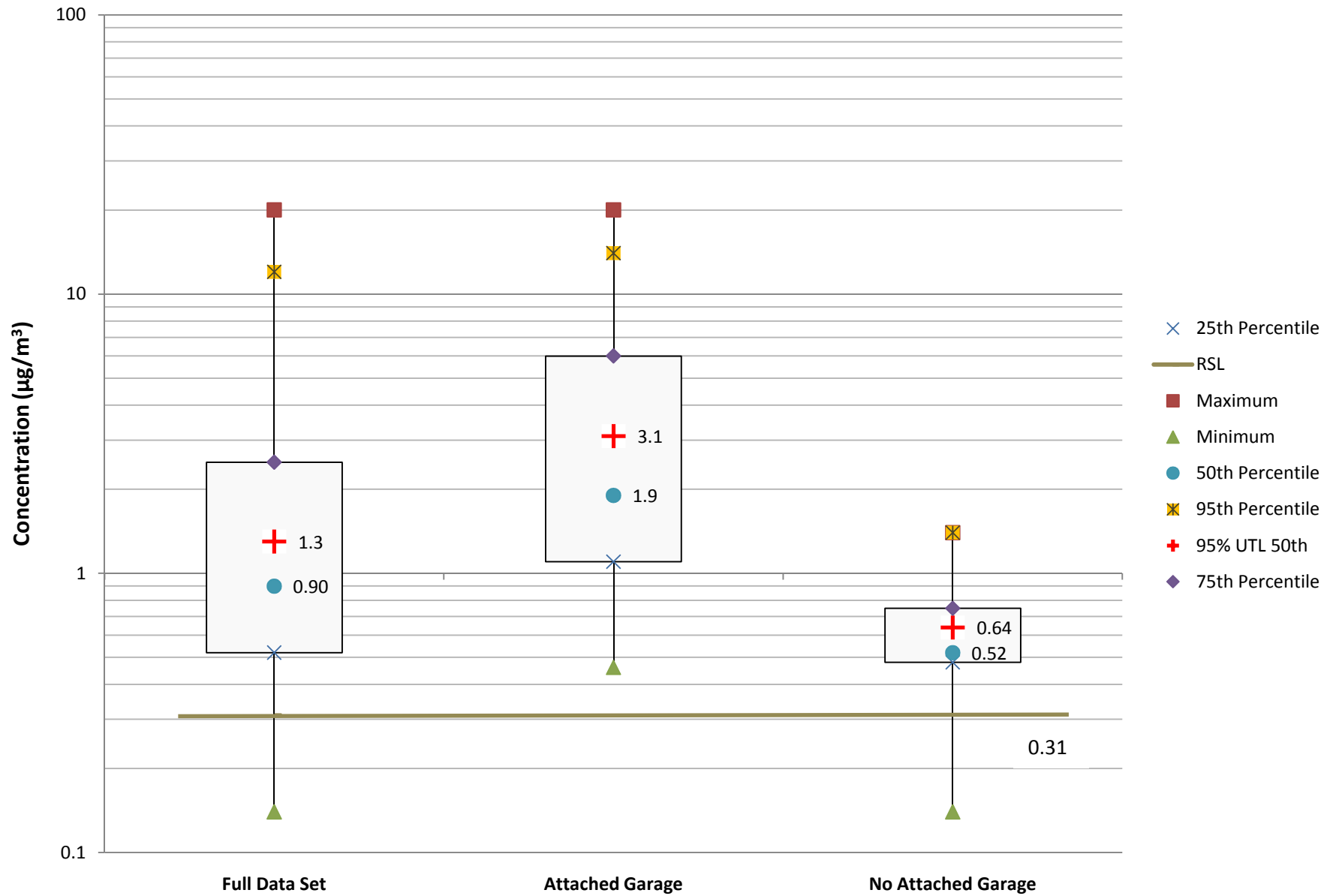




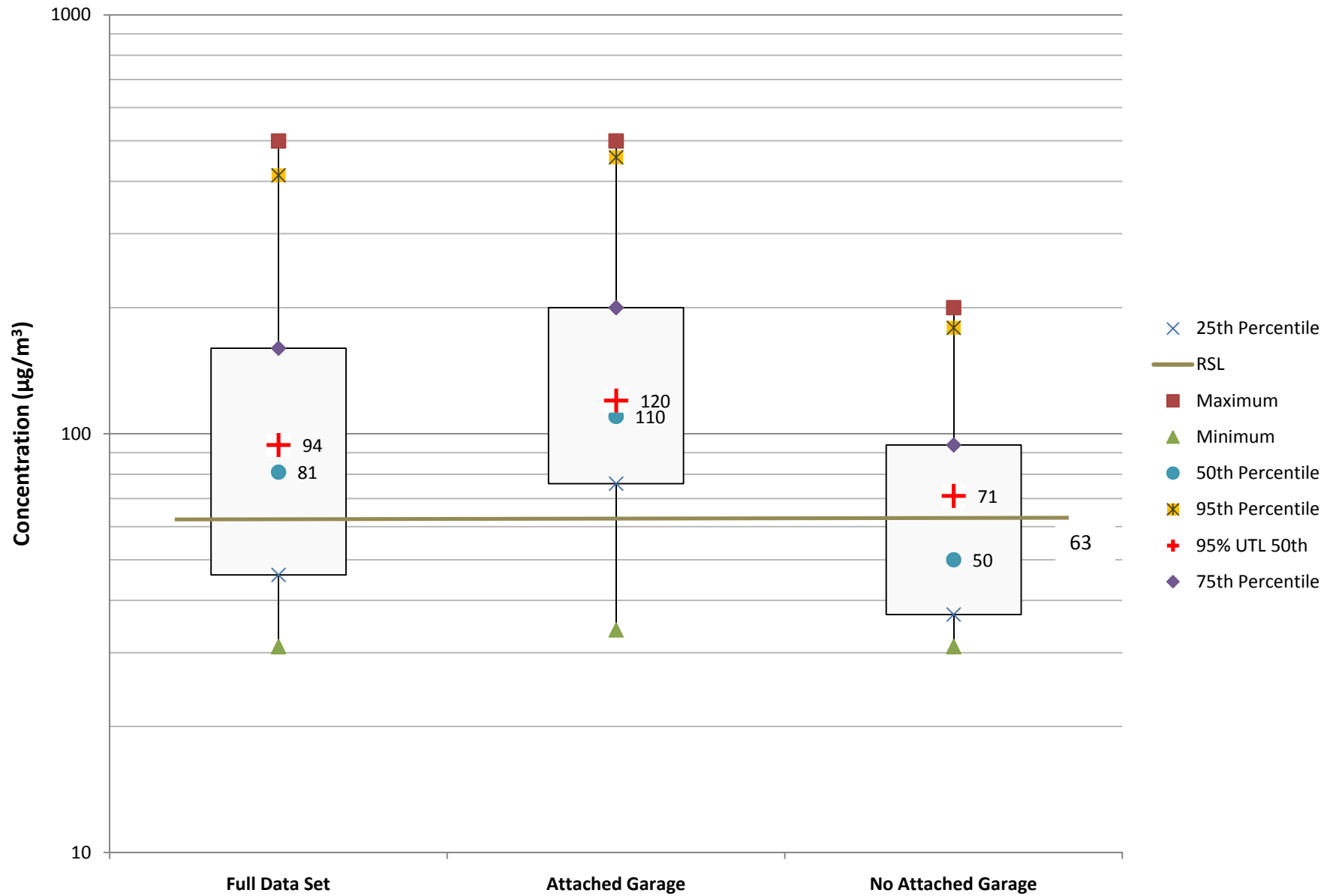




