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November 12, 2019

Mr. Michael Squire
New York State Department of Environmental Conservation
Division of Environmental Remediation
Remedial Bureau C
625 Broadway, 11th Floor
Albany, New York 12233-7014

Re: Soil Vapor Intrusion Investigation Report & Periodic Review Report
Congers Colonial Plaza
285 Route 303
Congers, Clarkstown, Rockland County, New York
NYSDEC # V00456
BEI Job # 060141

Dear Mr. Squire:

Please find the attached Soil Vapor Intrusion Investigation Report (SVIIR) and Periodic Review Report (PRR) for the above referenced subject property. The SVIIR was completed in accordance with BEI's Soil Vapor Extraction System (SVES) Decommissioning Work Plan dated February 9, 2018, approved in a letter correspondence from the New York State Department of Environmental Conservation (NYSDEC) dated March 23, 2018. This SVIIR documents the remedial activities conducted at the subject property to determine if it is appropriate to decommission the two (2) SVES operating at the subject property.

Based upon the analytical results of the vapor intrusion investigation, tetrachloroethene (PCE) and its degradation compounds were detected in soil gas at concentrations warranting continued operation of the SVES. Therefore, the annual inspection of the SVES was conducted and documented in the PRR portion of the report. The SVES were operating properly and the engineering and institutional controls are functioning in accordance with the requirements of the Site Management Plan (SMP) dated December 2013 and approved by NYSDEC and NYSDOH on June 12, 2014.

Congers Colonial Plaza
285 Route 303
Congers, New York
BEI Job # 060141

Page 2

If you have any questions, please contact me at our Summit, NJ office at (908) 918-1702.

Sincerely,



Jeff McCurdy, Project Manager
Brennan Environmental, Inc.

Cc: 285 Route 303, L.L.C., c/o UNLMTD Real Estate Group, Mr. Anthony Siniscalchi (via email)
Mr. Nathan Walz, New York State Dept. of Health, 547 River Street, Troy, NY 12180-2213

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BRENNAN ENVIRONMENTAL, INC.
ENVIRONMENTAL CONSULTING SERVICES

Soil Vapor Intrusion Investigation Report & Periodic Review Report

PER NYSDEC DER-10



Congers Colonial Plaza
285 Route 303
Congers, Town of Clarkstown, Rockland County, New York
Section 35.19 Block 2 Lot 11
September 2019
DEC Case #V00456
BEI Job #060141

Prepared for:
285 Route 303, LLC
c/o UNLMTD Real Estate
200 Washington Street
5th Floor
Hoboken, New Jersey 07030

Prepared by:
Brennan Environmental, Inc.
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TABLE OF CONTENTS

INTRODUCTION.....	1
SECTION I – SOIL VAPOR INTRUSION INVESTIGATION REPORT	1
1.0 HISTORICAL INFORMATION	1
2.0 PHYSICAL SETTING.....	2
2.1 REGIONAL GEOLOGY AND HYDROGEOLOGY	2
2.2 SITE TOPOGRAPHY.....	3
2.3 1,000 FOOT RADIUS LAND USE	3
2.4 BASELINE ECOLOGICAL ASSESSMENT	3
2.4.1 <i>Wellhead Assessment</i>	3
2.4.2 <i>Surface Water Body Assessment</i>	3
2.4.3 <i>Wetland Assessment</i>	4
3.0 TECHNICAL OVERVIEW.....	4
3.1 RELIABILITY OF ANALYTICAL DATA	4
3.2 SITE CONTAMINATION SUMMARY	5
4.0 FINDINGS/RECOMMENDATIONS.....	5
4.1 SVES DECOMMISSIONING ACTIVITIES CONDUCTED.....	5
4.2 SAMPLE RESULTS EVALUATION.....	6
4.3 DATA VALIDATION	7
5.0 SUMMARY	7
SECTION II – PERIODIC REVIEW REPORT	9
1.0 REMEDY PERFORMANCE, EFFECTIVENESS, & PROTECTIVENESS	9
1.1 <i>Engineering Control Systems</i>	9
1.2 <i>Institutional Control Systems</i>	9
1.3 <i>Conclusions</i>	10
2.0 MONITORING PLAN COMPLIANCE REPORT.....	11
3.0 OPERATION & MAINTENANCE PLAN COMPLIANCE.....	11
4.0 CONCLUSIONS AND RECOMMENDATIONS.....	12

TABLE OF CONTENTS

TABLES

SUB-SLAB SOIL VAPOR ANALYTICAL RESULTS – AUGUST 1, 2018	6
TETRACHLOROETHENE – MATRIX 2	7
TRICHLOROETHENE – MATRIX 1	7

FIGURES

SITE LOCATION MAP	1
U.S.G.S. TOPOGRAPHIC QUADRANGLE MAP	2
GIS WETLANDS MAP	3
VI SAMPLE LOCATION MAP	4

APPENDICES

NYSDEC CORRESPONDENCE	1
INDOOR AIR QUALITY QUESTIONNAIRE AND BUILDING INVENTORY	2
HISTORIC PCE CONCENTRATIONS: SOIL VAPOR AND INDOOR AIR	3
DATA USABILITY SUMMARY REPORT	4
INSTITUTIONAL & ENGINEERING CONTROLS CERTIFICATION FORM	5
GENERAL SITE-WIDE INSPECTION CHECKLIST	6
PHOTOGRAPHIC LOG	7
LABORATORY ANALYTICAL REPORT (E18-06141)	8



INTRODUCTION

Brennan Environmental, Inc. (BEI) was retained by 285 Route 303, LLC to prepare this Soil Vapor Intrusion Investigation Report (SVIIR) and Periodic Review Report (PRR) that document the remedial activities conducted at the site identified as 285 Route 303, Section 35.19 Block 2 Lot 11, Congers, Town of Clarkstown, Rockland County, NY (subject property). This SVIIR was prepared in accordance with the New York State Department of Environmental Conservation's (NYSDEC's) Technical Guidance for Site Investigation and Remediation (TGSIR), DER-10, and the New York State Department of Health's (NYSDOH's) Final Guidance for Evaluating Soil Vapor Intrusion in the State of New York dated October 2006 (NYSDOH Final Guidance). The PRR portion of the report was prepared in accordance with the NYSDEC's TGSIR, DER-10.

Section I of this report details the execution of the vapor sampling activities conducted to determine if the soil vapor extraction systems (SVES) operating at the subject property could be decommissioned. The sampling activities were conducted in accordance with BEI's SVES Decommissioning Workplan (DWP) dated February 9, 2018, which was approved by the NYSDEC in a correspondence dated March 23, 2018. Vapor intrusion (VI) investigation activities were conducted at three (3) tenancies at the subject property for which a SVES currently operates to mitigate the VI pathway; Carmen's Laundromat, the former First Class Dry Cleaners, and the former Tutor Time. The results from the VI investigation indicated that chlorinated volatile organic compounds (CVOCs), specifically tetrachloroethene (PCE) and its degradation compounds, remain present in soil vapor and require continued mitigation pursuant to the NYSDOH's Final Guidance.

Since the SVES will remain operational, the annual inspection was conducted. The results of the inspection are documented in the PRR portion of this report (Section II).

SECTION I – SOIL VAPOR INTRUSION INVESTIGATION REPORT

1.0 HISTORICAL INFORMATION

According to the Remedial Action Summary Report (RASR) dated August 2006 by RND Services Inc. (RND) of Nyack, New York, initial remedial activities were conducted at the subject property to address a discharge of residual product and waste during the dismantling of a dry cleaning unit at the former First Class Dry Cleaners in October 2000. CVOCs were reportedly present in these materials. Results of a subsequent indoor air sampling event conducted by RND in the former First Class Dry Cleaners indicated that PCE concentrations were detected above NYSDOH action guidelines. Two (2) SVES were installed in February and March 2001 at the subject building to mitigate the VI pathway. The locations of the SVES are depicted on Figure 4.

RND subsequently conducted soil and groundwater investigation activities. The concrete slab was removed from the northern side of the basement below the former dry cleaner tenancy and

test pits were installed to collect soil samples. Seven (7) overburden monitoring wells and three (3) bedrock monitoring wells were installed to investigate groundwater quality. Soil and groundwater analytical results indicated that PCE concentrations were detected above NYSDEC standards.

Approximately 200 tons of contaminated soil was excavated from below the building slab. However, some impacted soil remained in place due to structural limitations. Potassium permanganate injections were also conducted to remediate residual PCE contaminated groundwater.

Periodic indoor air and sub-slab soil vapor sampling was conducted to monitor PCE concentrations in the following six (6) tenancies at the Congers Colonial Plaza:

- Stay Fit Seniors:
- Hong Kong Kitchen:
- former Tutor Time which is currently vacant:
- DePaulis Enterprises IV, Inc.:
- former Launder Station Laundromat, currently Carmen's Laundromat: and
- former First Class Dry Cleaners which is currently vacant.

A majority of the sampling events were conducted with the SVES operational. Based on the results of the indoor air and soil vapor sampling, the SVES were effectively mitigating the VI pathway.

The NYSDEC authorized no further investigation for groundwater at the subject property in a correspondence dated November 30, 2005 and pursuant to NYSDEC's Decision Document dated December 2011, no further action was required for the subject property with the continued operation of the SVES to treat residual soil vapor contamination and the establishment of a deed restriction limiting the site use at the subject property to restricted residential, commercial or industrial use. A Site Management Plan (SMP) dated December 2013 was approved by NYSDEC and NYSDOH on June 12, 2014.

The SVES have been in operations for over 15 years. An investigation was proposed to determine if the SVES are still needed to mitigate the VI pathway. A SVES DWP detailing the proposed decommissioning sampling activities was submitted to the NYSDEC on February 9, 2018 and approved by NYSDEC on March 23, 2018. NYSDEC's approval letter is included as Appendix 1. The activities were executed in the summer of 2018 and the results of the activities are provided herein.

2.0 PHYSICAL SETTING

2.1 REGIONAL GEOLOGY AND HYDROGEOLOGY

According to RND's Results of the Investigation Work Plan dated September 2003, the subsurface geology at the subject property is composed of variable till deposited by glacier ice

underlain by the Brunswick Formation of the Newark Group. During RND's investigation, bedrock was encountered at depths ranging from 47 to 53 feet below site grade (bsg).

According to Environmental Data Resources' (EDR's) Radius Map Report, the soil component name for the area that includes the subject property is identified as "Watchaug", and is composed of fine sandy loam. This soil type has moderate infiltration rates, with moderately coarse textures. During the groundwater investigation at the subject property, depth to groundwater in monitoring wells was measured at approximately 5 feet below site grade (bsg). According to RND's Results of the Investigation Work Plan dated September 2003, the groundwater flow is to the southwest.

2.2 SITE TOPOGRAPHY

BEI reviewed the USGS Topographic Quadrangle for Haverstraw, New York. According to the topographic map, the subject property is located approximately 241 feet above mean sea level. The portion of the Haverstraw, New York quadrangle depicting the subject property is included as Figure 2.

2.3 1,000 FOOT RADIUS LAND USE

Based on visits to the site and the topographic map, land use within a 1,000-foot radius of the subject property appears to be a mixture of commercial and residential properties. The subject property is bounded by wooded areas to the north and west, Route 303 to the east, across which are commercial properties, and Meola Road to the south, across which are residential properties.

2.4 BASELINE ECOLOGICAL ASSESSMENT

2.4.1 Wellhead Assessment

One (1) supply well is located on the northwest side of the subject property, upgradient of the soil and groundwater contamination. The well is maintained to supply water to the washing machines operated in Carmen's Laundromat. RND collected a water sample from an inlet to the water heater located in the laundromat which was laboratory analyzed for CVOCs. All contaminant concentrations were reportedly not detected (ND) in the supply well sample. Based on this information, no further investigation is proposed for the supply well. No other potable wells were identified within close proximity to the subject property.

2.4.2 Surface Water Body Assessment

BEI reviewed available surface water mapping resources to identify surface water bodies located within one-half mile of the subject property. The nearest surface water body is the Toms Brook, which is located approximately 750 feet west of the subject property. Based upon the distance from the subject property, the Toms Brook is not considered to be a potential receptor.

2.4.3 Wetland Assessment

BEI reviewed the National Wetlands Inventory (NWI) map for wetlands located within one-half mile of the subject property. According to the NWI map, the subject property is mapped as an “Upland” area. Freshwater Emergent Wetlands (Classification PEM1Ed) were identified approximately 900 feet west of the subject property and Freshwater Forested/Shrub Wetlands (Classification PFO1C) were identified approximately 2,200 feet southeast and 2,600 feet west-southwest of the subject property. The portion of the NWI wetland map depicting the subject property is included as Figure 3. Based upon the distance from the subject property, wetlands are not considered to be a potential receptor.

3.0 TECHNICAL OVERVIEW

3.1 RELIABILITY OF ANALYTICAL DATA

Soil vapor sampling activities were performed in accordance with the requirements outlined in the NYSDOH’s Final Guidance. All samples were analyzed by Integrated Analytical Laboratories, Inc. (IAL) of Randolph, New Jersey, a New York certified laboratory (New York Certification #11402). The data quality assurance deliverables conformed to the Non-USEPA CLP Method as specified in the NYSDEC’s TGSIR, DER-10, and were validated by a third party pursuant to NYSDEC requirements. Proper chain of custody documentation were maintained for all samples until delivery to a New York certified laboratory. Minimum detection limits (MDLs) for all compounds analyzed were below the current applicable NYSDOH Air Guidance Values (AGV).

A summary of analytical methods and quality assurance is included in the Analysis Conformance/Nonconformance Summary pages in the laboratory packages. The locations of the sub-slab soil gas, indoor air, and ambient air samples are depicted on Figure 4.

The following “Key” provides the explanations necessary to understand the sampling result summary tables listed in the following report sections.

ABBREVIATION KEY

TVOCC - Indicates Total Volatile Organic Contaminant Concentration.
TOCC - Indicates Total Organic Contaminant Concentration.
MDL - Indicates Method Detection Limit / PQL - Indicates Practical Qualitative Limits.
NS - Indicates compound was Not Sampled for in the analysis.
ND - Indicates compound was analyzed for but not detected at the MDL.
~ - Indicates no standard is listed.
Bold & Italics - Indicate that compound was detected above NJDEP cleanup criteria or standard.
ppm - Indicates parts per million.
ppb - Indicates parts per billion.

All other compounds analyzed in the particular scan were not detected, therefore they are not listed in the summary table

3.2 SITE CONTAMINATION SUMMARY

During remedial investigation activities conducted at the subject property by RND, PCE concentrations were detected above the NYSDOH AGV in the indoor air samples collected from the Tutor Time tenancy and the former tenancy of the First Class Dry Cleaners. Therefore, the SVES were installed in February and March 2001. Subsequent indoor air sampling documented the effectiveness of the SVES in mitigating the VI pathway.

Four (4) sub-slab soil vapor sampling events were conducted between November 2005 and December 2010. The first two (2) events were conducted with the SVES turned off. Analytical results for these sampling event indicated that PCE concentrations were low level, less than 150 $\mu\text{g}/\text{m}^3$. However, soil vapor samples were not collected from below the former First Class Dry Cleaners during either event. Two (2) additional soil vapor sampling events were conducted in March and December 2010 while the SVES were operating. The sample collected from below the former First Class Dry Cleaners indicated that the PCE concentration in soil vapor was 3.1 $\mu\text{g}/\text{m}^3$ while the SVES were operating. A historic PCE concentration table for indoor air and soil vapor is provided as Appendix 3.

4.0 FINDINGS/RECOMMENDATIONS

Pursuant to BEI's approved SVES DWP, to satisfy the NYSDEC's and NYSDOH's requirements for decommissioning the SVES, nine (9) air samples were proposed to be collected over an 8-hour period and analyzed for CVOCs via US EPA Method TO-15. Specifically, five (5) indoor air samples, three (3) sub-slab soil vapor samples, and one (1) ambient air sample were proposed to be collected from the three (3) tenancies of the subject building. The samples were proposed to be collected after the SVES have been shut down for at least 30 days to allow conditions to acclimate. The results would be compared to the NYSDOH's Final Guidance VI Decision Matrices to evaluate if the SVES could be decommissioned. The indoor and ambient air samples would be conditionally analyzed based upon the results of the soil vapor samples. Following sampling, the SVES would be restarted. The NYSDEC's SVES DWP approval letter is provided as Appendix 1.

4.1 SVES DECOMMISSIONING ACTIVITIES CONDUCTED

On April 20, 2018, the SVES at the subject property were powered down in preparation for indoor air and sub-slab soil vapor sampling. On July 31, 2018, BEI was onsite to collect five (5) indoor air samples from the subject property. First-floor indoor air samples, identified as IA-FF-101, IA-FF-102, and IA-FF-103, were collected from Carmen's Laundromat, the former First Class Dry Cleaner, and the former Tutor Time, respectively. Basement-level indoor air samples, identified as IA-B-101 and IA-B-102, were collected from the former First Class Dry Cleaner and the former Tutor Time, respectively. One (1) ambient air sample, identified as AA-101, was collected from the northern exterior of the subject building to document background conditions. A six-liter Summa canister with an 8-hour regulator was placed at breathing zone height to collect each sample. The NYSDOH's Indoor Air Quality Questionnaire and Building Inventory

(Inventory) was completed which included a summary of hazardous substances stored in the tenancies. The completed Inventory is provided as Appendix 2.

On August 1, 2018, BEI returned to the site to collect three (3) sub-slab soil gas samples; SS-101 from Carmen’s Laundromat, SS-102 from the former First Class Dry Cleaner, and SS-103 from the former Tutor Time. A 3/4-inch diameter hole was installed through the concrete building slab at each sampling location to create temporary sampling ports. A six-liter Summa canister equipped with an 8-hour regulator was used to collect each sample. The sampling apparatus was purged with a negative pressure pump prior to sampling. All samples were sent to a New York certified laboratory to be analyzed for CVOCs via US EPA Method TO-15. Each temporary sampling point was repaired with concrete and the SVES were restarted following sampling. A summary of the analytical results for the sub-slab soil vapor samples is provided in Table 1 below and sample locations are depicted on the VI Sample Location Map included as Figure 4.

Table 1: Sub-Slab Soil Vapor Analytical Results – August 1, 2018

Sample ID:	SS-101	SS-102	SS-103
Lab ID:	E18-06141-07	E18-06141-08	E18-06141-09
Matrix:	Air	Air	Air
Date:	8/1/2018	8/1/2018	8/1/2018
Parameter (µg/m ³)			
1,1-Dichloroethene	ND	2	ND
Cis-1,2-Dichloroethene	ND	950	ND
Trans-1,2-Dichloroethene	ND	6.9	ND
Tetrachloroethene	10	10,000	950
Trichloroethene	ND	460	0.93

ND – Not Detected

The analytical results indicated that PCE was detected in sub-slab soil vapor at 10 µg/m³ in SS-101, at 950 µg/m³ in SS-103, and at 10,000 µg/m³ in SS-102. Trichloroethene (TCE), a degradation compound of PCE, was detected in sub-slab soil vapor at 0.93 µg/m³ in SS-103 and 460 µg/m³ in SS-102. The PCE concentrations in SS-101 and SS-103 are in line with previous soil vapor results collected from these locations while the SVES are turned off. SS-102 represents the first soil vapor sample collected from below the former First Class Dry Cleaner while the SVES are not operating. The full data analytical package is provided as Appendix 8.

4.2 SAMPLE RESULTS EVALUATION

In accordance with the NYSDOHs Final Guidance, the sub-slab soil gas concentrations of PCE and TCE were compared to the Soil Vapor/Indoor Air Decision Matrix 2 (Matrix 2) and Soil Vapor/Indoor Air Decision Matrix 1 (Matrix 1), respectively. Since the PCE and TCE concentrations in SS-102, which was collected from the former First Class Cleaner tenancy, exceeded 1,000 µg/m³ and 250 µg/m³, respectively, mitigation is the appropriate remedy regardless of the indoor air concentrations. Therefore, the indoor air and ambient air samples were not analyzed. Continued operation of the SVES are required to mitigate the VI pathway at

the subject property. Matrix 2 and Matrix 1 are provided below as Table 2 and Table 3, respectively.

Table 2: Tetrachloroethene - Matrix 2

		Indoor Air Concentration ($\mu\text{g}/\text{m}^3$)			
		< 3.0	3.0 - 30	30 - 100	> 100
Sub-Slab Vapor Concentration ($\mu\text{g}/\text{m}^3$)	< 100	NFA	Reasonable and Practical Actions	Reasonable and Practical Actions	Reasonable and Practical Actions
	100 - 1,000	Monitor	Monitor/Mitigate	Mitigate	Mitigate
	> 1,000	Mitigate	Mitigate	Mitigate	Mitigate

Table 3: Trichloroethene – Matrix 1

		Indoor Air Concentration ($\mu\text{g}/\text{m}^3$)			
		< 0.25	0.25 - 1.0	1.0 - 5	> 5
Sub-Slab Vapor Concentration ($\mu\text{g}/\text{m}^3$)	< 5.0	NFA	Reasonable and Practical Actions	Reasonable and Practical Actions	Reasonable and Practical Actions
	5.0 - 50.0	NFA	Monitor	Monitor	Mitigate
	50.0 - 250.0	Monitor	Monitor/Mitigate	Mitigate	Mitigate
	> 250	Mitigate	Mitigate	Mitigate	Mitigate

4.3 DATA VALIDATION

In accordance with NYSDEC requirements, a Data Usability Summary Report (DUSR) was generated for the air analytical results from the soil vapor intrusion investigation conducted at the subject property. The DUSR was generated by Alpha Geoscience (Alpha) of Clifton Park, NY. According to Alpha, the analytical data was complete and acceptable with no data flags. Therefore, the results meet the criteria for data quality and data use for the subject property. The DUSR and QA/QC Review generated by Alpha are included as Appendix 4.