# Benzene

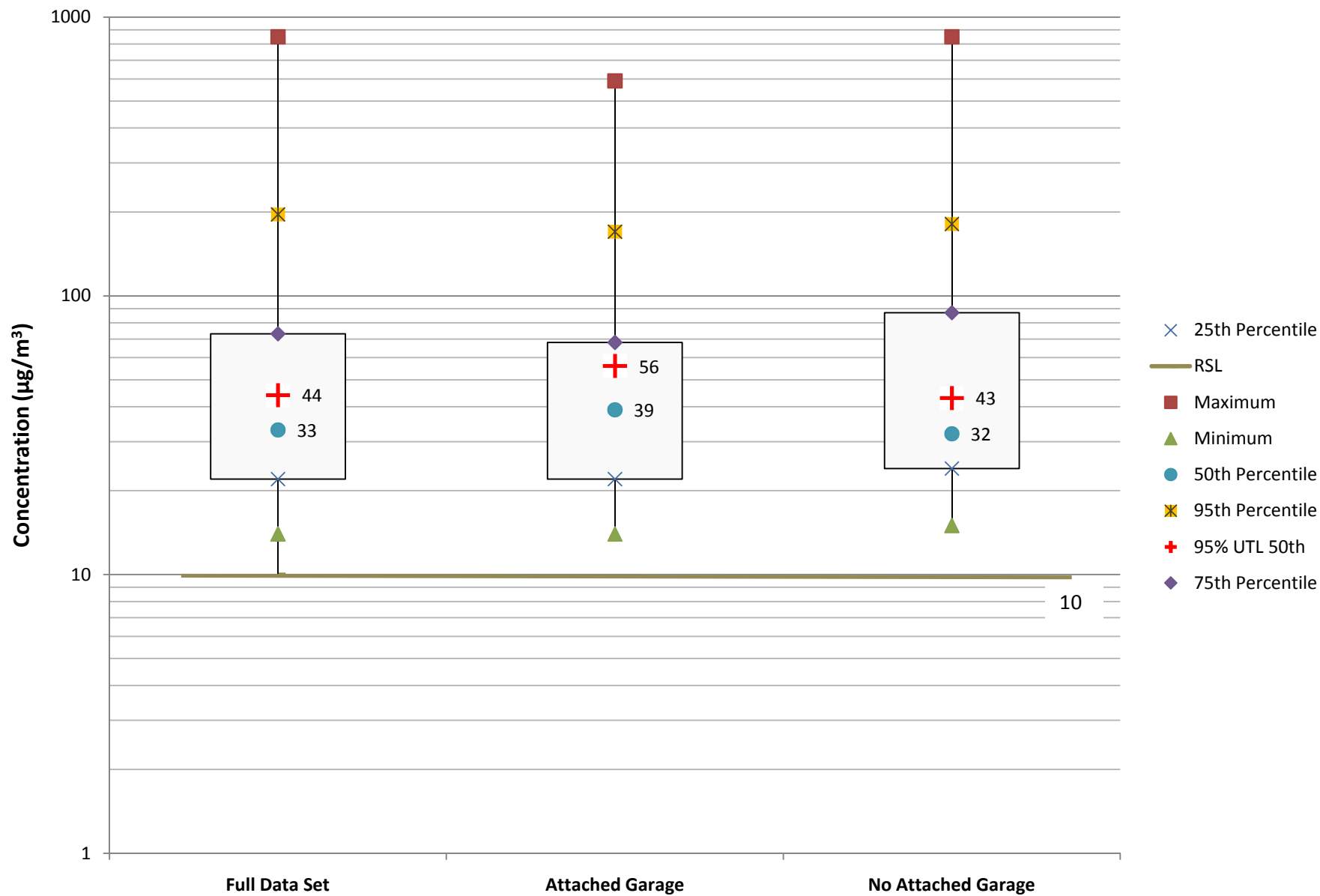