5.0 SUMMARY

In accordance with the February 9, 2018 SVES DWP, a soil vapor intrusion investigation was conducted to determine whether the SVES currently in operation at the subject property could be decommissioned. As part of the investigation, five (5) indoor air samples, three (3) soil gas samples, and one (1) ambient air sample were collected from three (3) tenancies of the subject building.

Based on the sub-slab soil vapor results, elevated concentrations of PCE and its degradation compounds remain present in soil gas. Pursuant to the NYSDOH Final Guidance Decision Matrices, mitigation remains the applicable remedy. Therefore, the SVES will remain operational. Since the SVES was observed to be functioning properly and has been demonstrated to be effective at mitigating the VI pathway, no changes are proposed to the monitoring plan. Section II below satisfies the requirements for the annual PRR.

SECTION II – PERIODIC REVIEW REPORT

1.0 REMEDY PERFORMANCE, EFFECTIVENESS, & PROTECTIVENESS

The remedies implemented at the subject property include ECs and ICs to protect human health and the environment. The ECs at the subject property consist of a soil cover system and the SVE system and the ICs consist of site use restrictions. Based on the August 1, 2018 site inspection, the remedy remains protective in accordance with the SMP. An evaluation of the ECs and ICs is provided below.

1.1 Engineering Control Systems

The ECs at the subject property consist of a soil cover system and two (2) SVES. The soil cover system prevents exposure to residual soil contamination left in-place below the former First Class Cleaners tenant space at the subject property. This cover system is comprised of the concrete building slab. The two (2) SVES, VES-1 and VES-2, are installed at the subject property to mitigate PCE-impacted sub-slab soil vapor. VES-1 is located in the basement below the former First Class Dry Cleaners and VES-2 is located in a garage adjacent to the western portion of the former Tutor Time tenancy. In accordance with the NYSDEC correspondence dated November 30, 2005, the SVES are required to be operated until PCE vapors in the indoor air, during non-operation of the VES, are below acceptable NYSDEC and NYSDOH levels.

During the August 1, 2018 site activities, the SVES were restarted following the decommissioning sampling event discussed in Section I of this report. The SVES were inspected and recommissioned as part of this event. VES-1 and VES-2 were both observed to be in good condition and functioning properly. Additionally, the basement floor below the former Tutor Time and former First Class Dry Cleaner was inspected and no breeches were observed. Only small hair-line cracks were observed in the concrete slab in the basement below the former dry cleaners tenancy. Therefore, the soil cover system remains protective of human health and the environment. Based on the inspection, the ECs are performing as designed and remain protective of human health and the environment. The EC certification form is attached as Appendix 5 and the General Site-Wide Inspection Checklist is attached as Appendix 6.

1.2 Institutional Control Systems

A series of ICs is required for the subject property to: (1) implement, maintain and monitor EC systems; (2) prevent future exposure to remaining contamination by controlling disturbances of the subsurface contamination; and (3) limit the use and development of the site to restricted residential, commercial or industrial uses only.

The subject property has a series of ICs in the form of site use restrictions. Adherence to these ICs is required by the Environmental Deed Restriction and is implemented under the SMP. The site use restrictions that apply to the Controlled Property are:

- The subject property may only be used for restricted residential, commercial, or industrial use provided that the long-term Engineering and Institutional Controls included in the SMP are employed;
- The subject property may not be used for a higher level of use, such as unrestricted residential use, without additional remediation and amendment of the Environmental Deed Restriction, as approved by the NYSDEC;
- All future activities on the property that will disturb remaining contaminated material must be conducted in accordance with the SMP;
- The use of the groundwater underlying the subject property is prohibited without treatment rendering it safe for intended use with the exception of the well currently used to provide water to washing machines at the on-site laundromat tenancy;
- The potential for vapor intrusion must be evaluated for any buildings developed at the subject property, and any potential impacts that are identified must be monitored or mitigated;
- Vegetable gardens and farming on the property are prohibited; and
- The site owner or remedial party will submit to NYSDEC a written statement that certifies, under penalty of perjury, that: (1) controls employed at the Controlled Property are unchanged from the previous certification or that any changes to the controls were approved by the NYSDEC; and, (2) nothing has occurred that impairs the ability of the controls to protect public health and environment or that constitute a violation or failure to comply with the SMP. NYSDEC retains the right to access such Controlled Property at any time in order to evaluate the continued maintenance of any and all controls. This certification shall be submitted annually, or an alternate period of time that NYSDEC may allow and will be made by an expert that the NYSDEC finds acceptable.

During BEI's August 1, 2018 site inspection, the site use at the subject property complied with the established ICs. The subject property was operated for commercial use, an acceptable use pursuant to the ICs described above. No new structures have been constructed at the subject property since the establishment of the Environmental Deed Restriction. No activities were conducted at the subject property that disturbed the remaining contaminated material. The IC certification is attached as Appendix 5.

1.3 Conclusions

BEI conducted a site inspection on August 1, 2018. All ECs were inspected and observed to be functioning in accordance with the SMP. The site use at the subject property was in compliance with the ICs detailed in the SMP. Based on this information, all SMP requirements are being achieved and no changes are required. The Institutional and Engineering Controls Certification Form is included as Appendix 5 and the General Site-Wide Inspection Checklist is included as Appendix 6.

2.0 MONITORING PLAN COMPLIANCE REPORT

The Monitoring Plan describes the measurements for evaluating the performance and effectiveness of the remedy to reduce or mitigate contamination at the site, the soil cover system, and all affected site media. Exposure to remaining contamination in soil at the subject property is prevented by a soil cover system placed over the residually contaminated soil. This cover system is comprised of the concrete building slab. Pursuant to the SMP, the Monitoring Plan for the subject property consists of an annual inspection of the soil cover at the subject property. During each inspection, the integrity of the basement floor in the former Tutor Time and former First Class Dry Cleaner is evaluated. Soil cover monitoring is covered separately from the other ECs because it is a passive component of the site remedy. The Monitoring Plan for the active ECs is included with the Operation and Maintenance Plan for these systems in Section 3.0 below.

A site inspection was conducted by BEI on August 1, 2018. The concrete slab basement floor of the former Tutor Time and former First Class Dry Cleaner was inspected and no breeches were observed. Only small hair-line cracks were observed in the concrete slab in the basement of the former Tutor Time, below the former dry cleaner tenant space. Based on the findings of the August 1, 2018 site inspection, the soil cover system remains protective of human health and the environment. The certification form is attached as Appendix 5 and the General Site-Wide Inspection Checklist is attached as Appendix 6.

3.0 OPERATION & MAINTENANCE PLAN COMPLIANCE

The Operation and Maintenance Plan for the subject property describes the measures necessary to operate, monitor, and maintain the mechanical components of the remedy selected for the subject property. Two (2) SVE systems are installed at the subject property to mitigate PCE-impacted soil vapors remaining below the subject building following soil excavation and in-situ remedial activities. VES-1 is located in the basement below the former First Class Dry Cleaners and VES-2 is located in the garage located adjacent to the western portion of the former Tutor Time tenancy. Pursuant to the SMP, the Operation and Maintenance Plan for the subject property consists of an annual inspection of the SVE systems. During each inspection, a visual inspection of the complete system will be conducted and the General Site-Wide Inspection Checklist will be prepared. SVE system components to be monitored include, but are not limited to, the vacuum blower and general system piping.

During the August 1, 2018 site inspection, the SVES were restarted following the vapor sampling event. The systems were inspected and the certification checklist and the General Site-Wide Inspection Checklist were completed. VES-1 and VES-2 were observed to be functioning properly during the site inspection.

The subject property is periodically inspected and maintained by the property owner. During these periodic inspections, the SVES will be checked by the owner's representative to confirm that the systems are operational.

4.0 CONCLUSIONS AND RECOMMENDATIONS

During the August 1, 2018 site inspection, the SVES were inspected and observed to be functioning properly. The site use at the subject property was in compliance with the ICs detailed in the SMP. Based on this information, all requirements set forth in the SMP are being met and no changes to the SMP are required. The Institutional and Engineering Controls Certification Form is included as Appendix 5 and the General Site-Wide Inspection Checklist is included as Appendix 6.

FIGURE 1

Site Location Map

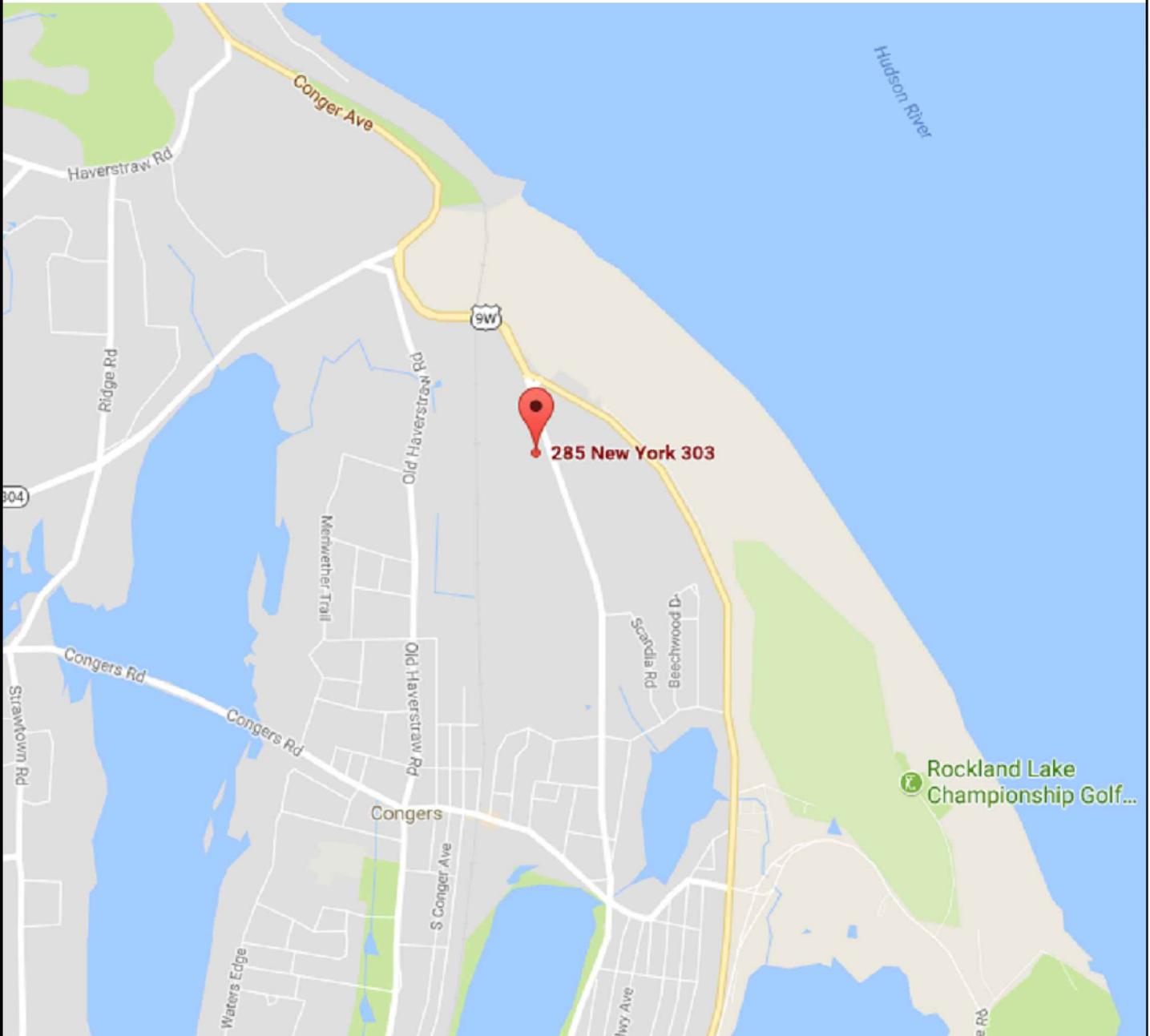


FIGURE: Site Location Map		1	Congers Colonial Plaza 285 Route 303 Congers, Rockland County, NY	
	SOURCE: Google Maps		Scale: As Shown Date: 9/26/2018	Drawn/Checked By: AS/JMc Project #060141
	 Brennan Environmental, Inc. 19 Chatham Road, Summit, New Jersey 07901 908.918.1702			

FIGURE 2

USGS Topographic Quadrangle Map



FIGURE:
USGS Topographic
Quadrangle Map

2

Congers Colonial Plaza
285 Route 303
Congers, Rockland County, NY



SOURCE:
Terrain Navigator

Scale: As Shown
Date: 9/26/2018

Drawn/Checked By: AS/JMc
Project #060141



Brennan Environmental, Inc.
19 Chatham Road, Summit, New Jersey 07901
908.918.1702

FIGURE 3

GIS Wetlands Map

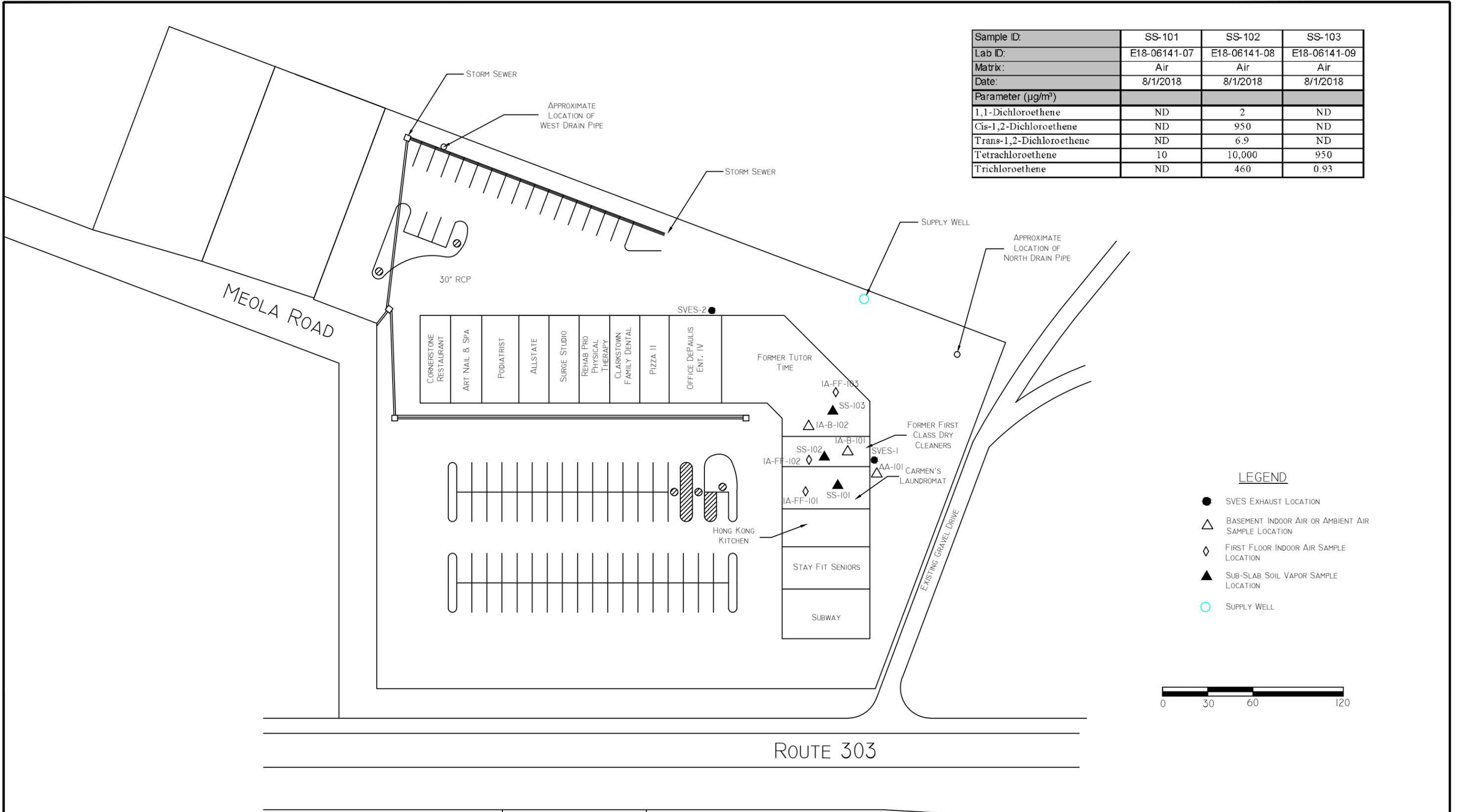


	FIGURE: NWI Wetlands Map	3	Congers Colonial Plaza 285 Route 303 Congers, Rockland County, NY	
			SOURCE: GIS	Scale: As Shown Date: 9/26/2018
			 Brennan Environmental, Inc. 19 Chatham Road, Summit, New Jersey 07901 908.918.1702	

FIGURE 4

VI Decommissioning Sample Location Map

Sample ID:	SS-101	SS-102	SS-103
Lab ID:	E18-06141-07	E18-06141-08	E18-06141-09
Matrix:	Air	Air	Air
Date:	8/1/2018	8/1/2018	8/1/2018
Parameter (µg/m³)			
1,1-Dichloroethene	ND	2	ND
Cis-1,2-Dichloroethene	ND	950	ND
Trans-1,2-Dichloroethene	ND	6.9	ND
Tetrachloroethene	10	10,000	950
Trichloroethene	ND	460	0.93



LEGEND

- SVES EXHAUST LOCATION
- △ BASEMENT INDOOR AIR OR AMBIENT AIR SAMPLE LOCATION
- ◇ FIRST FLOOR INDOOR AIR SAMPLE LOCATION
- ▲ SUB-SLAB SOIL VAPOR SAMPLE LOCATION
- SUPPLY WELL



FIGURE: VI DECOMMISSIONING SAMPLE LOCATION MAP	4	CONGERS COLONIAL PLAZA 285 ROUTE 303 CONGERS, NEW YORK	
		SCALE: 1" = 60' DATE: 9/26/18	DRAWN/CHECKED BY: AS/JMC PROJECT # 060141
SOURCE: RND SERVICES, INC. SAMPLING PLAN, SITE MEASUREMENTS		BRENNAN ENVIRONMENTAL, INC. 19 CHATHAM ROAD, SUMMIT, NEW JERSEY 07901 908.918.1702	

APPENDIX 1

NYSDEC Correspondence

NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION

Division of Environmental Remediation, Remedial Bureau C

625 Broadway, 11th Floor, Albany, NY 12233-7014

P: (518) 402-9662 | F: (518) 402-9679

www.dec.ny.gov

Mr. Anthony Siniscalchi
UNLMTD Real Estate Group
200 Washington Street
Hoboken, NJ 07030

Re: Congers Colonial Plaza
Site No. V00456
Congers, Rockland County, NY
SVE Decommissioning Work Plan

Dear Mr. Siniscalchi:

We have reviewed the Soil Vapor Extraction System Decommissioning Work Plan submitted by a letter of February 9, 2018 from your consultant, Jeff McCurdy of Brennan Environmental, Inc. for the referenced site. Based on our review, the task presented in the work plan appears to be consistent with the outline provided to Jeff McCurdy by Mark Sergott's (of NYSDOH) email of January 26, 2016. The work plan is hereby approved.

Please submit a letter report presenting the findings of the field activities. If you have any questions, please contact me at (518) 402-9662.

Sincerely,



Amen M. Omorogbe, P.E.
Remedial Bureau C
Division of Environmental Remediation

Jeff McCurdy, BEI, jmccurdy@bei-env.com
Mark Sergott, NYSDOH
G. Heitzman, NYSDEC



Department of
Environmental
Conservation

APPENDIX 2

Indoor Air Quality Questionnaire and Building Inventory

**NEW YORK STATE DEPARTMENT OF HEALTH
INDOOR AIR QUALITY QUESTIONNAIRE AND BUILDING INVENTORY
CENTER FOR ENVIRONMENTAL HEALTH**

This form must be completed for each residence involved in indoor air testing.

Preparer's Name Andrew Schmucker Date/Time Prepared 7/31/2018, 8:00 am

Preparer's Affiliation Brennan Environmental, Inc. Phone No. 908-918-1702

Purpose of Investigation To determine if SVES Decommissioning is appropriate

1. OCCUPANT:

The former First Class Dry Cleaners and Former Tutor Time were vacant at the time of sampling. Therefore, no occupants were present for interview during sampling.

Interviewed: Y N

Last Name: Carmen First Name: Roy

Address: 285 Route 303, Congers, NY

County: Rockland

Home Phone: 914-384-1671 Office Phone: _____

Number of Occupants/persons at this location 2 Age of Occupants Adults

2. OWNER OR LANDLORD: (Check if same as occupant)

Interviewed: Y N

285 Route 303, LLC c/o

Last Name: Siniscalchi First Name: Anthony

Address: 285 Route 303, Congers, New York 10920

County: Rockland

Home Phone: N/A Office Phone: _____

3. BUILDING CHARACTERISTICS

Type of Building: (Circle appropriate response)

Residential
Industrial

School
Church

Commercial Multi-use
Other: _____

If the property is residential, type? (Circle appropriate response)

- | | | |
|--------------|-----------------|-------------------|
| Ranch | 2-Family | 3-Family |
| Raised Ranch | Split Level | Colonial |
| Cape Cod | Contemporary | Mobile Home |
| Duplex | Apartment House | Townhouses/Condos |
| Modular | Log Home | Other: _____ |

If multiple units, how many? _____

If the property is commercial, type?

Business Type(s) Commerical Shopping Center

Does it include residences (i.e., multi-use)? Y N If yes, how many? _____

Other characteristics:

Number of floors 2

Building age ~ 30 years

Is the building insulated? Y N

How air tight? Tight Average Not Tight

4. AIRFLOW

Use air current tubes or tracer smoke to evaluate airflow patterns and qualitatively describe:

Airflow between floors

Stairwells, HVAC

Airflow near source

Exterior doorway

Outdoor air infiltration

Doors, windows, and subsurface utilities

Infiltration into air ducts

5. BASEMENT AND CONSTRUCTION CHARACTERISTICS (Circle all that apply)

- a. Above grade construction: wood frame concrete stone brick
- b. Basement type: full crawlspace slab other _____
- c. Basement floor: concrete dirt stone other _____
- d. Basement floor: uncovered covered covered with Linoleum Tile, Carpet
- e. Concrete floor: unsealed sealed sealed with _____
- f. Foundation walls: poured block stone other _____
- g. Foundation walls: unsealed sealed sealed with _____
- h. The basement is: wet damp dry moldy
- i. The basement is: finished unfinished partially finished
- j. Sump present? Y / N
- k. Water in sump? Y / N not applicable (sump sealed)

Basement/Lowest level depth below grade: 10 (feet)

Identify potential soil vapor entry points and approximate size (e.g., cracks, utility ports, drains)

Hairline cracks in basement slab and floor drains in former Tutor Time tenancy.

6. HEATING, VENTING and AIR CONDITIONING (Circle all that apply)

Type of heating system(s) used in this building: (circle all that apply – note primary)

- Hot air circulation
- Space Heaters
- Electric baseboard
- Heat pump
- Stream radiation
- Wood stove
- Hot water baseboard
- Radiant floor
- Outdoor wood boiler
- Other _____

The primary type of fuel used is:

- Natural Gas
- Electric
- Wood
- Fuel Oil
- Propane
- Coal
- Kerosene
- Solar

Domestic hot water tank fueled by: _____

Boiler/furnace located in: Basement Outdoors Main Floor Other _____

Air conditioning: Central Air Window units Open Windows None
Wall-mounted

Are there air distribution ducts present? Y N

Describe the supply and cold air return ductwork, and its condition where visible, including whether there is a cold air return and the tightness of duct joints. Indicate the locations on the floor plan diagram.

Unknown

7. OCCUPANCY

Is basement/lowest level occupied? Full-time Occasionally Seldom Almost Never

Level **General Use of Each Floor (e.g., familyroom, bedroom, laundry, workshop, storage)**

Basement Storage, Utilities, Former Tutor Time (Currently Vacant, Former Day Care facility)

1st Floor Laundromat, Former Day Care Facility, Former Retail Space (Currently Vacant)

2nd Floor _____

3rd Floor _____

4th Floor _____

8. FACTORS THAT MAY INFLUENCE INDOOR AIR QUALITY

a. Is there an attached garage?

Y N

b. Does the garage have a separate heating unit?

Y N NA

c. Are petroleum-powered machines or vehicles stored in the garage (e.g., lawnmower, atv, car)

Y N / NA
Please specify Truck

d. Has the building ever had a fire?

Y N When? _____

e. Is a kerosene or unvented gas space heater present?

Y N Where? _____

f. Is there a workshop or hobby/craft area?

Y N Where & Type? _____

g. Is there smoking in the building?

Y N How frequently? _____

h. Have cleaning products been used recently?

Y N When & Type? _____

i. Have cosmetic products been used recently?

Y N When & Type? _____

- j. Has painting/staining been done in the last 6 months? Y N _____ Where & When? _____
- k. Is there new carpet, drapes or other textiles? Y N _____ Where & When? _____
- l. Have air fresheners been used recently? Y N _____ When & Type? _____
- m. Is there a kitchen exhaust fan? Y N _____ If yes, where vented? _____
- n. Is there a bathroom exhaust fan? Y N _____ If yes, where vented? Outside
- o. Is there a clothes dryer? Y N _____ If yes, is it vented outside? Y N
- p. Has there been a pesticide application? Y N _____ When & Type? _____

Are there odors in the building? Y N
 If yes, please describe: _____ Light mold odors _____

Do any of the building occupants use solvents at work? Y N
 (e.g., chemical manufacturing or laboratory, auto mechanic or auto body shop, painting, fuel oil delivery, boiler mechanic, pesticide application, cosmetologist)

If yes, what types of solvents are used? _____

If yes, are their clothes washed at work? Y / N

Do any of the building occupants regularly use or work at a dry-cleaning service? (Circle appropriate response)

- Yes, use dry-cleaning regularly (weekly) No
- Yes, use dry-cleaning infrequently (monthly or less) Unknown
- Yes, work at a dry-cleaning service

Is there a radon mitigation system for the building/structure? Y N Date of Installation: _____
Is the system active or passive? Active/Passive

9. WATER AND SEWAGE

Water Supply: Public Water Drilled Well Driven Well Dug Well Other: _____

Sewage Disposal: Public Sewer Septic Tank Leach Field Dry Well Other: _____

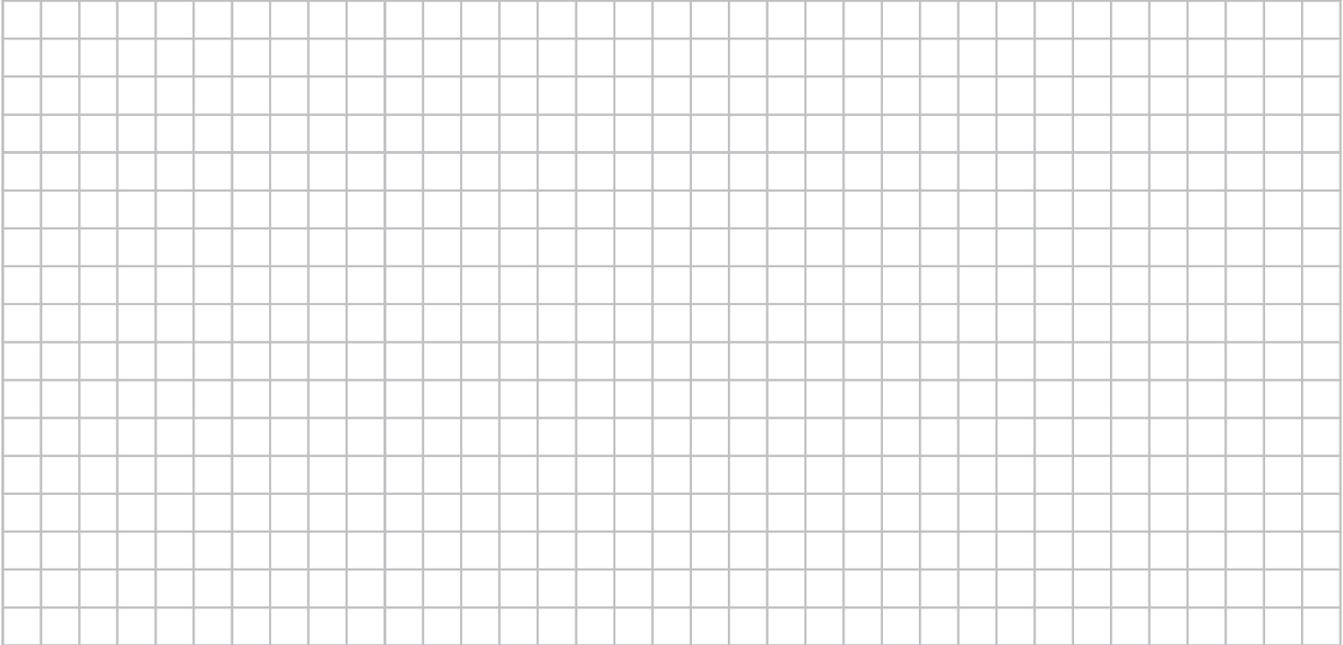
10. RELOCATION INFORMATION (for oil spill residential emergency)

- a. Provide reasons why relocation is recommended: _____
- b. Residents choose to: remain in home relocate to friends/family relocate to hotel/motel
- c. Responsibility for costs associated with reimbursement explained? Y / N
- d. Relocation package provided and explained to residents? Y / N

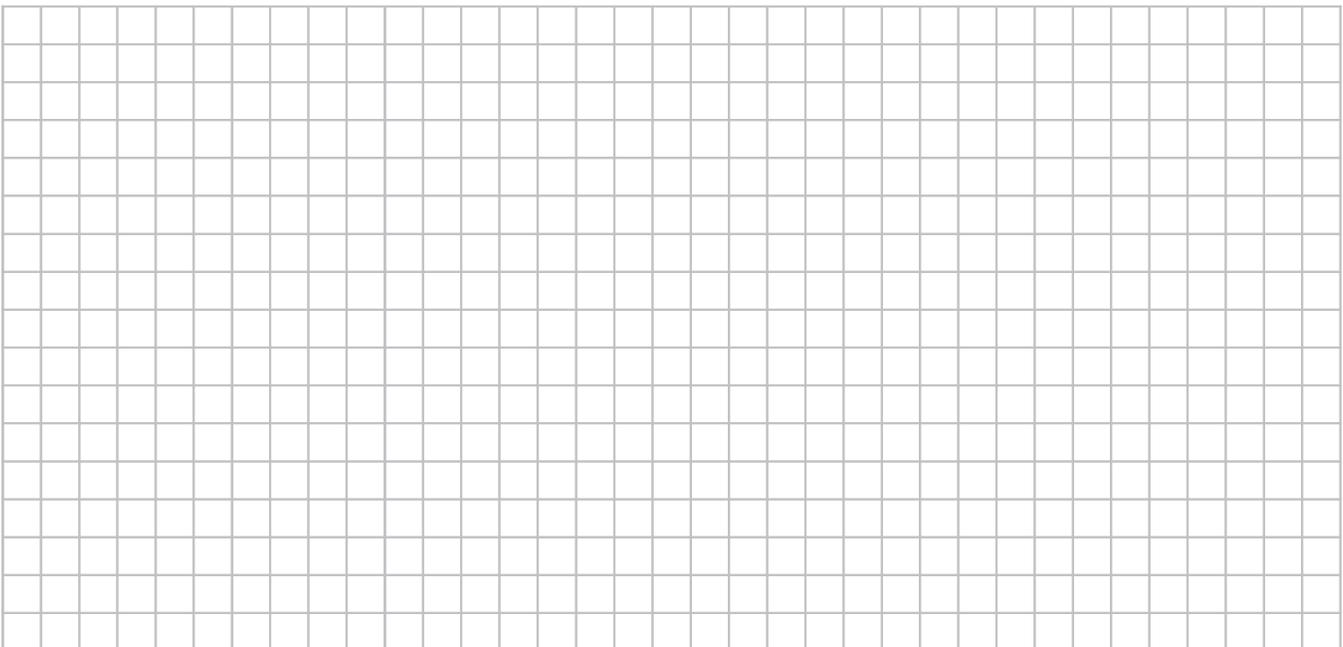
11. FLOOR PLANS

Draw a plan view sketch of the basement and first floor of the building. Indicate air sampling locations, possible indoor air pollution sources and PID meter readings. If the building does not have a basement, please note.

Basement: See attached figure.



First Floor: See attached figure.

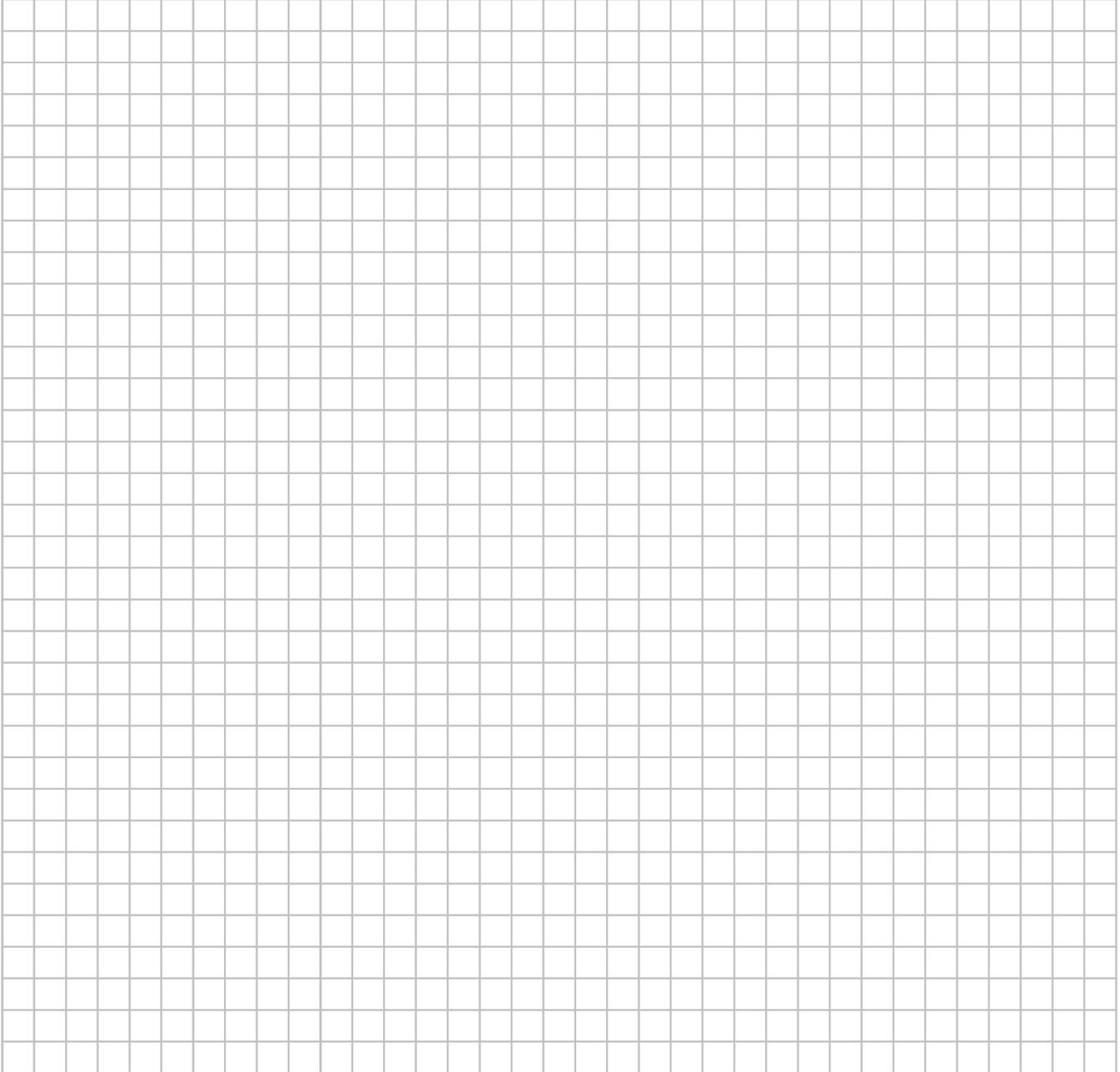


12. OUTDOOR PLOT

Draw a sketch of the area surrounding the building being sampled. If applicable, provide information on spill locations, potential air contamination sources (industries, gas stations, repair shops, landfills, etc.), outdoor air sampling location(s) and PID meter readings.

Also indicate compass direction, wind direction and speed during sampling, the locations of the well and septic system, if applicable, and a qualifying statement to help locate the site on a topographic map.

See attached figure.



13. PRODUCT INVENTORY FORM

Make & Model of field instrument used: N/A

List specific products found in the residence that have the potential to affect indoor air quality.

Location ^{****}	Product Description	Size ^{***} (units)	Condition [*]	Chemical Ingredients	Field Instrument Reading (units)	Photo ^{**} <u>Y/N</u>
Carmen's	Bleach	< 1 gal	UO		N/A	N
Carmen's	Toilet Scrubber	< 1 gal	U		N/A	N
Carmen's	Softsoap	< 1 gal	U		N/A	N
Carmen's	Propane	< 1 gal	U		N/A	N
Carmen's	CRC Lectra Clean	< 1 gal	U		N/A	N
Carmen's	Glade Air Freshener	< 1 gal	U		N/A	N
Carmen's	Air Duster	< 1 gal	U		N/A	N
Carmen's	Spray 9	< 1 gal	U		N/A	N
Carmen's	HDV Sprayer	< 1 gal	U		N/A	N
Carmen's	Spic & Span	< 1 gal	U		N/A	N
Carmen's	Lemon Pledge	< 1 gal	U		N/A	N
Carmen's	WD-40	< 1 gal	U		N/A	N
Carmen's	Fabuloso	< 1 gal	U		N/A	N
Carmen's	Germicidal Bleach	< 1 gal	U		N/A	N
Carmen's	ZEP Glass Cleaner	< 1 gal	U		N/A	N
Carmen's	HPV All Purpose Cleaning Vinegar	< 1 gal	U		N/A	N
Carmen's	Roto-Rooter Pipe Shield	< 1 gal	U		N/A	N
Carmen's	Laundry Detergent	< 1 gal	UO		N/A	N

* Describe the condition of the product containers as **Unopened (UO)**, **Used (U)**, or **Deteriorated (D)**

** Photographs of the **front and back** of product containers can replace the handwritten list of chemical ingredients. However, the photographs must be of good quality and ingredient labels must be legible.

*** Only small household quantities (<1 gallon per type) of these materials were identified in the tenancy.

**** The former First Class Dry Cleaners and Former Tutor Time were vacant at the time of sampling. No products were identified within these tenancies.

APPENDIX 3

Historic PCE Concentrations: Soil Vapor and Indoor Air

Historic PCE Concentrations: Soil Vapor and Indoor Air
 Congers Colonial Plaza
 285 Route 303
 Congers, New York
 BEI Job #060141

Sample Location	Sample ID	12/7/2000	12/18/2000	1/11/2001	1/18/2001	3/16/2001	4/11/2001	5/14/2001	6/16/2001	7/16/2001	10/29/2001	11/29/2001	12/3/2001	4/30/2002	7/22/2002	11/11/2002	2/4/2003	11/30/2003*	4/8/2004**	6/7/2004**	11/27/2004*	11/27-28/2005*	8/25/2009*	3/25/2010	12/6/2010	8/11/2018*
First Floor Indoor Air																										
Laundry Station Laundromat	Laundromat, IA-1	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	13	4.2	ND	~	~	
Hong Kong Kitchen	IA-2	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	1.6	ND	~	~	
Stay Fit Seniors	IA-3	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	ND	ND	~	~	
First Class Dry Cleaner	Up Dry, IA-5	1.435	0.9	0.14	0.016	0.05	41	66	~	~	~	~	~	~	~	~	~	~	~	~	13	~	8.1	~	~	
Tutor Time	IA-6	~	~	~	~	0.008	33	14	16	6	5.8	<2.1	<2.1	4	<2.0	<2.1	<2.1	118	86	5.2	7.7	~	~	~	ND	~
Depaulis IV	Depaulis IV	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	4.2	~	~	~	~	
Basement Indoor Air																										
Tutor Time	IA-4	~	~	~	~	0.012	46	4	32	13	7.3	<2.1	<2.1	4.7	9.5	2.8	<2.1	87	46	12	27	~	8.1	2.1	ND	~
First Class Dry Cleaner	Down Dry, IA-5 [^]	1.726	1.3	0.11	0.012	0.07	234	101	~	~	~	~	~	~	~	~	~	~	~	~	21	~	~	ND	~	
Depaulis IV	Garage	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	3.1	~	~	~	~	
Sub-Slab Soil Vapor																										
Laundry Station Laundromat	SS-1, S-101	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	17	1.7	~	10	
Hong Kong Kitchen	SS-2	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	5	1.8	~	~	
Stay Fit Seniors	SS-3	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	24	1.9	~	~	
Tutor Time	SS Toddler A, SS-4, SS-103	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	140	~	~	ND	950
First Class Dry Cleaners	SS-5, SS-102	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	3.1	10,000	
Depaulis IV	SS-VES2F	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	3.2	~	~	~	~	
VES-1 Exhaust	SS-VES1G	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	32	~	~	~	~
VES-2 Exhaust	SS-VES2G	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	0.69	~	~	~	~	
Ambient Air																										
VES-1 Exhaust	AA-1	<0.067	~	~	~	<0.0004	<2	~	19	20	2.9	<2.1	<2.1	<2.0	<2.1	<2.4	~	2	78	7.9	6.8	~	ND	ND	ND	~
VES-2 Exhaust	AA-2	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	ND	1.8	~	~
Outside Tutor Time	Background	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	~	ND	~	~	~	~

[^] - Please note that the indoor air sample collected from the basement of the former First Class Dry Cleaner was mistakenly identified by the identical sample name as the first floor, IA-5, on the chain of custody. In order to distinguish the results, the basement sample is identified herein as IA-5[^].

Sample results in µg/m³

ND - Not detected at laboratory MDL

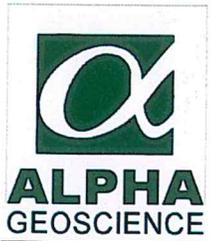
~ - Not sampled on this date

* - Samples collected with VES-1 and VES-2 turned off

** - Elevated results attributed to clogged filter; filters changed and locations resampled on June 7, 2004

APPENDIX 4

Data Usability Summary Report



Geology

Hydrology

Remediation

Water Supply

RECEIVED
SEP 24 2018

BY:

September 20, 2018

Mr. Jeff McCurdy
Brennan Environmental, Inc.
19 Chatham Road
Summit, New Jersey 07901

Re: Data Validation Report
Congers Colonial Plaza
July 2018 Sub Slab Air Samples

Dear Mr. McCurdy:

The data usability summary report (DUSR) and data validation summary are attached to this letter for Congers Colonial Plaza, July 2018 sub slab air samples. The data for Integrated Analytical Laboratories LLC, SDG# E18-06141, are acceptable with no issues that are identified and discussed in the validation summary. There are no data that are qualified as either estimated (J) or rejected (R) in the data pack.

A list of common data validation acronyms is attached to this letter to assist you in interpreting the validation summaries. If you have any questions concerning the work performed, please contact me at (518) 348-6995. Thank you for the opportunity to assist Brennan Environmental, Inc.