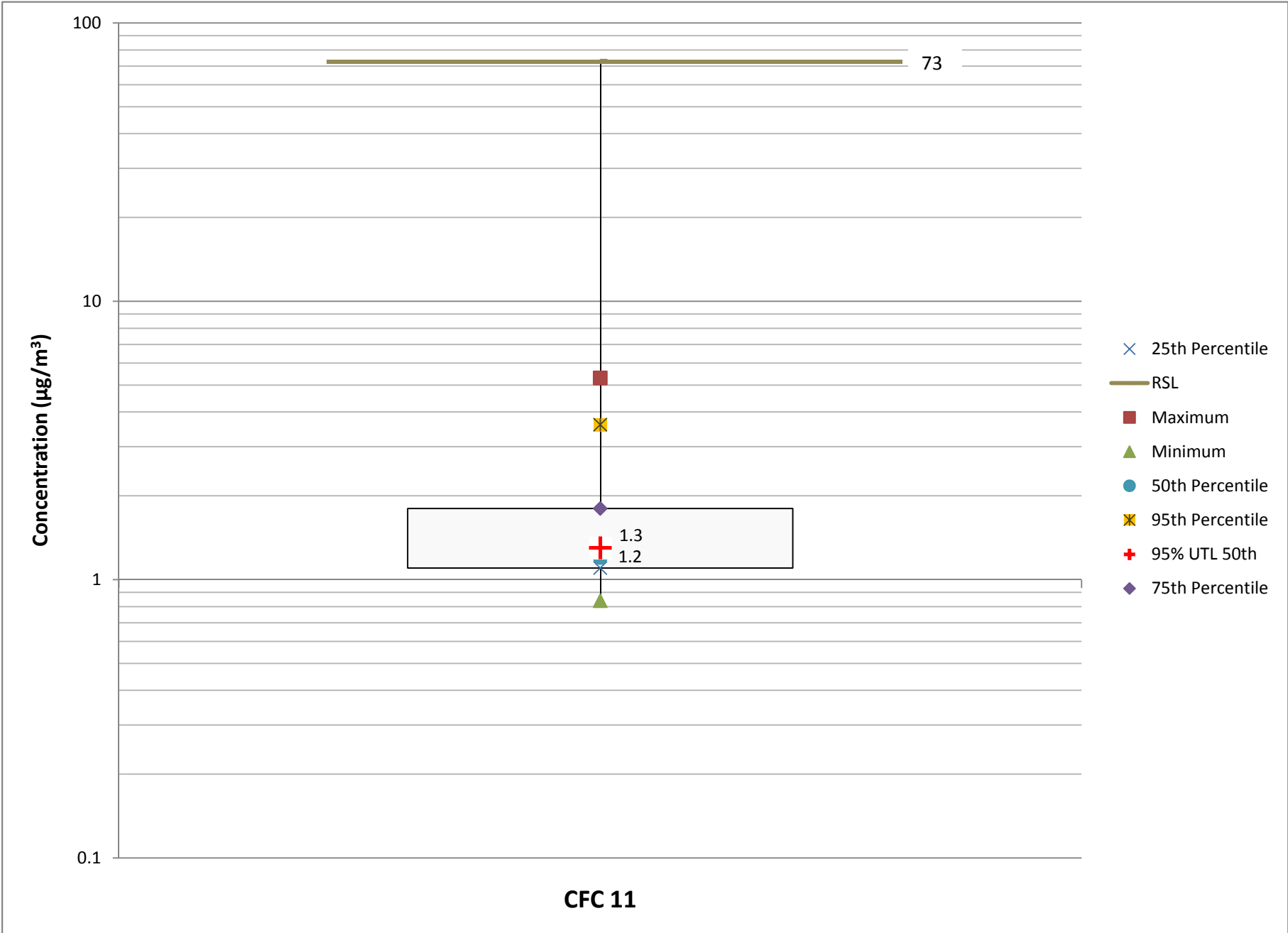
# C5-C8 