Sincerely,
Alpha Geoscience

Donald Anné
Senior Chemist

DCA:dca
attachments

Z:\projects\2010\10600 - 10620\10615-congers colonial plaza\2018\Congeners Colonial Plaza-181.ltr.wpd

Data Validation Acronyms

AA	Atomic absorption, flame technique
BHC	Hexachlorocyclohexane
BFB	Bromofluorobenzene
CCB	Continuing calibration blank
CCC	Calibration check compound
CCV	Continuing calibration verification
CN	Cyanide
CRDL	Contract required detection limit
CRQL	Contract required quantitation limit
CVAA	Atomic adsorption, cold vapor technique
DCAA	2,4-Dichlophenylacetic acid
DCB	Decachlorobiphenyl
DFTPP	Decafluorotriphenyl phosphine
ECD	Electron capture detector
FAA	Atomic absorption, furnace technique
FID	Flame ionization detector
FNP	1-Fluoronaphthalene
GC	Gas chromatography
GC/MS	Gas chromatography/mass spectrometry
GPC	Gel permeation chromatography
ICB	Initial calibration blank
ICP	Inductively coupled plasma-atomic emission spectrometer
ICV	Initial calibration verification
IDL	Instrument detection limit
IS	Internal standard
LCS	Laboratory control sample
LCS/LCSD	Laboratory control sample/laboratory control sample duplicate
MSA	Method of standard additions
MS/MSD	Matrix spike/matrix spike duplicate
PID	Photo ionization detector
PCB	Polychlorinated biphenyl
PCDD	Polychlorinated dibenzodioxins
PCDF	Polychlorinated dibenzofurans
QA	Quality assurance
QC	Quality control
RF	Response factor
RPD	Relative percent difference
RRF	Relative response factor
RRF(number)	Relative response factor at concentration of the number following
RT	Retention time
RRT	Relative retention time
SDG	Sample delivery group
SPCC	System performance check compound
TCX	Tetrachloro-m-xylene
%D	Percent difference
%R	Percent recovery
%RSD	Percent relative standard deviation

Data Validation Qualifiers Used in the QA/QC Reviews for USEPA Region II

- U = Not detected. The associated number indicates the approximate sample concentration necessary to be detected significantly greater than the level of the highest associated blank.
- R = Unreliable result; data is rejected or unusable. Analyte may or may not be present in the sample. Supporting data or information is necessary to confirm the result.
- N = Tentative identification. Analyte is considered present. Special methods may be needed to confirm its presence or absence during future sampling efforts.
- J = Analyte is present. Reported value may be associated with a higher level of uncertainty than is normally expected with the analytical method.
- J- = Analyte is present. Reported value may be biased low and associated with a higher level of uncertainty than is normally expected with the analytical method.
- J+ = Analyte is present. Reported value may be biased high and associated with a higher level of uncertainty than is normally expected with the analytical method.
- UJ = Not detected, quantitation limit may be inaccurate or imprecise.

Note: These qualifiers are used for data validation purposes. The data validation qualifiers may differ from the qualifiers that the laboratory assigns to the data. Refer to the laboratory analytical report for the definitions of the laboratory qualifiers.



Geology

Hydrology

Remediation

Water Supply

**Data Usability Summary Report for
Integrated Analytical Laboratories LLC
SDG # E18-06141**

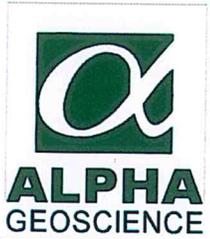
**3 Sub Slab Air Samples
Collected July 31, 2018**

Prepared by: Donald Anné
September 20, 2018

The data package contains the documentation required by NYSDEC ASP. The proper chain of custody procedures were followed by the samplers. All information appeared legible and complete. The data pack contained the results of TO15 volatile analyses for 3 sub slab air samples.

The overall performances of the analyses are acceptable. Integrated Analytical Laboratories LLC did fulfill the requirements of the analytical method.

The data are acceptable with no issues that are identified in the accompanying data validation reviews. There are no data that were qualified as either estimated (J) or rejected (R); therefore, all data are considered usable. Detailed information on data quality is included in the data validation review.



Geology

Hydrology

Remediation

Water Supply

**QA/QC Review of TO15 Volatiles Data for
Integrated Analytical Laboratories LLC
SDG # E18-06141**

**3 Sub Slab Air Samples
Collected July 31, 2018**

Prepared by: Donald Anné
September 20, 2018

Holding Times: Samples were analyzed within the EPA recommended holding times.

GC/MS Tuning and Mass Calibration: The BFB tuning criteria were within control limits.

Initial Calibration: The average RRFs for target compounds were above the allowable minimum (0.050), as required.

Continuing Calibration: The CCRFs for target compounds were above the allowable minimum (0.050) and the %Ds were below the allowable maximum (30%), as required.

Blanks: The analyses of the method blanks reported target compounds as not detected.

Internal Standard Area Summary: The internal standard areas and retention times were within control limits.

Laboratory Duplicate Sample: The relative percent differences for detected compounds were below the allowable maximum (30%) in laboratory duplicate samples E18-04189-01, E18-06204-01, and E18-06173-01, as required.

Compound ID: Checked compounds were within GC quantitation limits. The mass spectra for detected compounds contained the primary and secondary ions, as outlined in the method.

Clean Cannister Verification: The analysis of clean cannister verification samples reported target compounds as not detected, as required.

Canister Pressure: The laboratory reported all samples with residual vacuums, as required.

Z:\projects\2010\10600 - 10620\10615-congers colonial plaza\2018\E18-06141.t15.wpd



Integrated Analytical Laboratories LLC

Summary of Results

Brennan Environmental
 19 Chatham Road
 Summit, NJ 07901
 Attn: Jeff McCurdy
 Project: Congers CP / 060141
 Site: NY

Report Date: 08/13/18
 SDG Number: E18-06141
 Date Sampled: 08/01/18
 Date Received: 08/02/18
 Date Analyzed: 08/06/18
 Data File: AA8206
 Summa ID: 3014A
 DF: 1

Analysis: Volatile Organic Compounds by EPA Method TO-15

<u>Compound</u>	<u>CAS #</u>	<u>Sample Name:</u> <u>IAL ID:</u>		<u>Reporting Limits</u>	
		<u>ppbv</u>	<u>ug/m3</u>	<u>ppbv</u>	<u>ug/m3</u>
1,1-Dichloroethane	75-34-3	ND	ND	0.20	0.81
1,2-Dichloroethane	107-06-2	ND	ND	0.20	0.81
1,1-Dichloroethene	75-35-4	ND	ND	0.20	0.79
1,2-Dichloroethene (cis)	156-59-2	ND	ND	0.20	0.79
1,2-Dichloroethene (trans)	156-60-5	ND	ND	0.20	0.79
1,1,2,2-Tetrachloroethane	79-34-5	ND	ND	0.20	1.4
Tetrachloroethene	127-18-4	1.5	10	0.20	1.4
1,1,1-Trichloroethane	71-55-6	ND	ND	0.20	1.1
1,1,2-Trichloroethane	79-00-5	ND	ND	0.20	1.1
Trichloroethene	79-01-6	ND	ND	0.05	0.25
Vinyl chloride	75-01-4	ND	ND	0.20	0.51



Integrated Analytical Laboratories LLC

Summary of Results

Brennan Environmental
 19 Chatham Road
 Summit, NJ 07901
 Attn: Jeff McCurdy
 Project: Congers CP / 060141
 Site: NY

Report Date: 08/13/18
 SDG Number: E18-06141
 Date Sampled: 08/01/18
 Date Received: 08/02/18
 Date Analyzed: 08/06/18, 08/07/18
 Data File: AA8207, AA8232
 Summa ID: 2155
 DF: 1, 100

Analysis: Volatile Organic Compounds by EPA Method TO-15

<u>Compound</u>	<u>CAS #</u>	<u>IAL ID:</u>	<u>Sample Name:</u>		<u>Reporting Limits</u>	
			<u>SS-102</u>	<u>E18-06141-08</u>	<u>ppbv</u>	<u>ug/m3</u>
			<u>ppbv</u>	<u>ug/m3</u>	<u>ppbv</u>	<u>ug/m3</u>
1,1-Dichloroethane	75-34-3		ND	ND	0.20	0.81
1,2-Dichloroethane	107-06-2		ND	ND	0.20	0.81
1,1-Dichloroethene	75-35-4		0.50	2.0	0.20	0.79
1,2-Dichloroethene (cis)	156-59-2	D	240	950	20	79
1,2-Dichloroethene (trans)	156-60-5		1.8	6.9	0.20	0.79
1,1,2,2-Tetrachloroethane	79-34-5		ND	ND	0.20	1.4
Tetrachloroethene	127-18-4	D	1500	10000	20	136
1,1,1-Trichloroethane	71-55-6		ND	ND	0.20	1.1
1,1,2-Trichloroethane	79-00-5		ND	ND	0.20	1.1
Trichloroethene	79-01-6	D	86	460	4.6	25
Vinyl chloride	75-01-4		ND	ND	0.20	0.51



Integrated Analytical Laboratories LLC

Summary of Results

Brennan Environmental
 19 Chatham Road
 Summit, NJ 07901
 Attn: Jeff McCurdy
 Project: Congers CP / 060141
 Site: NY

Report Date: 08/13/18
 SDG Number: E18-06141
 Date Sampled: 08/01/18
 Date Received: 08/02/18
 Date Analyzed: 08/06/18, 08/07/18
 Data File: AA8208, AA8233
 Summa ID: 5089
 DF: 1, 10

Analysis: Volatile Organic Compounds by EPA Method TO-15

<u>Compound</u>	<u>CAS #</u>	<u>Sample Name:</u> <u>IAL ID:</u>		<u>Reporting Limits</u>	
		<u>ppbv</u>	<u>ug/m3</u>	<u>ppbv</u>	<u>ug/m3</u>
1,1-Dichloroethane	75-34-3	ND	ND	0.20	0.81
1,2-Dichloroethane	107-06-2	ND	ND	0.20	0.81
1,1-Dichloroethene	75-35-4	ND	ND	0.20	0.79
1,2-Dichloroethene (cis)	156-59-2	ND	ND	0.20	0.79
1,2-Dichloroethene (trans)	156-60-5	ND	ND	0.20	0.79
1,1,2,2-Tetrachloroethane	79-34-5	ND	ND	0.20	1.4
Tetrachloroethene	127-18-4	D	140	950	14
1,1,1-Trichloroethane	71-55-6	ND	ND	0.20	1.1
1,1,2-Trichloroethane	79-00-5	ND	ND	0.20	1.1
Trichloroethene	79-01-6	0.17	0.93	0.05	0.25
Vinyl chloride	75-01-4	ND	ND	0.20	0.51

APPENDIX 5

Institutional & Engineering Controls Certification Form



Enclosure 2
NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION
Site Management Periodic Review Report Notice
Institutional and Engineering Controls Certification Form



	Site Details	Box 1
Site No.	V00456	
Site Name Congers Colonial Plaza		
Site Address: 285 Route 303 Zip Code: 10920-		
City/Town: Congers		
County: Rockland		
Site Acreage: 2.8		
Reporting Period: July 31, 2017 to July 31, 2018		
		YES NO
1. Is the information above correct?		<input checked="" type="checkbox"/> <input type="checkbox"/>
If NO, include handwritten above or on a separate sheet.		
2. Has some or all of the site property been sold, subdivided, merged, or undergone a tax map amendment during this Reporting Period?		<input type="checkbox"/> <input checked="" type="checkbox"/>
3. Has there been any change of use at the site during this Reporting Period (see 6NYCRR 375-1.11(d))?		<input type="checkbox"/> <input checked="" type="checkbox"/>
4. Have any federal, state, and/or local permits (e.g., building, discharge) been issued for or at the property during this Reporting Period?		<input type="checkbox"/> <input checked="" type="checkbox"/>
If you answered YES to questions 2 thru 4, include documentation or evidence that documentation has been previously submitted with this certification form.		
5. Is the site currently undergoing development?		<input type="checkbox"/> <input checked="" type="checkbox"/>

	Box 2
	YES NO
6. Is the current site use consistent with the use(s) listed below? Restricted-Residential, Commercial, and Industrial	<input checked="" type="checkbox"/> <input type="checkbox"/>
7. Are all ICs/ECs in place and functioning as designed?	<input checked="" type="checkbox"/> <input type="checkbox"/>
IF THE ANSWER TO EITHER QUESTION 6 OR 7 IS NO, sign and date below and DO NOT COMPLETE THE REST OF THIS FORM. Otherwise continue.	
A Corrective Measures Work Plan must be submitted along with this form to address these issues.	
Signature of Owner, Remedial Party or Designated Representative	Date

Description of Institutional Controls

<u>Parcel</u>	<u>Owner</u>	<u>Institutional Control</u>
35.19-2-11	285 Route 303, LLC and 119 Route 46, LLC	Ground Water Use Restriction Landuse Restriction Monitoring Plan Site Management Plan O&M Plan Soil Management Plan IC/EC Plan

The owner of the property shall prohibit the property from ever being used for purposes other than for restricted residential, commercial or industrial use without the express written waiver of such prohibition by the Department or Relevant Agency.

The owner of the property shall prohibit the use of the groundwater underlying the property without treatment rendering it safe for drinking water or industrial purposes, as appropriate, unless the user first obtains permission to do so from the Department or Relevant Agency.

The owner of the property shall provide a periodic certification, prepared and submitted by a professional engineer or environmental professional acceptable to the Department or Relevant Agency, which will certify that the institutional and engineering controls put in place are unchanged from the previous certification, comply with the SMP, and have not been impaired.

Description of Engineering Controls

<u>Parcel</u>	<u>Engineering Control</u>
35.19-2-11	Vapor Mitigation Cover System
soil vapor extraction (SVE) Systems Cover System	

Periodic Review Report (PRR) Certification Statements

1. I certify by checking "YES" below that:

a) the Periodic Review report and all attachments were prepared under the direction of, and reviewed by, the party making the certification;

b) to the best of my knowledge and belief, the work and conclusions described in this certification are in accordance with the requirements of the site remedial program, and generally accepted engineering practices; and the information presented is accurate and complete.

YES NO

2. If this site has an IC/EC Plan (or equivalent as required in the Decision Document), for each Institutional or Engineering control listed in Boxes 3 and/or 4, I certify by checking "YES" below that all of the following statements are true:

(a) the Institutional Control and/or Engineering Control(s) employed at this site is unchanged since the date that the Control was put in-place, or was last approved by the Department;

(b) nothing has occurred that would impair the ability of such Control, to protect public health and the environment;

(c) access to the site will continue to be provided to the Department, to evaluate the remedy, including access to evaluate the continued maintenance of this Control;

(d) nothing has occurred that would constitute a violation or failure to comply with the Site Management Plan for this Control; and

(e) if a financial assurance mechanism is required by the oversight document for the site, the mechanism remains valid and sufficient for its intended purpose established in the document.

YES NO

**IF THE ANSWER TO QUESTION 2 IS NO, sign and date below and
DO NOT COMPLETE THE REST OF THIS FORM. Otherwise continue.**

A Corrective Measures Work Plan must be submitted along with this form to address these issues.

Signature of Owner, Remedial Party or Designated Representative

Date

IC CERTIFICATIONS
SITE NO. V00456

Box 6

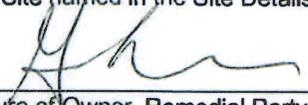
SITE OWNER OR DESIGNATED REPRESENTATIVE SIGNATURE

I certify that all information and statements in Boxes 1,2, and 3 are true. I understand that a false statement made herein is punishable as a Class "A" misdemeanor, pursuant to Section 210.45 of the Penal Law.

I Gabriella Lobente at 200 Washington Street Hirschon NJ 07030
print name print business address

am certifying as owner (Owner or Remedial Party)

for the Site named in the Site Details Section of this form.


Signature of Owner, Remedial Party, or Designated Representative
Rendering Certification

11/1/2019
Date

IC/EC CERTIFICATIONS

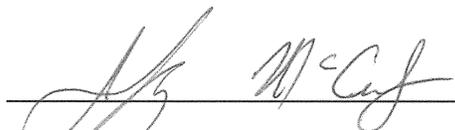
Box 7

Qualified Environmental Professional Signature

I certify that all information in Boxes 4 and 5 are true. I understand that a false statement made herein is punishable as a Class "A" misdemeanor, pursuant to Section 210.45 of the Penal Law.

I Jeff McCurdy at 19 Chatham Road, Summit, NJ 07901,
print name print business address

am certifying as a Qualified Environmental Professional for the Owner
(Owner or Remedial Party)



Signature of Qualified Environmental Professional, for
the Owner or Remedial Party, Rendering Certification

Stamp
(Required for PE)

9/13/19

Date

APPENDIX 6

General Site-Wide Inspection Checklist

General Site-Wide Inspection Checklist

Date of Inspection: August 1, 2018

Inspector Name: Andrew Schmucker

Inspector Company: Brennan Environmental, Inc.

Inspector Position: Environmental Scientist

Site Name: Congers Colonial Plaza

Site Address: 285 Route 303, Congers, NY

	Yes	No	NA
1.0 Is all PVC piping in good condition with no breaks, cracks or leaks?	<u>X</u>	<u> </u>	<u> </u>
2.0 Are openings around the suction point piping penetrations of the slab properly sealed?	<u>X</u>	<u> </u>	<u> </u>
3.0 Are accessible openings around utility penetrations of the foundation walls and slab, test holes, wells and other openings in slabs properly sealed?	<u>X</u>	<u> </u>	<u> </u>
4.0 Are openings / cracks sealed where the slab meets the foundation wall (if appropriate)?	<u>X</u>	<u> </u>	<u> </u>
5.0 Is there adequate access to service the fan and other electrical services?	<u>X</u>	<u> </u>	<u> </u>
6.0 Is there a sump pit in the basement?	<u>X</u>	<u> </u>	<u> </u>

	Yes	No	NA
6.1 Is the sump pit installed with a sealed impermeable cover?	<u> x </u>	<u> </u>	<u> </u>
6.2 Are the penetrations through the cover sealed?	<u> </u>	<u> </u>	<u> x </u>
6.3 Does the cover have a clear view port to permit observations of conditions in the sump pit?	<u> </u>	<u> x </u>	<u> </u>
7.0 Does each suction point have a mechanism to measure vacuum?	<u> x </u>	<u> </u>	<u> </u>
7.1 Is the pressure reading from the latest commissioning clearly marked on the vent pipe?	<u> </u>	<u> x </u>	<u> </u>
7.2 Are the current diagnostic measurements within a 20% difference as compared to the system commissioning (baseline) values?	<u> </u>	<u> </u>	<u> x </u>
8.0 Does the mitigation system include an operational audible alarm to inform occupants of a system malfunction?	<u> </u>	<u> x </u>	<u> </u>
9.0 Is the exterior PVC piping in good condition with no breaks, cracks or leaks?	<u> x </u>	<u> </u>	<u> </u>

10.0 Is the fan functioning properly?

Yes	No	NA
<u> x </u>	<u> </u>	<u> </u>

11.0 Are photographs from the inspection attached?

<u> x </u>	<u> </u>	<u> </u>
--------------	---------------	---------------

12.0 Are other documents such as copies of invoices for repair work, receipts for replacement equipment, etc. attached?

<u> </u>	<u> </u>	<u> x </u>
---------------	---------------	--------------

13.0 Were any maintenance activities required?

<u> </u>	<u> x </u>	<u> </u>
---------------	--------------	---------------

14.0 Are site records up to date?

<u> x </u>	<u> </u>	<u> </u>
--------------	---------------	---------------

15.0 Has site usage changed since the previous site inspection?

<u> </u>	<u> x </u>	<u> </u>
---------------	--------------	---------------

16.0 Additional Notes & Comments

The tenant space remains vacant.

17.0 Site Sketch (if applicable)

See attached.

APPENDIX 7

Photographic Log



Photo 1: View of SS-102 installed in the basement of the former First Class Dry Cleaners.



Photo 2: View of SS-103 installed in the basement of the former Tutor Time.



Photo 3: View of Carmen's Laundromat.



Photo 4: View of the former Tutor Time .

APPENDIX 8

Laboratory Analytical Report (E18-06141)

EPA TO-15 DATA PACKAGE

ANALYTICAL DATA PACKAGE FOR THE
NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION
ALBANY NEW YORK 12233

Integrated Analytical Laboratories, LLC
Project#: Congers CP / 060141
SDG #: E18-06141
Date of first sample receipt: 8/2/2018

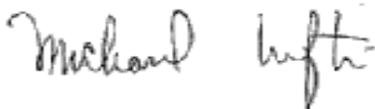
Randolph, NJ 07869
NY ELAP Certification#: 11402
NJDEP (Primary AB) Certification#: 14751
Date of last sample receipt: 8/2/2018

Client: Brennan Environmental
Project/Site: Congers CP / 060141/NY

Client Sample Number	Laboratory Sample	Sample Location	Date/Time of Collection
SS-101	E18-06141-07	NA	7/31/2018 9:17
SS-102	E18-06141-08	NA	7/31/2018 8:33
SS-103	E18-06141-09	NA	7/31/2018 8:38

This report shall not be reproduced, except in its entirety, without the written consent of Integrated Analytical Laboratories, LLC. The test results included in this report relate only to the samples analyzed. The results reported herein conform to the most current TNI standards and the laboratory's Quality Assurance Manual, where applicable, unless otherwise noted in the body of the report.

I certify that this data package is in compliance with the terms and conditions of this contract, both technically and for completeness, for other than the conditions detailed above. Release of data contained in this hardcopy data package and in the computer-readable data submitted on CD/diskette and by electronic mail has been authorized by the laboratory manager or his designee, as verified by the following signature.