Aliphatics

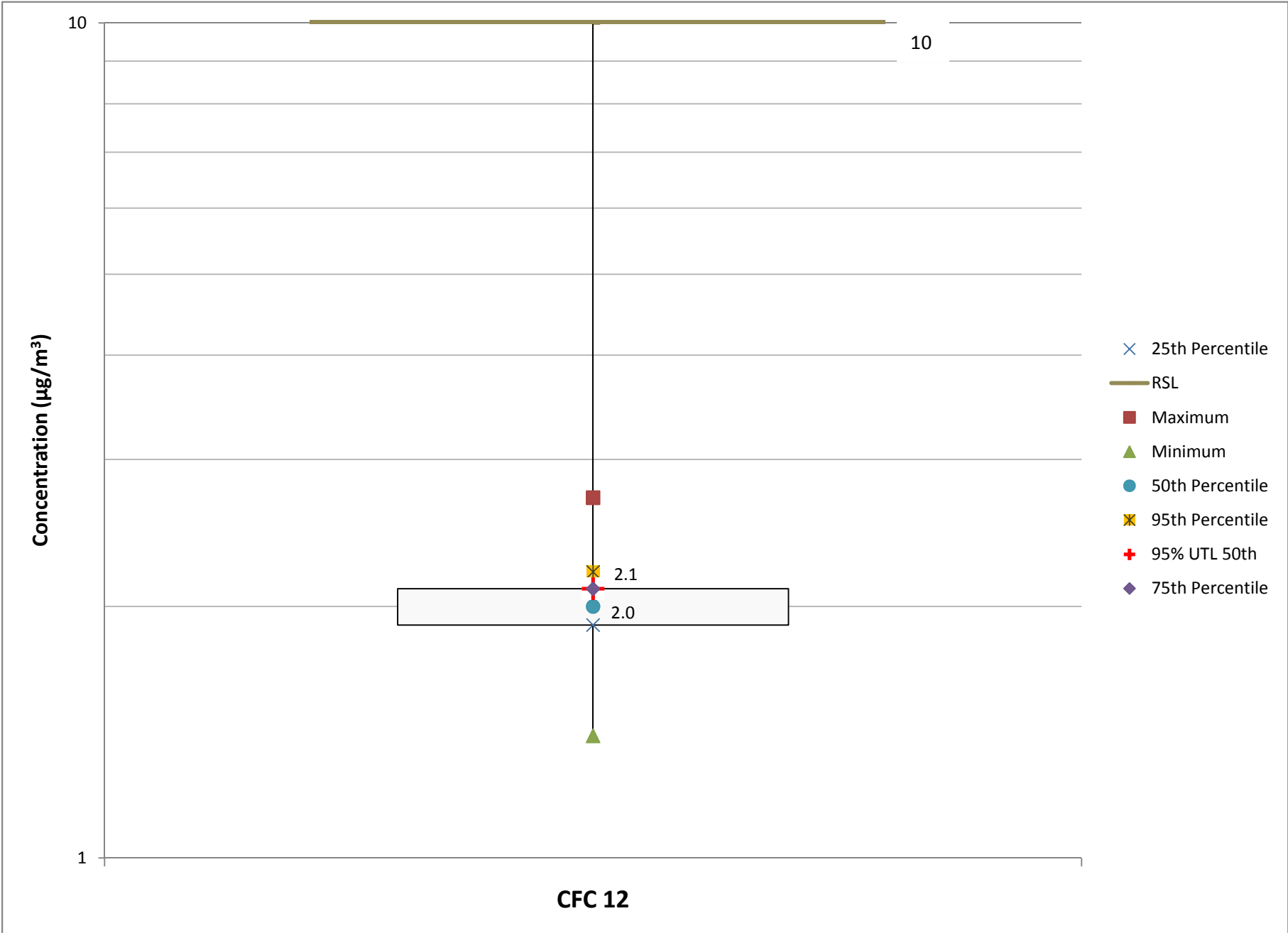


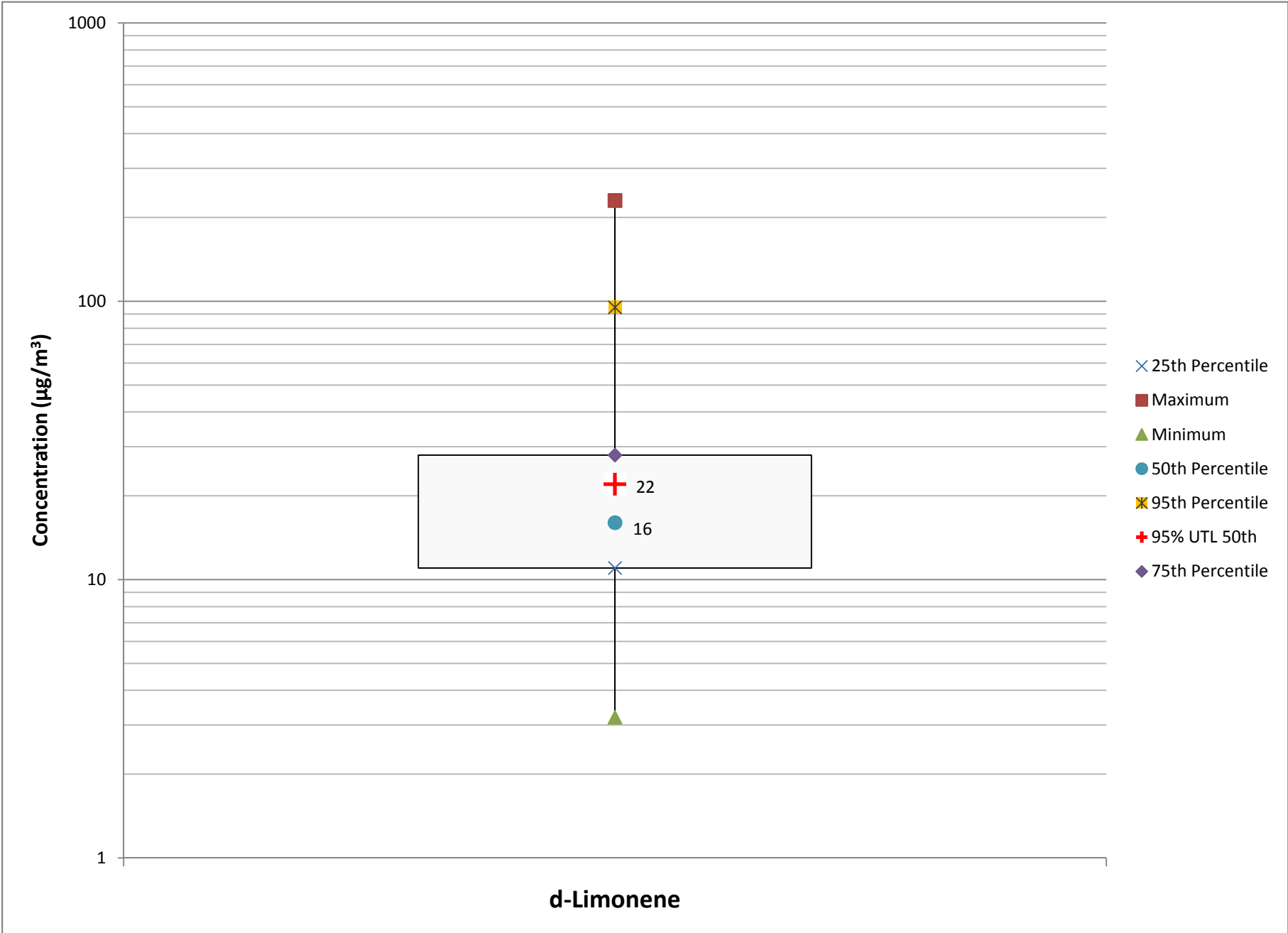


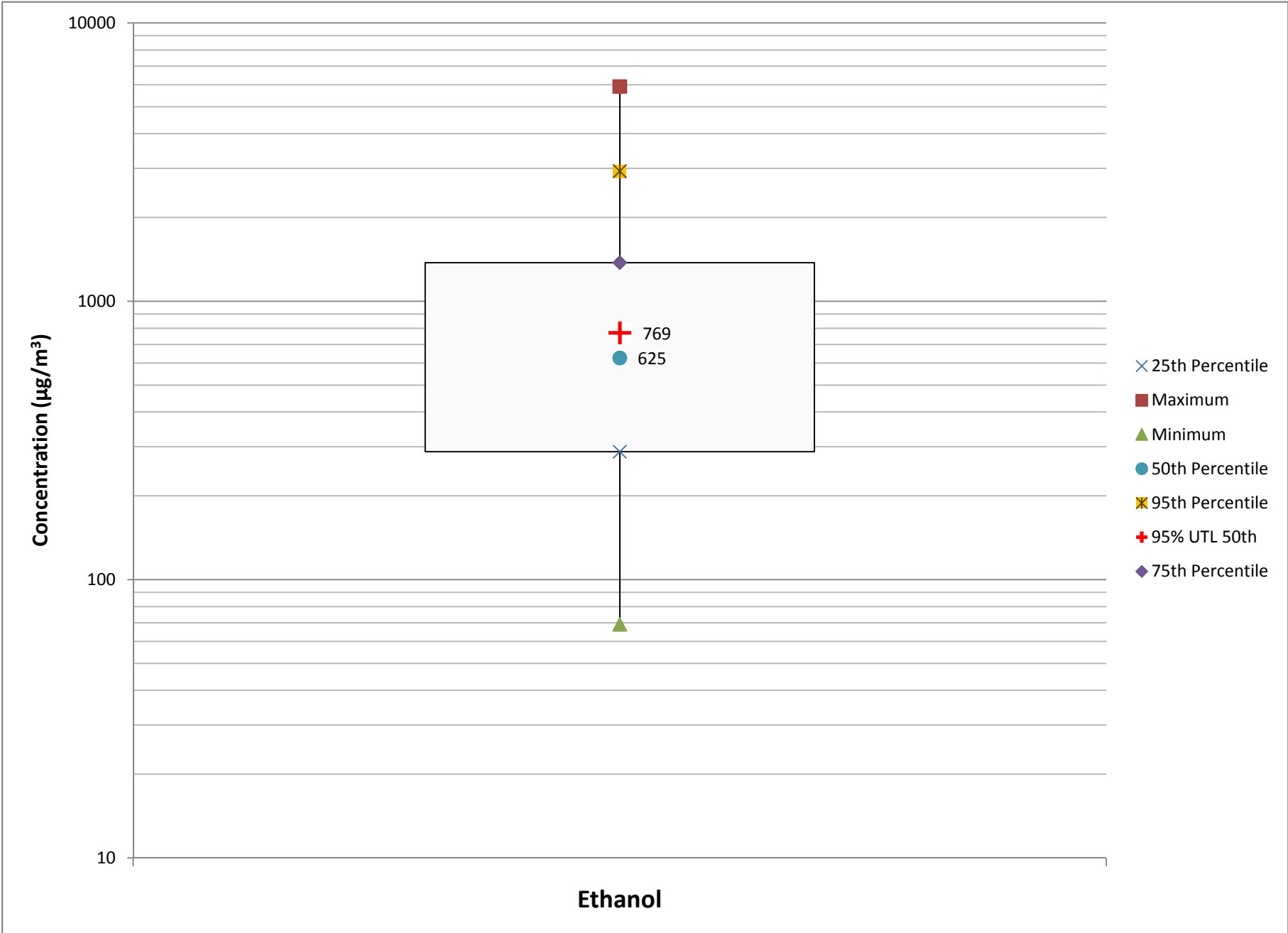