Michael H. Leftin, Ph.D.
Laboratory Director

Date: August 27, 2018



Lauren Jenkins
Air Division Quality Assurance Officer

Date: August 27, 2018

EPA Method TO-15 Table of Contents

Laboratory Acronyms.....	1
Section I: Chain of Custody.....	2
Section II: Methodology Review.....	8
Section III: Case Narrative.....	10
Section IV: Method Detection Limit Summary.....	17
Section V: Quality Control Data Summary.....	29
BFB Tune Summary.....	30
Method Blank.....	35
Laboratory Sample Duplicate.....	41
Internal Standard Area Summary.....	47
Section VI: Sample Data Summary.....	52
Certificate of Analysis.....	53
Results for Sample E18-06141-07.....	54
Results for Sample E18-06141-08.....	60
Results for Sample E18-06141-09.....	74
Section VII: Standards Data.....	85
Initial Calibration Data.....	86
Initial Calibration Verification Data.....	140
Continuing Calibration Data.....	150
Section VIII: Raw Quality Control Data Package.....	171
BFB Tune Spectra.....	172
Method Blank.....	182
Laboratory Sample Duplicate.....	194
Instrument Run Logs.....	244
Pressure Gauge Readings (initial and final).....	249
Example Calculations.....	251
Clean Canister Certification.....	252
LAST PAGE OF DOCUMENT.....	256

Laboratory Acronyms

The following is a list of laboratory acronyms commonly used in EPA Method TO-15 testing:

Acronym	Definition
BLK	Blank/Method Blank
BFB	4-Bromofluorobenzene (Tuning Standard)
CAS Number	Chemical Abstract Service Registry Number
cc	cubic centimeters
CCCVS	Closing Calibration Check Verification Standard
COC	Chain of Custody
DCVS	Daily Calibration Verification Standard
DF	Dilution Factor
EPA	U. S. Environmental Protection Agency (aka USEPA)
"Hg	Inches of Mercury
IA	Indoor Air
IASL	Indoor Air Screening Level
ICAL	Initial Calibration
ICVSS	Initial Calibration Verification Standard
ISTD	Internal Standard
GC	Gas Chromatography
GC/MS	Gas Chromatography/Mass Spectrometry
LCS	Laboratory Control Sample/Spike
LLTO-15	Low Level TO-15
MDL	Method Detection Limit
MDLV	Method Detection Limit Verification
ml	milliliters
ND	Not Detected (at or above RL)
NJDEP	New Jersey Department of Environmental Protection
PM	Project Manager
ppbv	parts per billion, volume-to-volume ratio
PQL	Practical Quantitation Limit - MDLx3
QA	Quality Assurance
QC	Quality Control
RAL	Rapid Action Limit
RL	Reporting Limit
RLLCS	Reporting Limit Laboratory Control Sample
RPD	Relative Percent Difference
RRF	Relative Response Factor
RSD	Relative Standard Deviation
SDG	Sample Delivery Group
SGSL	Soil Gas Screening Levels
SS	Sub Slab
TAT	Turnaround Time
TIC	Tentatively Identified Compound
µg/m ³	micrograms per cubic meter

Section I: Chain of Custody

Client Contact Information				Project Information				Carrier (check one): <input checked="" type="checkbox"/> IAL Courier <input type="checkbox"/> Client Courier <input type="checkbox"/> FedEx/UPS				pg <u>1</u> of <u>2</u>													
Company: <u>BEI</u>				Project Name: <u>Congers C.P. / 060141</u>				Invoice Information				Analysis		Report		Matrix									
Address: <u>19 Chatham Rd. Summit, NJ 07901</u>				Project Location (State): <u>NY</u>				Attn:				EPA TO-15 NJDEP LLTO-15 (includes 30 TICs) Library Search (10, 20, or 30 TICs) Other (Explain in Comments) Regulatory/ NY Cat B / Full (NJ Required) Reduced / NY Cat A Data Package Results Only Indoor Air Ambient / Outdoor Air Sub Slab / Soil Gas / Near Slab (Circle One) Stack Emission / SVE System High Concentrations Expected													
Phone: <u>908-918-1702</u>				Project Manager: <u>Jeff McCurdy</u>				Address:																	
Fax: <u>908-918-1707</u>				PM Signature: <u>[Signature]</u>				PO #:																	
Report to:				PM E-Mail: <u>Jmccurdy@bei-enr.com</u>				Quote #:																	
				Sampler: <u>AS</u>																					
Analysis Turnaround Time - IF NO TAT IS SPECIFIED, 2 WEEK TAT IS ASSUMED								Barometric Pressure																	
IAL Standard: 2 weeks (10 business days)								Start: <u>30.10</u> Stop: <u>30.10</u>																	
Rush (**pre-approved by lab):								Flow Regulator ID																	
		24hr**		48hr**		72hr**		96hr**		1wk**															
Sample Identification	Start DATE & TIME (24hr Clock)	End DATE & TIME (24hr Clock)	Starting Vacuum ("Hg)	Ending Vacuum ("Hg)	Starting Temp. (°F)	Ending Temp. (°F)	Outgoing Vacuum - Lab ("Hg)	Incoming Vacuum - Lab ("Hg)	Flow Regulator ID	Canister ID	Canister Size (1L or 6L)	Flow Controller Readout (cc/min)	EPA TO-15	NJDEP LLTO-15 (includes 30 TICs)	Library Search (10, 20, or 30 TICs)	Other (Explain in Comments)	Regulatory/ NY Cat B / Full (NJ Required)	Reduced / NY Cat A Data Package	Results Only	Indoor Air	Ambient / Outdoor Air	Sub Slab / Soil Gas / Near Slab (Circle One)	Stack Emission / SVE System	High Concentrations Expected	
1) AA-101	7/31/18 935	7/31/18 408	-30	-2.0	72	92	-29.0	-2.0	7342673	5076	6L	12.4	H				X					X			
2) IA-FF-101	837	411	-29.5	-2.5	69	69	-29.0	-2.5A	0702495-3	3013	6L	12.3	H				X				X				
3) IA-FF-102	839	355	-28.5	-2.0	68	69	-29.0	-2.0	7342708	3814	6L	12.4	H				X				X				
4) IA-FF-103	844	412	-30	-2.5	68	69	-29.0	-2.5A	0302495-5	2753	6L	12.5	H				X				X				
5) IA-B-101	840	348	-29	-2.5	69	68	-29.0	-2.5A	0121687-1	3277	6L	12.5	H				X				X				
6) IA-B-102	842	410	-29.5	-2.5	68	68	-29.0	-2.5	7340286	3026A	6L	12.6	H				X				X				
Comments/ Special Analysis Instructions / QC Requirements:													Note: Hold or contingent samples may be designated by writing an "H" or "C" in the appropriate analysis box. ALL FIELDS IN RED ARE REQUIRED												
3013, 2753, 3026A 06081801 06131822 3059 H = hold X = run Report VOCs only 5076, 3814, 3277 05221801 05291822 5078																									
Shipping Information / Canister Preparation (for laboratory use only)										Laboratory Canister Certification															
Individual Preparing Canisters / Title: P. Jenkins, J. Walukiewicz / Air Department Sample Custodians										GC/MS Analyst Signature															
Lab Affixed Seal Number(s): <u>IAL-20180200, -0201, -0202</u>										[Signature] IAL SDG#: <u>06141</u>															
Date/Time Shipping Container Sealed: <u>7-26-18 1130</u>																									
External Chain of Custody																									
Relinquished				Received				Date / Time				Reason for Change of External Custody													
[Signature]				[Signature]				7-26-18 1130				shipment from laboratory to client													
[Signature]				[Signature]				8/21/18 1606				[Signature]													
[Signature]				[Signature]				8/2/18 1659				[Signature]													
Name/Title Resealing Shipping Container Name:										NJDEP Affixed Seal Number:															
Date/Time Sample Shipping Container Resealed:										Individual Opening Sample Shipping Container: Padraic Jenkins / Joseph Walukiewicz															
Date/Time Sample Shipping Container Opened:										Date/Time Internal Chain of Custody Initiated:															

Client Contact Information				Project Information				Carrier (check one): <input checked="" type="checkbox"/> IAL Courier <input type="checkbox"/> Client Courier <input type="checkbox"/> FedEx/UPS				pg <u>2</u> of <u>2</u>															
Company: <u>BEI</u>				Project Name: <u>Congers C.P./066141</u>				Invoice Information				Analysis		Report		Matrix											
Address: <u>19 Chatham Rd. Summit, NJ 07901</u>				Project Location (State): <u>NY</u>				Attn:				EPA TO - 15 NJDEP LLTO-15 (Includes 30 TICs) Library Search (10, 20, or 30 TICs) Other (Explain in Comments) Regulatory/ NY Cat B / Full (NJ Required) Reduced / NY Cat A Data Package Results Only Indoor Air Ambient / Outdoor Air Sub Slab / Soil Gas / Near Slab (Circle One) Stack Emission / SVE System High Concentrations Expected															
Phone: <u>908-918-1702</u>				Project Manager: <u>Jeff McCurdy</u>				Address:																			
Fax: <u>908-918-1707</u>				PM Signature: <u>[Signature]</u>				PO #:																			
Report to:				PM E-Mail: <u>jmccurdy@be-enr.com</u>				Quote #:																			
Sampler: <u>AS</u>																											
Analysis Turnaround Time - IF NO TAT IS SPECIFIED, 2 WEEK TAT IS ASSUMED										Barometric Pressure																	
IAL Standard: 2 weeks (10 business days)										Start		30.10		Stop		30.10											
Rush (**pre-approved by lab):										24hr**		48hr**		72hr**		96hr**		1wk**									
Sample Identification	Start DATE & TIME (24hr Clock)	End DATE & TIME (24hr Clock)	Starting Vacuum ("Hg)	Ending Vacuum ("Hg)	Starting Temp. (°F)	Ending Temp. (°F)	Outgoing Vacuum - Lab ("Hg)	Incoming Vacuum - Lab ("Hg)	Flow Regulator ID	Canister ID	Canister Size (1L or 6L)	Flow Controller Readout (cc/min)															
7) 55-101	8/1/18 9:17	8/1/18 4:15	-29.0	-2.5	68	68	-29.0	-2.5	A00988641-6	3014A	6L	12.7	X					X									
8) 55-102	↓ 8:33	↓ 4:19	-30.0	-2.5	68	68	-29.0	-2.5	A0302498-2	2155	6L	12.6	X					X									
9) 55-103	↓ 8:39	↓ 4:11	-30.0	-2.5	68	69	-29.0	-2.5	A0302497	5089	6L	12.5	X					X									
Comments/ Special Analysis Instructions / QC Requirements: 3014A, 2155, 5089 06091801 061318aa 3059 H=hold X=run Report CVOCS only													Note: Hold or contingent samples may be designated by writing an "H" or "C" in the appropriate analysis box. ALL FIELDS IN RED ARE REQUIRED														
Shipping Information / Canister Preparation (for laboratory use only)										Laboratory Canister Certification																	
Individual Preparing Canisters / Title: P. Jenkins, J. Walukiewicz / Air Department Sample Custodians										GC/MS Analyst Signature																	
Lab Affixed Seal Number(s): <u>IAL-20180200, -0201, -0202</u>										[Signature] IAL SDG#: <u>06141</u>																	
Date/Time Shipping Container Sealed: <u>7-26-18 11:30</u>																											
External Chain of Custody																											
Relinquished				Received				Date / Time				Reason for Change of External Custody															
[Signature]				[Signature]				7-26-18 11:30				shipment from laboratory to client															
								8/2/18 1606																			
								8/2/18 1659				Rec'd @ air lab															
Name/Title Resealing Shipping Container Name:										NJDEP Affixed Seal Number:																	
Date/Time Sample Shipping Container Resealed:										Individual Opening Sample Shipping Container: Padraic Jenkins / Joseph Walukiewicz																	
Date/Time Sample Shipping Container Opened:										Date/Time Internal Chain of Custody Initiated:																	

PROJECT INFORMATION

RUSH

E18-06141: CONGERS CP / 060141

To: Jeff McCurdy
Brennan Environmental, Inc.
Fax: 1(908) 918-1707
EMail: jmccurdy@bei-env.com

Report To

Brennan Environmental, Inc.
19 Chatham Road
Summit, NJ 07901
Attn: Jeff McCurdy

Bill To

Brennan Environmental, Inc.
19 Chatham Rd.
Summit, NJ 07901
Attn: Jeff McCurdy

Report Format	P.O. #	Received At Lab	TPHC Due	Verbal Due	Hardcopy Due
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Air Regulatory		Aug 02, 2018 @ 16:59	NA	Aug 16, 2018	Aug 30, 2018 *
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* Any *Conditional or Hold* status will delay final hardcopy report sent date.

Diskette Req. Not Required

Lab ID	Client Sample ID	Depth	Sampling Time	Matrix	Unit	Field pH/Temp
06141-001	AA-101	NA	07/31/18@16:08	Air-Ambient	ppbV	
06141-002	IA-FF-101	NA	07/31/18@16:11	Air-Indoor	ppbV	
06141-003	IA-FF-102	NA	07/31/18@15:55	Air-Indoor	ppbV	
06141-004	IA-FF-103	NA	07/31/18@16:12	Air-Indoor	ppbV	
06141-005	IA-B-101	NA	07/31/18@15:48	Air-Indoor	ppbV	
06141-006	IA-B-102	NA	07/31/18@16:10	Air-Indoor	ppbV	
06141-007	SS-101	NA	08/01/18@16:15	Air-Other	ppbV	
06141-008	SS-102	NA	08/01/18@16:19	Air-Other	ppbV	
06141-009	SS-103	NA	08/01/18@16:11	Air-Other	ppbV	

Sample #	Test	Status	QA Method	TAT	Holding Time Expires
001	TO-15 analysis	Cancel	TO-15	STD/2 WKS	8/30/2018
002	TO-15 analysis	Cancel	TO-15	STD/2 WKS	8/30/2018
003	TO-15 analysis	Cancel	TO-15	STD/2 WKS	8/30/2018
004	TO-15 analysis	Cancel	TO-15	STD/2 WKS	8/30/2018
005	TO-15 analysis	Cancel	TO-15	STD/2 WKS	8/30/2018
006	TO-15 analysis	Cancel	TO-15	STD/2 WKS	8/30/2018
007	TO-15 analysis	Analyze	TO-15	STD/2 WKS	8/31/2018
008	TO-15 analysis	Analyze	TO-15	STD/2 WKS	8/31/2018
009	TO-15 analysis	Analyze	TO-15	STD/2 WKS	8/31/2018

Project Notes:

NOTE 1 taken by kfalconer on 08/03/2018 09:24
REPORT CVOCS ONLY



PROJECT INFORMATION

RUSH

E18-06141: CONGERS CP / 060141

REV 1 taken by lauren on 08/23/2018 11:02

As per Jeff McCurdy, cancel TO-15 analysis for sample # 1,2,3,4,5,6





Integrated Analytical Laboratories Internal Chain of Custody

Instructions: Use 1 form for each 20 samples of aliquot.

Laboratory Person Accepting Responsibility for Sample(s)			
Laboratory: Integrated Analytical Laboratories	Location: 273 Franklin Rd Randolph, NJ 07869		
Name: Joseph Walukiewicz	Title: Air Department Receiving		
Case No.: E18-06141	Analytical Parameter/Fraction: (check one) <input type="checkbox"/> NJDEP LLTO-15 <input checked="" type="checkbox"/> EPA TO-15		

Sample No.	Aliquot/Extract No.
22-101	E18-06141-001
22-FF-101	E18-06141-002
22-FF-102	E18-06141-003
22-FF-103	E18-06141-004
22-B-101	E18-06141-005
22-B-102	E18-06141-006
22-101	E18-06141-007
22-102	E18-06141-008
22-103	E18-06141-009
	E18-

Sample No.	Aliquot/Extract No.
	E18-

Date	Time	Relinquished By	Received By	Purpose of Change of Custody
08 02 18	17 00	SIGNATURE	SIGNATURE <i>Joseph Walukiewicz</i>	1. Sample log-in 2. Pressure Check 3. Pre-analysis storage
		PRINTED NAME	PRINTED NAME JOSEPH WALUKIEWICZ	
08 02 18	17 15	SIGNATURE <i>Joseph Walukiewicz</i>	SIGNATURE	
		PRINTED NAME JOSEPH WALUKIEWICZ	PRINTED NAME	
08 06 18	12 30	SIGNATURE	SIGNATURE <i>Jeff Schmitt</i>	TO-15/LLTO-15 analysis on: E18-06141-01, 02, 03, 04, 05 06, 07, 08, 09
		PRINTED NAME	PRINTED NAME JEFF SCHMITT	
		SIGNATURE	SIGNATURE	
		PRINTED NAME	PRINTED NAME	
		SIGNATURE	SIGNATURE	
		PRINTED NAME	PRINTED NAME	
		SIGNATURE	SIGNATURE	
		PRINTED NAME	PRINTED NAME	
		SIGNATURE	SIGNATURE	
		PRINTED NAME	PRINTED NAME	

Section II: Methodology Review

Methodology Summary for Air Collected from Hazardous Waste Site Contract

Laboratory:	Integrated Analytical Lab, LLC	Project No:	Congers CP / 060141
Location:	Randolph, NJ	SDG No:	E18-06141

Name	Required Methodology	Indicate Method
Volatile Organics	US EPA TO-15	US EPA Method TO-15

Section III: Case Narrative

CASE NARRATIVE

ANALYTICAL DATA PACKAGE FOR THE NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION ALBANY NEW YORK 12233

Integrated Analytical Laboratories, LLC
 Project #: Congers CP / 060141
 SDG #: E18-06141
 Date of first sample receipt: 8/2/2018

Randolph, NJ 07869
 NY ELAP Certification #: 11402
 NJDEP (Primary AB) Certification#: 14751
 Date of last sample receipt: 8/2/2018

Client: Brennan Environmental
 Project/Site: Congers CP / 060141 / NY

Client ID	Lab ID	Receipt Date	Analysis Date	DF	Diluted For
AA-101	E18-06141-01	08/02/2018	Cancelled as per client request		NA
IA-FF-101	E18-06141-02	08/02/2018	Cancelled as per client request		NA
IA-FF-102	E18-06141-03	08/02/2018	Cancelled as per client request		NA
IA-FF-103	E18-06141-04	08/02/2018	Cancelled as per client request		NA
IA-B-101	E18-06141-05	08/02/2018	Cancelled as per client request		NA
IA-B-102	E18-06141-06	08/02/2018	Cancelled as per client request		NA
SS-101	E18-06141-07	08/02/2018	08/06/2018	1.0	NA
SS-102	E18-06141-08	08/02/2018	08/07/2018	100.0	1,2-Dichloroethene (cis) Trichloroethene Tetrachloroethene
SS-102	E18-06141-08	08/02/2018	08/06/2018	1.0	NA
SS-103	E18-06141-09	08/02/2018	08/07/2018	10.0	Tetrachloroethene
SS-103	E18-06141-09	08/02/2018	08/06/2018	1.0	NA

IAL Sample ID	Canister ID	Outgoing Pressure ("Hg)	Incoming Pressure ("Hg)	Flow Controller ID	Outgoing Flow Rate (cc/min)	Incoming Flow Rate (cc/min)	Flow Rate RPD*
E18-06141-01	5076	-29	-2.0	7342673	12.40	12.60	1.60
E18-06141-02	3013	-29	-2.5	A0302495-3	12.30	12.50	1.61
E18-06141-03	3814	-29	-2.0	7342708	12.40	12.40	0.00
E18-06141-04	2753	-29	-2.5	A03024955-5	12.50	12.30	1.61
E18-06141-05	3277	-29	-2.5	A0121687-1	12.50	12.70	1.59
E18-06141-06	3026A	-29	-2.5	7340286	12.60	12.60	0.00
E18-06141-07	3014A	-29	-2.5	A00988641-6	12.70	12.80	0.78
E18-06141-08	2155	-29	-2.5	A0302498-2	12.60	12.90	2.35
E18-06141-09	5089	-29	-2.5	A0302497-4	12.50	12.80	2.37

*Pre-sampling and Post-sampling Flow Controller calibration check RPD ≤ 20%

Flow Controller Note: none

CASE NARRATIVE

ANALYTICAL DATA PACKAGE FOR THE NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION ALBANY NEW YORK 12233

Integrated Analytical Laboratories, LLC
 Project #: Congers CP / 060141
 SDG #: E18-06141
 Date of first sample receipt: 8/2/2018

Randolph, NJ 07869
 NY ELAP Certification #: 11402
 NJDEP (Primary AB) Certification#: 14751
 Date of last sample receipt: 8/2/2018

Client: Brennan Environmental
 Project/Site: Congers CP / 060141 / NY

Sample Receipt: Samples were received in good condition. Documentation was in order.
 Samples were received at IAL by: Joseph Walukiewicz

Sample Preparation: None required.

Sample Analysis:

Hold Time: All within recommended hold times.
Instrument Calibration: Meets method criteria.
Analysis performed by: Jeff Schmitt
SDG Non-Conformances: none
Tentatively Identified Compounds: Tentatively Identified Compounds (TICs) are determined using a NIST library search. TICs are reported at 10% of the applicable internal standard. Dilution factors are calculated into the final reported result. Since the compounds found are tentatively identified, the conversion from ppbv to ug/m3 may not be made.

Canister-to-Canister dilutions: none

Dilutions: Dilutions, if necessary, will be conducted directly on the instrument up to a 500x dilution. When dilutions of 1000x or higher are necessary, the laboratory must inject a volume of sample into another certified clean canister and add humidified Z-1 zero air to the remainder of the canister volume. Tedlar bags are not used for dilutions.

If a sample is received with historically high levels of analytes, a 100x can-to-can dilution may be used from the start. A 100x canister-to-canister dilution may be also be used at the analyst's discretion.

On-instrument dilutions are conducted as follows:

Dilution Factor	Sample Volume Injected (cc)
1	500
2.5	200
5	100
10	50
20	25
25	20
50	10
100	5
200	2.5
250	2
500	1

CASE NARRATIVE

ANALYTICAL DATA PACKAGE FOR THE NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION ALBANY NEW YORK 12233

Integrated Analytical Laboratories, LLC
 Project #: Congers CP / 060141
 SDG #: E18-06141
 Date of first sample receipt: 8/2/2018

Randolph, NJ 07869
 NY ELAP Certification #: 11402
 NJDEP (Primary AB) Certification#: 14751
 Date of last sample receipt: 8/2/2018

Client: Brennan Environmental
 Project/Site: Congers CP / 060141 / NY

Canister-to-canister dilutions are conducted as follows:

A certified clean canister is obtained and evacuated to approximately -30"Hg. Both the clean/dilution canister and sample canister are fitted with a ¼" Swagelok® nut fitting equipped with septa. Depending on dilution factor necessary, a sample aliquot is removed from the canister and injected into the clean canister using 30cc Multifit gas-tight syringe. Once the correct sample aliquot has been transferred, the dilution canister should be connected to the humidified Z-1 zero air supply and filled to ambient pressure (0"Hg).