# C9-C12 Aliphatics

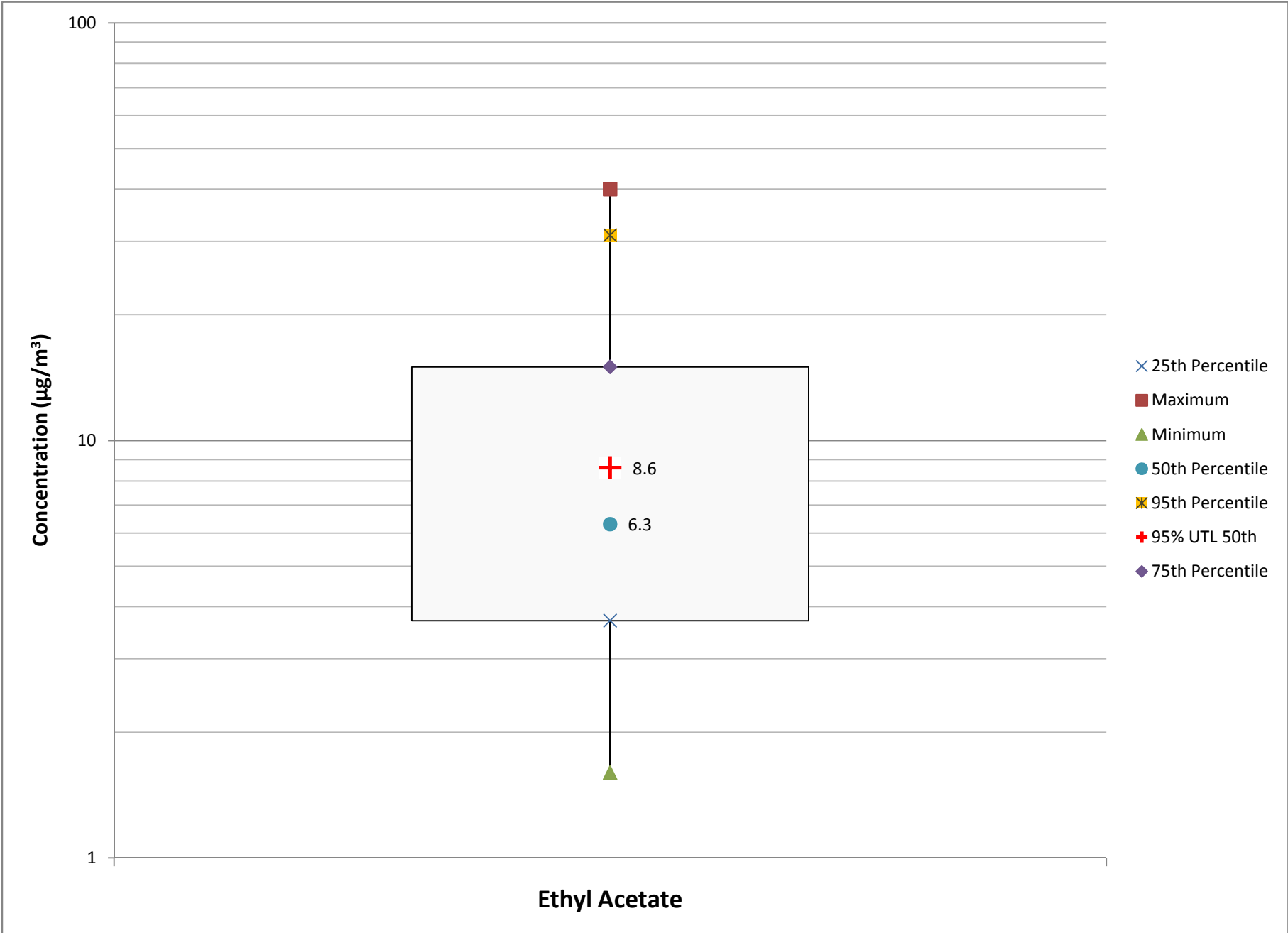