Dilution Factor	Sample Aliquot	Z-1 Make-up Added
100	60ml	5940ml
1000	6ml	5994ml

If further dilutions need to be made from the dilution canister, they may be made on-instrument. Using a 100x dilution canister, the following on-instrument dilutions can be produced:

Dilution Factor	Sample Volume Injected
100	500ml
250	200ml
500	100ml
1000	50ml
2000	25ml
2500	20ml
5000	10ml

Using a 1000x dilution canister, the following on-instrument dilutions can be produced:

Dilution Factor	Sample Volume Injected
1000	500ml
2500	200ml
5000	100ml
10,000	50ml
20,000	25ml
25,000	20ml
50,000	10ml

If further dilutions need to be made from the dilution canister, beyond 50,000x, a subsequent canister-to-canister dilution must be made using the above prescribed protocol.

CASE NARRATIVE

ANALYTICAL DATA PACKAGE FOR THE NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION ALBANY NEW YORK 12233

Integrated Analytical Laboratories, LLC
Project #: Congers CP / 060141
SDG #: E18-06141
Date of first sample receipt: 8/2/2018

Randolph, NJ 07869
NY ELAP Certification #: 11402
NJDEP (Primary AB) Certification#: 14751
Date of last sample receipt: 8/2/2018

Client: Brennan Environmental
Project/Site: Congers CP / 060141 / NY

GC Column and ID: RTX-1 SN 1119138 or equivalent

Calibration Standards: Only gas phase standards were used. Primary and second-source standards provided by Scott Specialty Gases or Airgas Specialty Gases/ Air Liquide

Working Standards: Primary source standards* are created from:
- Airgas Specialty Gases Cylinder #CC483586, valid 3/6/17 through 4/6/19 @ approximately 100ppb per compound, with exception of m&p-xylenes @ 200ppb. Standard is directly introduced into the instrument for all calibration standard concentrations. Dilutions are made accordingly, on instrument, with humidified clean air. The 10ppbv standard is also used for the Daily Calibration Verification Standard (DCVS) and Closing Calibration Verification Standard (CCCVS).

The second source standard*, used as the Initial Calibration Verification Standard (ICVSS), is introduced into the instrument in the same manner as the primary source standard, using:

- Airgas Specialty Gases Cylinder #CC483422, valid 5/4/17 through 4/27/19 @ approximately 100ppb per compound, with exception of m&p-xylenes @ 200ppb.

Internal standards* are created from:

- Scott Gas, Cylinder #ALM029426, valid 5/11/2016 through 5/12/2019 @ 5ppm per compound. Standard is directly introduced into the instrument to reach the 10ppbv concentrations. 1:500 Dilutions are made on instrument with humidified clean air. 1cc of internal standard is added to every standard, method blank, instrument blank, and sample run.

*Standard may be used past its expiration date provided that concentrations are verified by a current/unexpired second source standard.

05/18/2018

100 ppbv internal standard mix - prepared in cylinder #ALM029426
10 ppbv per standard/sample - 50 ml injected

100 ppbv calibration standard - prepared in cylinder #CC483586
40 ppbv standard - 200 ml injected
20 ppbv standard - 100 ml injected
10 ppbv standard* - 50 ml injected

*Standard also used for CCCVS

2 ppbv standard - 10 ml injected
0.20 ppbv standard* - 1 ml injected

*Standard also used for RLLCS

06/13/2018

100 ppbv internal standard mix - prepared in cylinder #ALM029426
10 ppbv per standard/sample - 50 ml injected
100 ppbv calibration standard - prepared in cylinder #CC483586

CASE NARRATIVE

ANALYTICAL DATA PACKAGE FOR THE NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION ALBANY NEW YORK 12233

Integrated Analytical Laboratories, LLC
Project #: Congers CP / 060141
SDG #: E18-06141
Date of first sample receipt: 8/2/2018

Randolph, NJ 07869
NY ELAP Certification #: 11402
NJDEP (Primary AB) Certification#: 14751
Date of last sample receipt: 8/2/2018

Client: Brennan Environmental
Project/Site: Congers CP / 060141 / NY

10 ppbv standard* - 50 ml injected
*Standard also used for DCVS & CCCVS

0.20 ppbv standard* - 1 ml injected
*Standard also used for RLLCS

Method Blank - prepared in canister #1127
500 ml injected

07/25/2018

100 ppbv internal standard mix - prepared in cylinder #ALM029426
10 ppbv per standard/sample - 50 ml injected
100 ppbv calibration standard - prepared in cylinder #CC483586
40 ppbv standard - 200 ml injected
20 ppbv standard - 100 ml injected
10 ppbv standard* - 50 ml injected
*Standard also used for CCCVS
2 ppbv standard - 10 ml injected
0.20 ppbv standard* - 1 ml injected
*Standard also used for RLLCS

08/06/2018

100 ppbv internal standard mix - prepared in cylinder #ALM029426
10 ppbv per standard/sample - 50 ml injected
100 ppbv calibration standard - prepared in cylinder #CC483586
10 ppbv standard* - 50 ml injected
*Standard also used for DCVS & CCCVS
0.20 ppbv standard* - 1 ml injected
*Standard also used for RLLCS
Method Blank - prepared in canister #1127
500 ml injected
Sample E18-06141-07 - sample taken in canister #3014A
500 ml sample volume injected, 1x dilution
Sample E18-06141-08 - sample taken in canister #2155
500 ml sample volume injected, 1x dilution
Sample E18-06141-09 - sample taken in canister #5089
500 ml sample volume injected, 1x dilution

08/07/2018

100 ppbv internal standard mix - prepared in cylinder #ALM029426
10 ppbv per standard/sample - 50 ml injected
100 ppbv calibration standard - prepared in cylinder #CC483586
10 ppbv standard* - 50 ml injected
*Standard also used for DCVS & CCCVS

CASE NARRATIVE

**ANALYTICAL DATA PACKAGE FOR THE
NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION
ALBANY NEW YORK 12233**

Integrated Analytical Laboratories, LLC
Project #: Congers CP / 060141
SDG #: E18-06141
Date of first sample receipt: 8/2/2018

Randolph, NJ 07869
NY ELAP Certification #: 11402
NJDEP (Primary AB) Certification#: 14751
Date of last sample receipt: 8/2/2018

Client: Brennan Environmental
Project/Site: Congers CP / 060141 / NY

0.20 ppbv standard* - 1 ml injected
*Standard also used for RLLCS
Method Blank - prepared in canister #1127
500 ml injected
Sample E18-06141-08 - sample taken in canister #2155
5 ml sample volume injected, 100x dilution
Sample E18-06141-09 - sample taken in canister #5089
50 ml sample volume injected, 10x dilution

All work recorded herein has been done in accordance with normal professional standards using accepted testing methodologies, quality assurance and quality control procedures except where otherwise agreed to by the client and testing company in writing. All conversions are based upon a room temperature of 77°F(25°C) and room pressure of 101.325 kPa (1atm).

I certify that this data package is in compliance with the terms and conditions of this contract, both technically and for completeness, for other than the conditions detailed above. Release of data contained in this hardcopy data package and in the computer-readable data submitted on CD/diskette and by electronic mail has been authorized by the laboratory manager or his designee, as verified by the following signature.



Michael H. Leftin, Ph.D.
Laboratory Director

August 27, 2018
Date

Section IV: Method Detection Limit Summary

METHOD DETECTION LIMIT (MDL) REPORT

Integrated Analytical Laboratories - Randolph, NJ

Matrix: Air

Column ID: Restek Rtx-1, 60 meter, 0.32mm ID, 1 um

Instrument ID: GC - Agilent 7890A / MS - Agilent 5975C (IAL ID: Instrument AA)

Report Prepared by: Jeff Schmitt

Preparation Date: 6/27/17

MDL Effective Date: 6/27/2017

MDL Expiration Date: 6/27/2018

MDL Analysis Date: 6/27/2017

Analyst: Jeff Schmitt

Filename	Run #	Date	Time	Filename	Run #	Date	Time
aa2539	Run 1	6/27/2017	11:54	aa2544	Run 6	6/27/2017	14:44
aa2540	Run 2	6/27/2017	12:28	aa2545	Run 7	6/27/2017	15:18
aa2541	Run 3	6/27/2017	13:02				
aa2542	Run 4	6/27/2017	13:36				
aa2543	Run 5	6/27/2017	14:10				

Compound Name	Run 1*	Run 2*	Run 3*	Run 4*	Run 5*	Run 6*	Run 7*	MEAN Value	TRUE Value	Percent Recovery	Std Dev Conc	MDL ppbv	PQL ppbv	NY/LLTO-15/PA RL ppbv	NY/LLTO-15/PA RL µg/m³	NJ TO-15 RL ppbv	NJ TO-15 RL µg/m³	True value/MDL
Propene	0.18	0.15	0.10	0.08	0.15	0.05	0.13	0.12	0.20	61	0.045	0.14	0.43	0.20	0.34	0.40	0.69	1.4
Dichlorodifluoromethane	0.18	0.20	0.18	0.16	0.21	0.21	0.18	0.19	0.20	94	0.016	0.050	0.15	0.20	0.99	0.40	2.0	4.0
Chloromethane	0.08	0.17	0.12	0.08	0.13	0.10	0.17	0.12	0.20	61	0.038	0.12	0.36	0.20	0.41	0.40	0.83	1.7
1,2-Dichlorotetrafluoroethane	0.31	0.32	0.31	0.28	0.34	0.34	0.32	0.32	0.20	159	0.019	0.059	0.18	0.20	1.4	0.40	2.8	3.4
Vinyl chloride	0.12	0.14	0.13	0.11	0.15	0.14	0.12	0.13	0.20	65	0.015	0.046	0.14	0.20	0.51	0.40	1.0	4.3
1,3-Butadiene	0.18	0.21	0.19	0.15	0.21	0.20	0.18	0.19	0.20	96	0.022	0.069	0.21	0.20	0.44	0.40	0.88	2.9
n-Butane	0.28	0.29	0.27	0.24	0.30	0.29	0.26	0.28	0.20	139	0.022	0.068	0.20	0.20	0.47	0.40	0.95	3.0
Bromomethane	0.12	0.12	0.12	0.09	0.12	0.06	0.11	0.11	0.20	53	0.023	0.073	0.22	0.20	0.78	0.40	1.6	2.7
Chloroethane	0.08	0.16	0.08	0.16	0.24	0.20	0.17	0.15	0.20	77	0.058	0.18	0.54	0.20	0.53	0.40	1.1	1.1
Ethanol	0.24	0.11	0.17	0.20	0.27	0.29	0.22	0.21	0.20	107	0.061	0.19	0.57	0.20	0.38	0.40	0.75	1.0
Vinyl bromide	0.14	0.14	0.14	0.12	0.16	0.17	0.14	0.14	0.20	72	0.015	0.047	0.14	0.20	0.87	0.40	1.7	4.3
Acrolein	0.20	0.24	0.21	0.19	0.10	0.22	0.20	0.19	0.20	97	0.043	0.14	0.41	0.20	0.46	0.40	0.92	1.5
Acetone	0.26	0.26	0.25	0.23	0.28	0.19	0.24	0.24	0.20	122	0.028	0.088	0.26	0.20	0.48	0.40	0.95	2.3
Trichlorofluoromethane	0.33	0.35	0.35	0.32	0.38	0.38	0.35	0.35	0.20	176	0.023	0.072	0.22	0.20	1.1	0.40	2.2	3
Isopropanol	0.34	0.35	0.33	0.30	0.40	0.34	0.32	0.34	0.20	170	0.032	0.10	0.30	0.20	0.49	0.40	0.98	2.0
n-Pentane	0.31	0.33	0.32	0.28	0.31	0.32	0.30	0.31	0.20	155	0.015	0.046	0.14	0.20	0.59	0.40	1.2	4.3
1,1-Dichloroethene	0.21	0.23	0.22	0.20	0.25	0.25	0.22	0.23	0.20	113	0.017	0.054	0.16	0.20	0.79	0.40	1.6	3.7
Methylene chloride	0.27	0.27	0.26	0.25	0.29	0.29	0.28	0.27	0.20	137	0.014	0.045	0.13	0.20	0.69	0.40	1.4	4.5
Tert-butyl alcohol	0.36	0.36	0.35	0.30	0.34	0.38	0.31	0.34	0.20	172	0.030	0.093	0.28	0.20	0.61	0.40	1.2	2.2
Allyl Chloride	0.22	0.22	0.21	0.18	0.22	0.23	0.19	0.21	0.20	105	0.017	0.054	0.16	0.20	0.63	0.40	1.3	3.7
1,1,2-Trichloro-1,2,2-trifluoroethane	0.34	0.36	0.35	0.31	0.37	0.38	0.35	0.35	0.20	175	0.022	0.068	0.20	0.20	1.5	0.40	3.1	2.9
Carbon disulfide	0.17	0.19	0.18	0.15	0.20	0.19	0.17	0.18	0.20	89	0.017	0.052	0.16	0.20	0.62	0.40	1.2	3.8
1,2-Dichloroethene (trans)	0.21	0.24	0.22	0.20	0.24	0.24	0.21	0.22	0.20	112	0.018	0.057	0.17	0.20	0.79	0.40	1.6	3.5
1,1-Dichloroethane	0.29	0.32	0.29	0.28	0.33	0.34	0.30	0.31	0.20	155	0.022	0.069	0.21	0.20	0.81	0.40	1.6	2.9
Methyl tert-butyl ether	0.26	0.28	0.25	0.22	0.27	0.27	0.24	0.26	0.20	128	0.020	0.064	0.19	0.20	0.72	0.40	1.4	3.1
Methyl ethyl ketone	0.35	0.37	0.34	0.31	0.38	0.37	0.34	0.35	0.20	176	0.024	0.077	0.23	0.20	0.59	0.40	1.2	2.6
1,2-Dichloroethene (cis)	0.25	0.27	0.25	0.22	0.27	0.27	0.24	0.25	0.20	127	0.019	0.058	0.17	0.20	0.79	0.40	1.6	3.4
Ethyl acetate	0.46	0.48	0.49	0.44	0.51	0.54	0.45	0.48	0.20	241	0.035	0.11	0.33	0.20	0.72	0.40	1.4	1.8
n-Hexane	0.33	0.35	0.34	0.31	0.38	0.35	0.32	0.34	0.20	170	0.022	0.069	0.21	0.20	0.70	0.40	1.4	2.9
Chloroform	0.29	0.33	0.31	0.29	0.34	0.34	0.31	0.31	0.20	157	0.022	0.068	0.20	0.20	0.98	0.40	2.0	2.9
Tetrahydrofuran	0.39	0.41	0.41	0.35	0.43	0.43	0.38	0.40	0.20	200	0.026	0.082	0.24	0.20	0.59	0.40	1.2	2.5
1,2-Dichloroethane	0.32	0.34	0.33	0.30	0.37	0.37	0.34	0.34	0.20	170	0.026	0.082	0.25	0.20	0.81	0.40	1.6	2.4
1,1,1-Trichloroethane	0.31	0.33	0.31	0.28	0.34	0.34	0.31	0.32	0.20	158	0.020	0.064	0.19	0.20	1.1	0.40	2.2	3.1
Benzene	0.33	0.35	0.33	0.30	0.35	0.36	0.32	0.33	0.20	167	0.021	0.065	0.20	0.20	0.64	0.40	1.3	3
Carbon tetrachloride	0.25	0.27	0.24	0.22	0.26	0.27	0.25	0.25	0.20	126	0.016	0.050	0.15	0.20	1.3	0.40	2.5	4.0
Cyclohexane	0.27	0.29	0.27	0.24	0.29	0.28	0.25	0.27	0.20	134	0.019	0.061	0.18	0.20	0.69	0.40	1.4	3.3
1,2-Dichloropropane	0.40	0.41	0.41	0.38	0.44	0.44	0.40	0.41	0.20	206	0.023	0.071	0.21	0.20	0.92	0.40	1.8	2.8
Bromodichloromethane	0.32	0.34	0.33	0.30	0.36	0.35	0.32	0.33	0.20	165	0.021	0.066	0.20	0.20	1.3	0.40	2.7	3.0
2,2,4-Trimethylpentane	0.22	0.25	0.24	0.20	0.25	0.26	0.22	0.23	0.20	117	0.021	0.065	0.20	0.20	0.93	0.40	1.9	3.1
Trichloroethene	0.36	0.38	0.37	0.32	0.39	0.39	0.35	0.37	0.20	183	0.026	0.080	0.24	0.20	1.1	0.40	2.1	2.5
1,4-Dioxane	0.39	0.41	0.38	0.35	0.41	0.41	0.36	0.39	0.20	194	0.024	0.076	0.23	0.20	0.72	0.40	1.4	2.6
Methyl methacrylate	0.39	0.42	0.38	0.37	0.42	0.41	0.38	0.40	0.20	198	0.023	0.071	0.21	0.20	0.82	0.40	1.6	2.8
n-Heptane	0.35	0.37	0.35	0.31	0.37	0.36	0.34	0.35	0.20	175	0.022	0.068	0.20	0.20	0.82	0.40	1.6	3.0
1,3-Dichloropropene (cis)	0.33	0.35	0.33	0.31	0.36	0.36	0.32	0.34	0.20	169	0.020	0.063	0.19	0.20	0.91	0.40	1.8	3.2
Methyl isobutyl ketone	0.44	0.48	0.45	0.42	0.51	0.50	0.44	0.46	0.20	232	0.034	0.11	0.32	0.20	0.82	0.40	1.6	1.9
1,3-Dichloropropene (trans)	0.37	0.39	0.36	0.34	0.40	0.39	0.35	0.37	0.20	186	0.020	0.063	0.19	0.20	0.91	0.40	1.8	3.2
1,1,2-Trichloroethane	0.40	0.42	0.39	0.36	0.42	0.45	0.41	0.41	0.20	203	0.027	0.085	0.26	0.20	1.1	0.40	2.2	2.4

METHOD DETECTION LIMIT (MDL) REPORT

Integrated Analytical Laboratories - Randolph, NJ

Matrix: Air
 Column ID: Restek Rtx-1, 60 meter, 0.32mm ID, 1 um
 Instrument ID: GC - Agilent 7890A / MS - Agilent 5975C (IAL ID: Instrument AA)
 Report Prepared by: Jeff Schmitt
 Preparation Date: 6/27/17

MDL Effective Date: 6/27/2017
 MDL Expiration Date: 6/27/2018
 MDL Analysis Date: 6/27/2017
 Analyst: Jeff Schmitt

Filename	Run #	Date	Time	Filename	Run #	Date	Time
aa2539	Run 1	6/27/2017	11:54	aa2544	Run 6	6/27/2017	14:44
aa2540	Run 2	6/27/2017	12:28	aa2545	Run 7	6/27/2017	15:18
aa2541	Run 3	6/27/2017	13:02				
aa2542	Run 4	6/27/2017	13:36				
aa2543	Run 5	6/27/2017	14:10				

Compound Name	Run 1*	Run 2*	Run 3*	Run 4*	Run 5*	Run 6*	Run 7*	MEAN Value	TRUE Value	Percent Recovery	Std Dev Conc	MDL ppbv	PQL ppbv	NY/LLTO-	NY/LLTO-	NJ TO-15	NJ TO-15	True value/ MDL
														15/ PA RL ppbv	15/ PA RL µg/m³	RL ppbv	RL µg/m³	
Toluene	0.43	0.45	0.43	0.38	0.45	0.44	0.42	0.43	0.20	214	0.025	0.077	0.23	0.20	0.75	0.40	1.5	2.6
Methyl n-butyl ketone	0.48	0.50	0.47	0.43	0.51	0.52	0.46	0.48	0.20	240	0.031	0.10	0.30	0.20	0.82	0.40	1.6	2.0
Dibromochloromethane	0.27	0.28	0.27	0.24	0.29	0.29	0.27	0.27	0.20	136	0.016	0.051	0.15	0.20	1.7	0.40	3.4	4
1,2-Dibromoethane	0.38	0.40	0.38	0.34	0.41	0.41	0.36	0.38	0.20	192	0.025	0.078	0.23	0.20	1.5	0.40	3.1	2.6
Tetrachloroethene	0.41	0.43	0.43	0.39	0.46	0.46	0.41	0.43	0.20	213	0.025	0.078	0.24	0.20	1.4	0.40	2.7	2.5
Chlorobenzene	0.39	0.41	0.40	0.36	0.43	0.43	0.39	0.40	0.20	200	0.026	0.080	0.24	0.20	0.92	0.40	1.8	2.5
Ethylbenzene	0.36	0.38	0.35	0.32	0.38	0.37	0.34	0.36	0.20	179	0.023	0.071	0.21	0.20	0.87	0.40	1.7	3
Xylenes (m&p)	0.75	0.76	0.72	0.67	0.77	0.76	0.68	0.73	0.40	182	0.040	0.13	0.38	0.40	1.7	0.40	1.7	1.6
Bromoform	0.22	0.23	0.22	0.20	0.24	0.24	0.22	0.23	0.20	113	0.014	0.045	0.13	0.20	2.1	0.40	4.1	4
Styrene	0.33	0.34	0.32	0.28	0.35	0.33	0.30	0.32	0.20	161	0.022	0.069	0.21	0.20	0.85	0.40	1.7	3
1,1,2,2-Tetrachloroethane	0.37	0.38	0.37	0.32	0.38	0.38	0.34	0.36	0.20	182	0.023	0.073	0.22	0.20	1.4	0.40	2.7	2.8
Xylene (o)	0.40	0.41	0.39	0.35	0.43	0.43	0.38	0.40	0.20	199	0.026	0.083	0.25	0.20	0.87	0.40	1.7	2.4
n-Nonane	0.34	0.35	0.32	0.30	0.36	0.36	0.31	0.33	0.20	167	0.023	0.072	0.22	0.20	1.4	0.40	2.1	3
Cumene	0.34	0.36	0.34	0.31	0.37	0.37	0.33	0.35	0.20	173	0.024	0.075	0.23	0.20	1.0	0.40	2.0	3
2-Chlorotoluene	0.36	0.36	0.34	0.32	0.37	0.37	0.34	0.35	0.20	175	0.019	0.059	0.18	0.20	0.98	0.40	2.1	3.4
n-Propyl benzene	0.35	0.37	0.34	0.31	0.37	0.36	0.33	0.35	0.20	173	0.023	0.072	0.22	0.20	1.0	0.40	2.0	2.8
4-Ethyltoluene	0.32	0.34	0.32	0.29	0.34	0.34	0.30	0.32	0.20	161	0.020	0.062	0.19	0.20	0.98	0.40	2.0	3.2
1,3,5-Trimethylbenzene	0.32	0.34	0.32	0.28	0.34	0.34	0.30	0.32	0.20	160	0.022	0.069	0.21	0.20	0.98	0.40	2.0	2.9
1,2,4-Trimethylbenzene	0.31	0.32	0.30	0.27	0.32	0.31	0.28	0.30	0.20	151	0.022	0.069	0.21	0.20	0.98	0.40	2.0	3
Benzyl chloride	0.27	0.27	0.25	0.23	0.27	0.27	0.24	0.26	0.20	129	0.018	0.056	0.17	0.20	1.0	0.40	2.1	3.6
1,3-Dichlorobenzene	0.38	0.39	0.37	0.34	0.40	0.40	0.36	0.38	0.20	189	0.023	0.072	0.22	0.20	1.2	0.40	2.4	2.8
1,4-Dichlorobenzene	0.36	0.38	0.35	0.32	0.38	0.38	0.34	0.36	0.20	179	0.024	0.076	0.23	0.20	1.2	0.40	2.4	2.6
1,2-Dichlorobenzene	0.36	0.37	0.36	0.33	0.38	0.38	0.34	0.36	0.20	179	0.020	0.063	0.19	0.20	1.2	0.40	2.4	3.2
1,2,4-Trichlorobenzene	0.41	0.41	0.39	0.35	0.41	0.41	0.38	0.40	0.20	198	0.024	0.077	0.23	0.20	1.5	0.40	3.0	3
Naphthalene	0.41	0.41	0.39	0.36	0.41	0.40	0.35	0.39	0.20	194	0.025	0.079	0.24	0.20	1.0	0.40	2.1	2.5
1,3-Hexachlorobutadiene	0.40	0.42	0.40	0.38	0.43	0.45	0.40	0.41	0.20	206	0.024	0.075	0.22	0.20	2.1	0.40	4.3	2.7