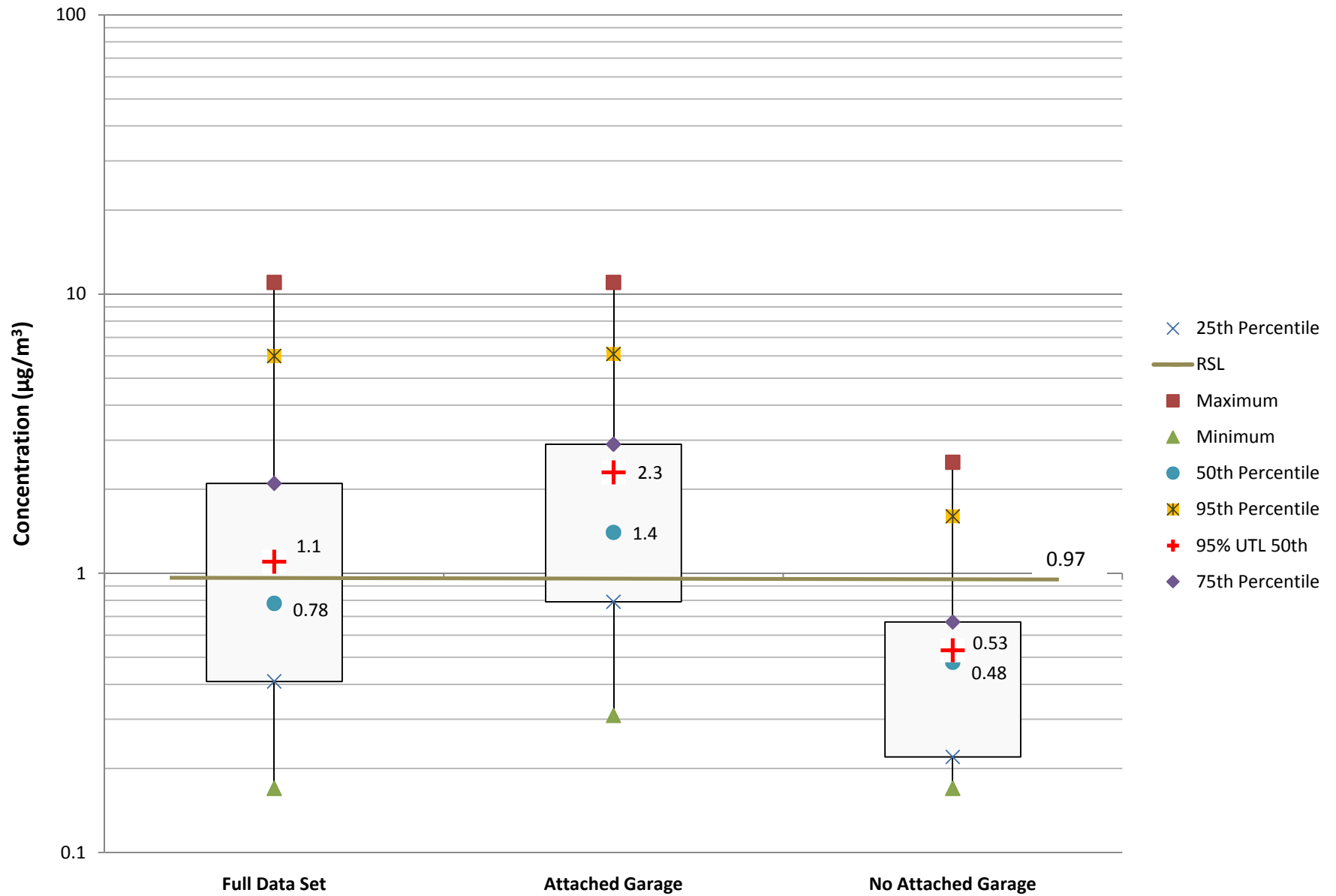