Processing Method: C:\MSDCHEM\1\METHODS\AA0612.M
 Initial Calibration: C:\MSDCHEM\1\METHODS\AA0612.M
 Location of this file: P:\PAL Reports\2015\LLTO-15 and TO-15 Common Files\Agilent MDL

Michael Leftin
 Michael Leftin, Ph.D.
 Laboratory Director

Lauren Jenkins
 Lauren Jenkins
 Air Division Quality Assurance Officer

Date: July 5, 2017

Date: July 5, 2017

Instrument used for Clean Canister Certification Analysis? YES

METHOD DETECTION LIMIT VERIFICATION (MDLV) REPORT

Integrated Analytical Laboratories - Randolph, NJ

Analysis Level: 0.20 ppbv, 0.40 for m&p-xylenes
 Matrix: Air
 Column ID: Restek Rtx-1, 60 meter, 0.32mm ID, 1 um
 Instrument Identification: AA
 Date of Verification Study: 6/27/2017
 Study Identification File #: aa2546
 Analyst: Jeff Schmitt
 Analysis/Processing Method: C:\MSDCHEM1\METHODS\AA0612.M
 Cleanup Method: Not Applicable

Lauren Jenkins
 Lauren Jenkins
 Air Division QA Officer

Date: 6/28/17

Compound Name	CAS #	MDLV Source	Source Study (File ID)	Source Instrument	Source Analysis Date	MDLV (ppbv)	RL (ppbv)	RL/MDLV Ratio
Propene	115-07-1	AAL071156	aa4155rlcs	AA	6/22/15	0.09	0.20	2.22
Dichlorodifluoromethane	124-48-1	AAL071156	aa4155rlcs	AA	6/22/15	0.2	0.20	1.00
Chloromethane	74-87-3	AAL071156	aa4155rlcs	AA	6/22/15	0.18	0.20	1.1
1,2-Dichlorotetrafluoroethane	76-14-2	AAL071156	aa4155rlcs	AA	6/22/15	0.34	0.20	0.59
Vinyl chloride	75-01-4	AAL071156	aa4155rlcs	AA	6/22/15	0.14	0.20	1.4
1,3-Butadiene	106-99-0	AAL071156	aa4155rlcs	AA	6/22/15	0.2	0.20	1.0
n-Butane	106-97-8	AAL071156	aa4155rlcs	AA	6/22/15	0.29	0.20	0.7
Bromomethane	74-83-9	AAL071156	aa4155rlcs	AA	6/22/15	0.14	0.20	1.43
Chloroethane	75-00-3	AAL071156	aa4155rlcs	AA	6/22/15	0.13	0.20	1.5
Ethanol	64-17-5	AAL071156	aa4155rlcs	AA	6/22/15	0.29	0.20	0.7
Vinyl bromide	593-60-2	AAL071156	aa4155rlcs	AA	6/22/15	0.16	0.20	1.3
Acrolein	107-02-8	AAL071156	aa4155rlcs	AA	6/22/15	0.19	0.20	1.1
Acetone	67-64-1	AAL071156	aa4155rlcs	AA	6/22/15	0.31	0.20	0.65
Trichlorofluoromethane	75-69-4	AAL071156	aa4155rlcs	AA	6/22/15	0.37	0.20	0.54
Isopropanol	67-63-0	AAL071156	aa4155rlcs	AA	6/22/15	0.34	0.20	0.6
n-Pentane	109-66-0	AAL071156	aa4155rlcs	AA	6/22/15	0.33	0.20	0.6
1,1-Dichloroethene	75-35-4	AAL071156	aa4155rlcs	AA	6/22/15	0.24	0.20	0.8
Methylene chloride	75-09-2	AAL071156	aa4155rlcs	AA	6/22/15	0.28	0.20	0.71
Tert-butyl alcohol	75-65-0	AAL071156	aa4155rlcs	AA	6/22/15	0.36	0.20	0.6
Allyl Chloride	107-05-1	AAL071156	aa4155rlcs	AA	6/22/15	0.21	0.20	1.0
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	AAL071156	aa4155rlcs	AA	6/22/15	0.37	0.20	0.54
Carbon disulfide	75-15-0	AAL071156	aa4155rlcs	AA	6/22/15	0.19	0.20	1.05
1,2-Dichloroethene (trans)	156-60-5	AAL071156	aa4155rlcs	AA	6/22/15	0.24	0.20	0.8
1,1-Dichloroethane	75-34-3	AAL071156	aa4155rlcs	AA	6/22/15	0.33	0.20	0.6
Methyl tert-butyl ether	1634-04-4	AAL071156	aa4155rlcs	AA	6/22/15	0.26	0.20	0.8
Methyl ethyl ketone	78-93-3	AAL071156	aa4155rlcs	AA	6/22/15	0.34	0.20	0.6
1,2-Dichloroethene (cis)	156-59-2	AAL071156	aa4155rlcs	AA	6/22/15	0.26	0.20	0.8
Ethyl acetate	141-78-6	AAL071156	aa4155rlcs	AA	6/22/15	0.49	0.20	0.4
n-Hexane	110-54-3	AAL071156	aa4155rlcs	AA	6/22/15	0.35	0.20	0.6
Chloroform	67-66-3	AAL071156	aa4155rlcs	AA	6/22/15	0.34	0.20	0.59
Tetrahydrofuran	109-99-9	AAL071156	aa4155rlcs	AA	6/22/15	0.37	0.20	0.5
1,2-Dichloroethane	106-93-4	AAL071156	aa4155rlcs	AA	6/22/15	0.36	0.20	0.56
1,1,1-Trichloroethane	71-55-6	AAL071156	aa4155rlcs	AA	6/22/15	0.33	0.20	0.61
Benzene	71-43-2	AAL071156	aa4155rlcs	AA	6/22/15	0.34	0.20	0.6
Carbon tetrachloride	56-23-5	AAL071156	aa4155rlcs	AA	6/22/15	0.26	0.20	0.77
Cyclohexane	110-82-7	AAL071156	aa4155rlcs	AA	6/22/15	0.27	0.20	0.7
1,2-Dichloropropane	78-87-5	AAL071156	aa4155rlcs	AA	6/22/15	0.45	0.20	0.44
Bromodichloromethane	75-27-4	AAL071156	aa4155rlcs	AA	6/22/15	0.35	0.20	0.57
2,2,4-Trimethylpentane	540-84-1	AAL071156	aa4155rlcs	AA	6/22/15	0.24	0.20	0.8
Trichloroethene	79-01-6	AAL071156	aa4155rlcs	AA	6/22/15	0.4	0.20	0.5
1,4-Dioxane	123-91-1	AAL071156	aa4155rlcs	AA	6/22/15	0.41	0.20	0.5
Methyl methacrylate	80-62-6	AAL071156	aa4155rlcs	AA	6/22/15	0.4	0.20	0.5
n-Heptane	142-82-5	AAL071156	aa4155rlcs	AA	6/22/15	0.37	0.20	0.5
1,3-Dichloropropene (cis)	10061-01-5	AAL071156	aa4155rlcs	AA	6/22/15	0.35	0.20	0.6
Methyl isobutyl ketone	108-10-1	AAL071156	aa4155rlcs	AA	6/22/15	0.49	0.20	0.4
1,3-Dichloropropene (trans)	10061-02-6	AAL071156	aa4155rlcs	AA	6/22/15	0.38	0.20	0.5
1,1,2-Trichloroethane	79-00-5	AAL071156	aa4155rlcs	AA	6/22/15	0.43	0.20	0.47
Toluene	108-88-3	AAL071156	aa4155rlcs	AA	6/22/15	0.44	0.20	0.5
Methyl n-butyl ketone	591-78-6	AAL071156	aa4155rlcs	AA	6/22/15	0.5	0.20	0.4
Dibromochloromethane	75-71-8	AAL071156	aa4155rlcs	AA	6/22/15	0.29	0.20	0.69
1,2-Dibromoethane	107-06-2	AAL071156	aa4155rlcs	AA	6/22/15	0.41	0.20	0.5
Tetrachloroethene	127-18-4	AAL071156	aa4155rlcs	AA	6/22/15	0.43	0.20	0.5
Chlorobenzene	108-90-7	AAL071156	aa4155rlcs	AA	6/22/15	0.21	0.20	0.95
Ethylbenzene	100-41-4	AAL071156	aa4155rlcs	AA	6/22/15	0.42	0.20	0.5
Xylenes (m&p)	179601-23-1	AAL071156	aa4155rlcs	AA	6/22/15	0.37	0.40	1.1
Bromoform	75-25-2	AAL071156	aa4155rlcs	AA	6/22/15	0.77	0.20	0.26
Styrene	100-42-5	AAL071156	aa4155rlcs	AA	6/22/15	0.23	0.20	0.9
Xylene (o)	95-47-6	AAL071156	aa4155rlcs	AA	6/22/15	0.32	0.20	0.6
1,1,2,2-Tetrachloroethane	79-34-5	AAL071156	aa4155rlcs	AA	6/22/15	0.37	0.20	0.54
n-Nonane	111-84-2	AAL071156	aa4155rlcs	AA	6/22/15	0.41	0.20	0.5
Cumene	98-82-8	AAL071156	aa4155rlcs	AA	6/22/15	0.36	0.20	0.13
2-Chlorotoluene	95-49-8	AAL071156	aa4155rlcs	AA	6/22/15	0.36	0.20	0.12
4-Ethyltoluene	622-96-8	AAL071156	aa4155rlcs	AA	6/22/15	0.32	0.20	0.10
1,3,5-Trimethylbenzene	108-67-8	AAL071156	aa4155rlcs	AA	6/22/15	0.33	0.20	0.10
1,2,4-Trimethylbenzene	95-63-6	AAL071156	aa4155rlcs	AA	6/22/15	0.31	0.20	0.090
1,3-Dichlorobenzene	541-73-1	AAL071156	aa4155rlcs	AA	6/22/15	0.39	0.20	0.14
1,4-Dichlorobenzene	106-46-7	AAL071156	aa4155rlcs	AA	6/22/15	0.38	0.20	0.12
1,2-Dichlorobenzene	95-50-1	AAL071156	aa4155rlcs	AA	6/22/15	0.36	0.20	0.13
1,2,4-Trichlorobenzene	120-82-1	AAL071156	aa4155rlcs	AA	6/22/15	0.4	0.20	0.10
Naphthalene	91-20-3	AAL071156	aa4155rlcs	AA	6/22/15	0.36	0.20	0.030
1,3-Hexachlorobutadiene	87-68-3	AAL071156	aa4155rlcs	AA	6/22/15	0.44	0.20	0.22

Location of this file: P:\Paldata\Pal Reports\TO-15 MDLS

REPORTING METHOD DETECTION LIMIT (MDL) REPORT

Integrated Analytical Laboratories - Randolph, NJ

Matrix: Air
 Column ID: Restek Rtx-1, 60 meter, 0.32mm ID, 1 um
 Instrument ID: GC - Agilent 7890A / MS - Agilent 5975C (IAL ID: Instrument AA)
 Report Prepared by: Lauren Jenkins
 Preparation Date: 5/22/18

MDL Effective Date: 5/22/2018
 MDL Expiration Date: 4/9/2019
 Analyst: Jeff Schmitt

Compound Name	CAS #	Molecular Weight	MDL ppbv	PQL ppbv	NY/ LLTO-15/ PA RL ppbv	NY/ LLTO-15/ PA RL µg/m ³	NJ TO-15 RL ppbv	NJ TO-15 RL µg/m ³	True value/ MDL
Propene	115-07-1	42.08	0.13	0.40	0.20	0.34	0.40	0.69	1.7
Dichlorodifluoromethane	124-48-1	120.9	0.15	0.44	0.20	0.99	0.40	2.0	1.5
Chloromethane	74-87-3	50.49	0.20	0.59	0.20	0.41	0.40	0.83	1.1
1,2-Dichlorotetrafluoroethane	76-14-2	170.9	0.12	0.36	0.20	1.4	0.40	2.8	1.7
Vinyl chloride	75-01-4	62.50	0.10	0.31	0.20	0.51	0.40	1.0	2.1
1,3-Butadiene	106-99-0	54.09	0.075	0.22	0.20	0.44	0.40	0.88	2.9
n-Butane	106-97-8	58.12	0.074	0.22	0.20	0.48	0.40	0.95	3.0
Bromomethane	74-83-9	94.94	0.14	0.43	0.20	0.78	0.40	1.6	1.4
Chloroethane	75-00-3	64.52	0.087	0.26	0.20	0.53	0.40	1.1	2.3
Ethanol	64-17-5	46.07	0.068	0.20	0.20	0.38	0.40	0.75	2.8
Vinyl bromide	593-60-2	106.9	0.082	0.25	0.20	0.87	0.40	1.7	2.7
Acrolein	107-02-8	56.06	0.059	0.18	0.20	0.46	0.40	0.92	4.1
Acetone	67-64-1	58.08	0.10	0.29	0.20	0.48	0.40	0.95	2.3
Trichlorofluoromethane	75-69-4	137.4	0.11	0.34	0.20	1.1	0.40	2.2	1.9
Isopropanol	67-63-0	60.10	0.068	0.20	0.20	0.49	0.40	0.98	2.9
n-Pentane	109-66-0	72.15	0.057	0.17	0.20	0.59	0.40	1.2	3.8
1,1-Dichloroethene	75-35-4	96.94	0.068	0.21	0.20	0.79	0.40	1.6	3.2
Methylene chloride	75-09-2	84.94	0.11	0.32	0.20	0.69	0.40	1.4	2.0
Tert-butyl alcohol	75-65-0	74.12	0.060	0.18	0.20	0.61	0.40	1.2	4.0
Allyl Chloride	107-05-1	76.53	0.075	0.23	0.20	0.63	0.40	1.3	2.9
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	187.4	0.13	0.40	0.20	1.5	0.40	3.1	1.6
Carbon disulfide	75-15-0	76.14	0.12	0.36	0.20	0.62	0.40	1.2	1.8
1,2-Dichloroethene (trans)	156-60-5	96.94	0.067	0.20	0.20	0.79	0.40	1.6	3.3
1,1-Dichloroethane	75-34-3	98.96	0.092	0.27	0.20	0.81	0.40	1.6	2.4
Methyl tert-butyl ether	1634-04-4	88.15	0.036	0.11	0.20	0.72	0.40	1.4	6.1
Methyl ethyl ketone	78-93-3	72.11	0.043	0.13	0.20	0.59	0.40	1.2	5.1
1,2-Dichloroethene (cis)	156-59-2	96.94	0.064	0.19	0.20	0.79	0.40	1.6	3.5
Ethyl acetate	141-78-6	88.11	0.090	0.27	0.20	0.72	0.40	1.4	2.4
n-Hexane	110-54-3	86.17	0.084	0.25	0.20	0.70	0.40	1.4	2.6
Chloroform	67-66-3	119.4	0.10	0.30	0.20	0.98	0.40	2.0	2.2
Tetrahydrofuran	109-99-9	72.11	0.047	0.14	0.20	0.59	0.40	1.2	5.1
1,2-Dichloroethane	106-93-4	98.96	0.089	0.27	0.20	0.81	0.40	1.6	2.5
1,1,1-Trichloroethane	71-55-6	133.4	0.10	0.31	0.20	1.1	0.40	2.2	2.1
Benzene	71-43-2	78.11	0.049	0.15	0.20	0.64	0.40	1.3	4.5
Carbon tetrachloride	56-23-5	153.8	0.11	0.33	0.20	1.3	0.40	2.5	2.0
Cyclohexane	110-82-7	84.16	0.040	0.12	0.20	0.69	0.40	1.4	5.4
1,2-Dichloropropane	78-87-5	113.0	0.094	0.28	0.20	0.92	0.40	1.8	2.3
Bromodichloromethane	75-27-4	163.8	0.14	0.41	0.20	1.3	0.40	2.7	1.6
2,2,4-Trimethylpentane	540-84-1	114.2	0.082	0.25	0.20	0.93	0.40	1.9	2.7
Trichloroethene	79-01-6	131.4	0.091	0.27	0.20	1.1	0.40	2.1	2.4
1,4-Dioxane	123-91-1	88.12	0.057	0.17	0.20	0.72	0.40	1.4	3.5
Methyl methacrylate	80-62-6	100.12	0.057	0.17	0.20	0.82	0.40	1.6	3.8
n-Heptane	142-82-5	100.2	0.066	0.20	0.20	0.82	0.40	1.6	3.3
1,3-Dichloropropene (cis)	10061-01-5	111.0	0.062	0.19	0.20	0.91	0.40	1.8	3.5
Methyl isobutyl ketone	108-10-1	100.2	0.080	0.24	0.20	0.82	0.40	1.6	2.8
1,3-Dichloropropene (trans)	10061-02-6	111.0	0.06	0.18	0.20	0.91	0.40	1.8	3.3
1,1,2-Trichloroethane	79-00-5	133.4	0.12	0.36	0.20	1.1	0.40	2.2	1.9
Toluene	108-88-3	92.14	0.057	0.17	0.20	0.75	0.40	1.5	3.9
Methyl n-butyl ketone	591-78-6	100.16	0.072	0.22	0.20	0.82	0.40	1.6	3.1
Dibromochloromethane	75-71-8	208.3	0.11	0.34	0.20	1.7	0.40	3.4	2.0
1,2-Dibromoethane	107-06-2	187.9	0.095	0.28	0.20	1.5	0.40	3.1	2.3
Tetrachloroethene	127-18-4	165.8	0.082	0.25	0.20	1.4	0.40	2.7	2.7
Chlorobenzene	108-90-7	112.6	0.15	0.44	0.20	0.92	0.40	1.8	1.5
Ethylbenzene	100-41-4	106.2	0.091	0.27	0.20	0.87	0.40	1.7	2.4
Xylenes (m&p)	179601-23-1	106.2	0.084	0.25	0.40	1.7	0.40	1.7	5.2
Bromoform	75-25-2	252.8	0.067	0.20	0.20	2.1	0.40	4.1	3.3
Styrene	100-42-5	104.1	0.045	0.14	0.20	0.85	0.40	1.7	4.9
1,1,2,2-Tetrachloroethane	79-34-5	167.9	0.057	0.17	0.20	1.4	0.40	2.7	3.8
Xylene (o)	95-47-6	106.2	0.16	0.48	0.20	0.87	0.40	1.7	1.4
n-Nonane	111-84-2	128.2	0.057	0.17	0.20	1.0	0.40	2.1	3.9
Cumene (Isopropylbenzene)	98-82-8	120.2	0.068	0.20	0.20	0.98	0.40	2.0	3.3
2-Chlorotoluene	95-49-8	126.6	0.055	0.17	0.20	1.0	0.40	2.1	4.0
n-Propyl benzene	103-65-1	120.19	0.059	0.18	0.20	0.98	0.40	2.0	3.8
4-Ethyltoluene	622-96-8	120.2	0.059	0.18	0.20	0.98	0.40	2.0	3.7
1,3,5-Trimethylbenzene	108-67-8	120.2	0.058	0.17	0.20	0.98	0.40	2.0	3.8
1,2,4-Trimethylbenzene	95-63-6	120.2	0.041	0.12	0.20	0.98	0.40	2.0	5.3

REPORTING METHOD DETECTION LIMIT (MDL) REPORT

Integrated Analytical Laboratories - Randolph, NJ

Matrix: Air
 Column ID: Restek Rtx-1, 60 meter, 0.32mm ID, 1 um
 Instrument ID: GC - Agilent 7890A / MS - Agilent 5975C (IAL ID: Instrument AA)
 Report Prepared by: Lauren Jenkins
 Preparation Date: 5/22/18