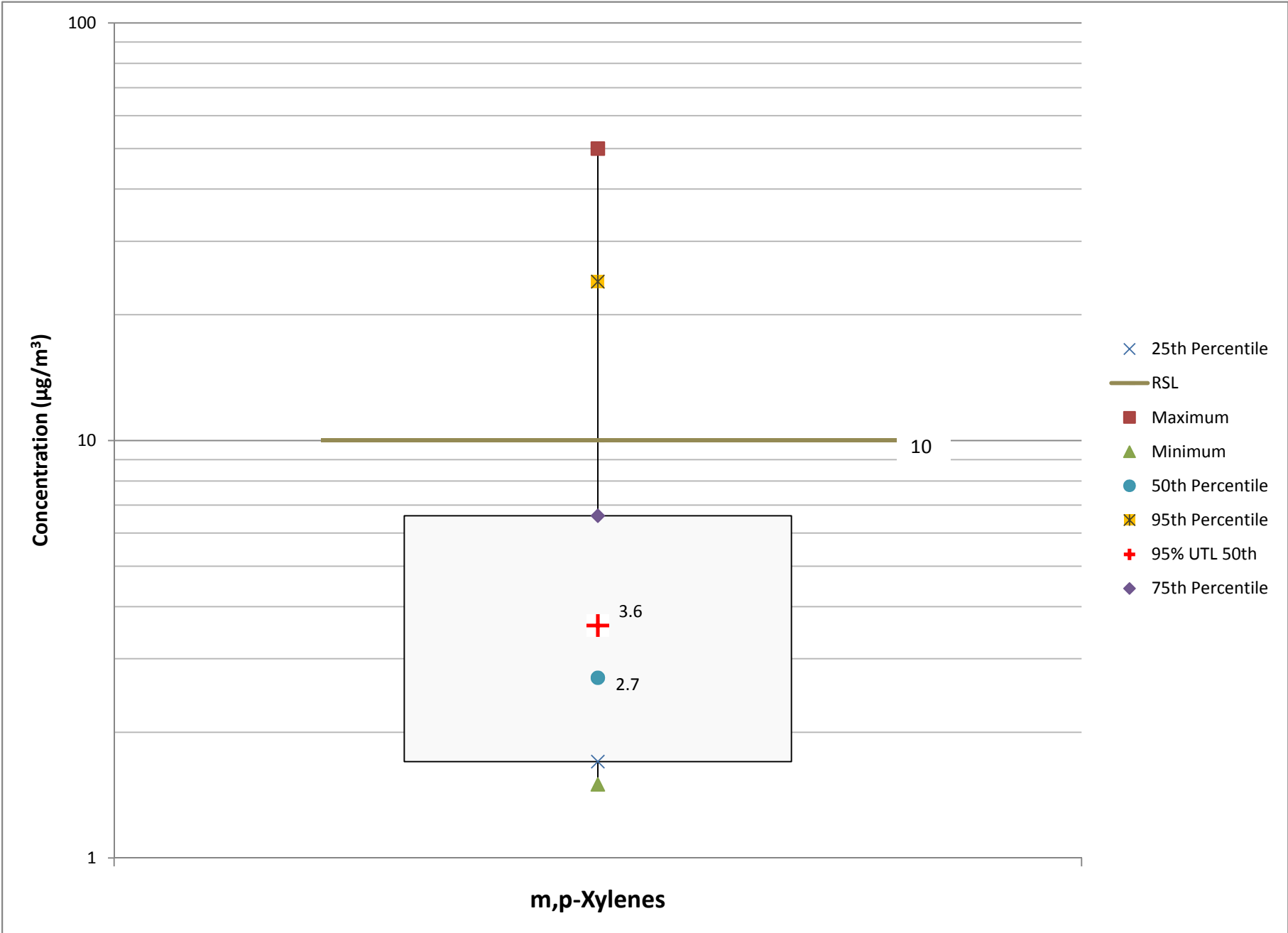






# Ethylbenzene







# Naphthalene