MDL Effective Date: 5/22/2018
 MDL Expiration Date: 4/9/2019
 Analyst: Jeff Schmitt

Compound Name	CAS #	Molecular Weight	MDL ppbv	PQL ppbv	NY/ LLTO-15/ PA RL ppbv	NY/ LLTO-15/ PA RL µg/m ³	NJ TO-15 RL ppbv	NJ TO-15 RL µg/m ³	True value/ MDL
Benzyl chloride	100-44-7	126.59	0.066	0.20	0.20	1.0	0.40	2.1	3.0
1,3-Dichlorobenzene	541-73-1	147.0	0.071	0.21	0.20	1.2	0.40	2.4	3.1
1,4-Dichlorobenzene	106-46-7	147.0	0.065	0.19	0.20	1.2	0.40	2.4	3.4
1,2-Dichlorobenzene	95-50-1	147.0	0.070	0.21	0.20	1.2	0.40	2.4	3.1
1,2,4-Trichlorobenzene	120-82-1	181.5	0.080	0.24	0.20	1.5	0.40	3.0	3.0
Naphthalene	91-20-3	128.2	0.075	0.23	0.20	1.0	0.40	2.1	3.2
1,3-Hexachlorobutadiene	87-68-3	260.8	0.16	0.47	0.20	2.1	0.40	4.3	1.4

Where:

MDL is defined as the higher of the MDL Spike and MDL Blank

PQL is MDLx3

ppbv is parts per billion by volume and is how results come off the instrument

µg/m³ = ppbv x molecular weight / 24.45

Processing Method: C:\MSDCHEM\1\METHODS\AA0302.M

Initial Calibration: C:\MSDCHEM\1\METHODS\AA0302.M

Location of this file: P:\PAL Reports\2018\LLTO-15 and TO-15 Common Files\Agilent MDL

Instrument used for Clean Canister Certification Analysis? YES



Michael Leftin, Ph.D.
 Laboratory Director

Date: May 24, 2018



Lauren Jenkins
 Quality Assurance Officer

Date: May 24, 2018

METHOD DETECTION LIMIT SPIKE (MDL_s) REPORT

Integrated Analytical Laboratories - Randolph, NJ

Matrix: Air
Method: EPA TO-15/NJDEP LLTO-15

Filename	Run #/Type	Date	Time
aa6444	Spike 1	4/10/2018	22:33
aa6445	Spike 2	4/10/2018	23:08
aa6446	Spike 3	4/10/2018	23:44
aa6479	Spike 4	4/12/2018	13:07
aa6480	Spike 5	4/12/2018	13:42
aa6543	Spike 6	4/16/2018	18:36
aa6544	Spike 7	4/16/2018	19:15

Compound Name	CAS #	Spike 1	Spike 2	Spike 3	Spike 4	Spike 5	Spike 6	Spike 7	Mean Value	Spike Value	Percent Recovery	Standard Deviation	MDLs ppbv	Ratio
Propene	115-07-1	0.27	0.25	0.27	0.26	0.28	0.35	0.35	0.29	0.22	132	0.042	0.13	1.67
Dichlorodifluoromethane	124-48-1	0.34	0.32	0.30	0.30	0.28	0.40	0.39	0.33	0.22	151	0.047	0.147	1.49
Chloromethane	74-87-3	0.25	0.25	0.25	0.26	0.24	0.40	0.34	0.28	0.22	129	0.063	0.20	1.11
1,2-Dichlorotetrafluoroethane	76-14-2	0.43	0.41	0.43	0.39	0.38	0.48	0.46	0.42	0.20	212	0.038	0.121	1.65
Vinyl chloride	75-01-4	0.19	0.16	0.18	0.18	0.18	0.24	0.24	0.20	0.22	89	0.033	0.104	2.11
1,3-Butadiene	106-99-0	0.13	0.11	0.11	0.09	0.12	0.16	0.15	0.12	0.22	56	0.024	0.075	2.95
n-Butane	106-97-8	0.18	0.19	0.19	0.16	0.16	0.22	0.21	0.19	0.22	85	0.024	0.074	2.97
Bromomethane	74-83-9	0.26	0.24	0.23	0.23	0.22	0.32	0.33	0.26	0.20	132	0.046	0.144	1.38
Chloroethane	75-00-3	0.25	0.25	0.21	0.24	0.20	0.26	0.28	0.24	0.20	120	0.028	0.09	2.31
Ethanol	64-17-5	0.10	0.09	0.11	0.10	0.10	0.13	0.15	0.11	0.19	60	0.022	0.07	2.76
Vinyl bromide	593-60-2	0.15	0.16	0.13	0.14	0.12	0.18	0.19	0.15	0.22	70	0.026	0.082	2.67
Acrolein	107-02-8	0.11	0.14	0.13	0.12	0.14	0.15	0.16	0.14	0.24	56	0.019	0.06	4.09
Acetone	67-64-1	0.21	0.21	0.19	0.18	0.21	0.25	0.27	0.22	0.22	99	0.031	0.097	2.27
Trichlorofluoromethane	75-69-4	0.35	0.35	0.34	0.32	0.30	0.41	0.39	0.35	0.22	159	0.036	0.114	1.93
Isopropanol	67-63-0	0.10	0.10	0.11	0.10	0.10	0.14	0.15	0.11	0.20	57	0.022	0.07	2.93
n-Pentane	109-66-0	0.13	0.15	0.13	0.12	0.12	0.17	0.16	0.14	0.22	65	0.018	0.057	3.83
1,1-Dichloroethene	75-35-4	0.15	0.16	0.15	0.14	0.14	0.19	0.19	0.16	0.22	72	0.022	0.068	3.21
Methylene chloride	75-09-2	0.30	0.30	0.29	0.30	0.28	0.37	0.36	0.31	0.22	142	0.034	0.108	2.03
Tert-butyl alcohol	75-65-0	0.15	0.14	0.13	0.13	0.13	0.17	0.18	0.15	0.24	62	0.019	0.060	3.99
Allyl Chloride	107-05-1	0.14	0.15	0.13	0.12	0.10	0.18	0.15	0.14	0.22	63	0.024	0.075	2.93
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	0.42	0.41	0.40	0.38	0.36	0.48	0.46	0.41	0.22	188	0.043	0.134	1.64
Carbon disulfide	75-15-0	0.26	0.25	0.25	0.24	0.23	0.32	0.32	0.27	0.22	122	0.038	0.119	1.85
1,2-Dichloroethene (trans)	156-60-5	0.18	0.18	0.18	0.17	0.16	0.22	0.21	0.18	0.22	84	0.021	0.067	3.26
1,1-Dichloroethane	75-34-3	0.28	0.27	0.27	0.25	0.26	0.33	0.31	0.28	0.22	128	0.029	0.092	2.40
Methyl tert-butyl ether	1634-04-4	0.13	0.13	0.13	0.12	0.11	0.15	0.14	0.13	0.22	58	0.012	0.036	6.08
Methyl ethyl ketone	78-93-3	0.15	0.16	0.16	0.15	0.15	0.18	0.18	0.16	0.22	73	0.014	0.043	5.09
1,2-Dichloroethene (cis)	156-59-2	0.16	0.17	0.15	0.15	0.14	0.20	0.19	0.17	0.22	75	0.020	0.064	3.46
Ethyl acetate	141-78-6	0.24	0.21	0.21	0.22	0.17	0.25	0.25	0.22	0.22	100	0.029	0.09	2.45
n-Hexane	110-54-3	0.16	0.16	0.16	0.15	0.13	0.21	0.20	0.17	0.22	76	0.027	0.084	2.62
Chloroform	67-66-3	0.33	0.33	0.32	0.30	0.29	0.38	0.37	0.33	0.22	150	0.032	0.100	2.19
Tetrahydrofuran	109-99-9	0.11	0.12	0.12	0.11	0.12	0.14	0.15	0.12	0.24	52	0.015	0.047	5.09
1,2-Dichloroethane	106-93-4	0.27	0.26	0.26	0.25	0.24	0.31	0.31	0.27	0.22	124	0.028	0.089	2.47
1,1,1-Trichloroethane	71-55-6	0.32	0.30	0.31	0.29	0.27	0.37	0.35	0.32	0.22	144	0.033	0.105	2.10
Benzene	71-43-2	0.20	0.21	0.20	0.19	0.19	0.23	0.23	0.21	0.22	94	0.016	0.049	4.46
Carbon tetrachloride	56-23-5	0.34	0.32	0.32	0.30	0.29	0.39	0.37	0.33	0.22	151	0.035	0.109	2.01
Cyclohexane	110-82-7	0.13	0.12	0.12	0.12	0.11	0.14	0.14	0.13	0.22	57	0.013	0.040	5.45
1,2-Dichloropropane	78-87-5	0.34	0.33	0.32	0.30	0.30	0.36	0.38	0.33	0.22	150	0.030	0.094	2.33
Bromodichloromethane	75-27-4	0.39	0.38	0.36	0.32	0.33	0.44	0.43	0.38	0.22	172	0.044	0.137	1.60
2,2,4-Trimethylpentane	540-84-1	0.16	0.16	0.16	0.15	0.15	0.21	0.20	0.17	0.22	77	0.026	0.082	2.68
Trichloroethene	79-01-6	0.30	0.28	0.27	0.26	0.24	0.33	0.29	0.28	0.22	128	0.029	0.091	2.42
1,4-Dioxane	123-91-1	0.19	0.21	0.21	0.16	0.19	0.22	0.21	0.20	0.20	100	0.018	0.057	3.50
Methyl methacrylate	80-62-6	0.16	0.15	0.15	0.14	0.13	0.18	0.17	0.15	0.22	70	0.018	0.057	3.85
n-Heptane	142-82-5	0.18	0.16	0.18	0.15	0.15	0.20	0.20	0.17	0.22	79	0.021	0.066	3.35
1,3-Dichloropropene (cis)	10061-01-5	0.20	0.19	0.19	0.17	0.17	0.22	0.22	0.19	0.22	88	0.020	0.062	3.52
Methyl isobutyl ketone	108-10-1	0.18	0.18	0.19	0.16	0.17	0.24	0.21	0.19	0.22	87	0.025	0.08	2.76
1,3-Dichloropropene (trans)	10061-02-6	0.18	0.19	0.17	0.16	0.16	0.21	0.21	0.18	0.20	92	0.019	0.060	3.34
1,1,2-Trichloroethane	79-00-5	0.40	0.39	0.38	0.36	0.35	0.45	0.43	0.39	0.22	180	0.038	0.119	1.85
Toluene	108-88-3	0.18	0.18	0.17	0.17	0.16	0.20	0.21	0.18	0.22	83	0.018	0.057	3.89
Methyl n-butyl ketone	591-78-6	0.16	0.16	0.16	0.14	0.13	0.19	0.19	0.16	0.22	74	0.023	0.07	3.06
Dibromochloromethane	75-71-8	0.36	0.34	0.35	0.33	0.32	0.42	0.40	0.36	0.22	164	0.036	0.112	1.96
1,2-Dibromoethane	107-06-2	0.29	0.28	0.28	0.27	0.25	0.34	0.32	0.29	0.22	132	0.030	0.095	2.32
Tetrachloroethene	127-18-4	0.29	0.31	0.28	0.27	0.26	0.34	0.32	0.30	0.22	134	0.026	0.082	2.69
Chlorobenzene	108-90-7	0.52	0.51	0.50	0.47	0.45	0.58	0.56	0.51	0.22	233	0.046	0.145	1.52
Ethylbenzene	100-41-4	0.28	0.27	0.28	0.25	0.24	0.33	0.30	0.28	0.22	127	0.029	0.091	2.42
Xylenes (m&p)	179601-23-1	0.26	0.25	0.25	0.22	0.21	0.29	0.28	0.25	0.44	57	0.027	0.08	5.23
Bromoform	75-25-2	0.19	0.18	0.18	0.16	0.16	0.21	0.21	0.18	0.22	83	0.021	0.067	3.28
Styrene	100-42-5	0.18	0.17	0.16	0.16	0.15	0.19	0.18	0.17	0.22	78	0.014	0.045	4.87
1,1,2,2-Tetrachloroethane	79-34-5	0.15	0.16	0.16	0.14	0.14	0.18	0.18	0.16	0.22	72	0.018	0.057	3.83
Xylene (o)	95-47-6	0.44	0.43	0.42	0.38	0.37	0.51	0.48	0.43	0.22	197	0.051	0.161	1.37
n-Nonane	111-84-2	0.14	0.14	0.13	0.12	0.12	0.16	0.16	0.14	0.22	64	0.018	0.057	3.87

METHOD DETECTION LIMIT SPIKE (MDL_s) REPORT

Integrated Analytical Laboratories - Randolph, NJ

Matrix: Air

Method: EPA TO-15/NJDEP LLTO-15

Filename	Run #/Type	Date	Time
aa6444	Spike 1	4/10/2018	22:33
aa6445	Spike 2	4/10/2018	23:08
aa6446	Spike 3	4/10/2018	23:44
aa6479	Spike 4	4/12/2018	13:07
aa6480	Spike 5	4/12/2018	13:42
aa6543	Spike 6	4/16/2018	18:36
aa6544	Spike 7	4/16/2018	19:15

Compound Name	CAS #	Spike 1	Spike 2	Spike 3	Spike 4	Spike 5	Spike 6	Spike 7	Mean Value	Spike Value	Percent Recovery	Standard Deviation	MDLs ppbv	Ratio
Cumene (Isopropylbenzene)	98-82-8	0.22	0.21	0.21	0.19	0.18	0.24	0.23	0.21	0.22	97	0.022	0.068	3.25
2-Chlorotoluene	95-49-8	0.18	0.16	0.17	0.15	0.15	0.19	0.19	0.17	0.22	77	0.018	0.055	3.99
n-Propyl benzene	103-65-1	0.16	0.16	0.15	0.14	0.14	0.19	0.18	0.16	0.22	74	0.019	0.059	3.76
4-Ethyltoluene	622-96-8	0.14	0.14	0.14	0.12	0.12	0.16	0.17	0.14	0.22	64	0.019	0.059	3.71
1,3,5-Trimethylbenzene	108-67-8	0.15	0.15	0.14	0.13	0.12	0.17	0.17	0.15	0.22	66	0.019	0.058	3.77
1,2,4-Trimethylbenzene	95-63-6	0.13	0.12	0.13	0.11	0.11	0.14	0.14	0.12	0.22	56	0.013	0.041	5.33
Benzyl chloride	100-44-7	0.18	0.17	0.17	0.15	0.15	0.21	0.19	0.17	0.20	87	0.021	0.066	3.03
1,3-Dichlorobenzene	541-73-1	0.23	0.22	0.22	0.20	0.20	0.26	0.25	0.23	0.22	102	0.023	0.071	3.08
1,4-Dichlorobenzene	106-46-7	0.19	0.18	0.18	0.17	0.17	0.21	0.22	0.19	0.22	86	0.021	0.065	3.40
1,2-Dichlorobenzene	95-50-1	0.20	0.20	0.20	0.18	0.17	0.23	0.22	0.20	0.22	91	0.022	0.070	3.12
1,2,4-Trichlorobenzene	120-82-1	0.11	0.11	0.10	0.11	0.10	0.16	0.16	0.12	0.24	50	0.025	0.080	3.00
Naphthalene	91-20-3	0.06	0.04	0.04	0.05	0.05	0.10	0.09	0.06	0.24	25	0.024	0.075	3.20
1,3-Hexachlorobutadiene	87-68-3	0.36	0.34	0.32	0.31	0.31	0.44	0.41	0.36	0.22	162	0.050	0.157	1.40

METHOD DETECTION LIMIT BLANK (MDL_B) REPORT

Integrated Analytical Laboratories - Randolph, NJ

Matrix: Air
 Method: EPA TO-15/
 NJDEP LLTO-15

Filename	Run #/Type	Date	Time
aa6439	Blank 1	4/10/2018	19:42
aa6443	Blank 2	4/10/2018	21:57
aa6476	Blank 3	4/12/2018	23:24
aa6477	Blank 4	4/12/2018	11:58
aa6478	Blank 5	4/12/2018	12:31
aa6541	Blank 6	4/16/2018	17:22
aa6542	Blank 7	4/16/2018	17:56

Compound Name	CAS #	Blank 1	Blank 2	Blank 3	Blank 4	Blank 5	Blank 6	Blank 7	Mean Value	Standard Deviation	MDLb ppbv
Propene	115-07-1	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Dichlorodifluoromethane	124-48-1	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Chloromethane	74-87-3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,2-Dichlorotetrafluoroethane	76-14-2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Vinyl chloride	75-01-4	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,3-Butadiene	106-99-0	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
n-Butane	106-97-8	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Bromomethane	74-83-9	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Chloroethane	75-00-3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Ethanol	64-17-5	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Vinyl bromide	593-60-2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Acrolein	107-02-8	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Acetone	67-64-1	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Trichlorofluoromethane	75-69-4	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Isopropanol	67-63-0	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
n-Pentane	109-66-0	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,1-Dichloroethene	75-35-4	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Methylene chloride	75-09-2	0.071	0.073	0.063	0.067	0.047	0.065	0.079	0.066	0.010	0.031
Tert-butyl alcohol	75-65-0	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Allyl Chloride	107-05-1	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Carbon disulfide	75-15-0	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,2-Dichloroethene (trans)	156-60-5	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,1-Dichloroethane	75-34-3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Methyl tert-butyl ether	1634-04-4	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Methyl ethyl ketone	78-93-3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,2-Dichloroethene (cis)	156-59-2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Ethyl acetate	141-78-6	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
n-Hexane	110-54-3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Chloroform	67-66-3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Tetrahydrofuran	109-99-9	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,2-Dichloroethane	106-93-4	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,1,1-Trichloroethane	71-55-6	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Benzene	71-43-2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Carbon tetrachloride	56-23-5	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Cyclohexane	110-82-7	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,2-Dichloropropane	78-87-5	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Bromodichloromethane	75-27-4	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
2,2,4-Trimethylpentane	540-84-1	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Trichloroethene	79-01-6	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,4-Dioxane	123-91-1	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Methyl methacrylate	80-62-6	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
n-Heptane	142-82-5	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,3-Dichloropropene (cis)	10061-01-5	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Methyl isobutyl ketone	108-10-1	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,3-Dichloropropene (trans)	10061-02-6	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,1,2-Trichloroethane	79-00-5	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Toluene	108-88-3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Methyl n-butyl ketone	591-78-6	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Dibromochloromethane	75-71-8	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,2-Dibromoethane	107-06-2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Tetrachloroethene	127-18-4	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Chlorobenzene	108-90-7	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Ethylbenzene	100-41-4	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Xylenes (m&p)	179601-23-1	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Bromoform	75-25-2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Styrene	100-42-5	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,1,2,2-Tetrachloroethane	79-34-5	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Xylene (o)	95-47-6	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
n-Nonane	111-84-2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00

METHOD DETECTION LIMIT BLANK (MDL_B) REPORT

Integrated Analytical Laboratories - Randolph, NJ

Matrix: Air

Method: EPA TO-15/
NJDEP LLTO-15

Filename	Run #/Type	Date	Time
aa6439	Blank 1	4/10/2018	19:42
aa6443	Blank 2	4/10/2018	21:57
aa6476	Blank 3	4/12/2018	23:24
aa6477	Blank 4	4/12/2018	11:58
aa6478	Blank 5	4/12/2018	12:31
aa6541	Blank 6	4/16/2018	17:22
aa6542	Blank 7	4/16/2018	17:56

Compound Name	CAS #	Blank 1	Blank 2	Blank 3	Blank 4	Blank 5	Blank 6	Blank 7	Mean Value	Standard Deviation	MDLb ppbv
Cumene (Isopropylbenzene)	98-82-8	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
2-Chlorotoluene	95-49-8	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
n-Propyl benzene	103-65-1	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
4-Ethyltoluene	622-96-8	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,3,5-Trimethylbenzene	108-67-8	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,2,4-Trimethylbenzene	95-63-6	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Benzyl chloride	100-44-7	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,3-Dichlorobenzene	541-73-1	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,4-Dichlorobenzene	106-46-7	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,2-Dichlorobenzene	95-50-1	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,2,4-Trichlorobenzene	120-82-1	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
Naphthalene	91-20-3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
1,3-Hexachlorobutadiene	87-68-3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00

METHOD DETECTION LIMIT VERIFICATION (MDLV) REPORT

Integrated Analytical Laboratories - Randolph, NJ

Analysis Level: 0.20 ppbv, 0.40 for m&p-xylenes
 Matrix: Air
 Column ID: Restek Rtx-1, 60 meter, 0.32mm ID, 1 um
 Instrument Identification: AA
 Date of Verification Study: 4/17/2018
 Study Identification File #: aa6565rlcs
 Analyst: Jeff Schmitt
 Analysis/Processing Method: C:\MSDCHEM\1\METHODS\AA0302.M
 Cylinder ID: CC483586

Lauren Jenkins

Lauren Jenkins
 QA Officer

Date: May 22, 2018

Compound Name	CAS #	MDLV (ppbv)	RL (ppbv)	RL/MDLV Ratio
Propene	115-07-1	0.43	0.20	0.47
Dichlorodifluoromethane	124-48-1	0.16	0.20	1.25
Chloromethane	74-87-3	0.34	0.20	0.59
1,2-Dichlorotetrafluoroethane	76-14-2	0.25	0.20	0.80
Vinyl chloride	75-01-4	0.26	0.20	0.77
1,3-Butadiene	106-99-0	0.20	0.20	1.00
n-Butane	106-97-8	0.25	0.20	0.80
Bromomethane	74-83-9	0.15	0.20	1.33
Chloroethane	75-00-3	0.25	0.20	0.80
Ethanol	64-17-5	0.15	0.20	1.33
Vinyl bromide	593-60-2	0.16	0.20	1.25
Acrolein	107-02-8	0.15	0.20	1.33
Acetone	67-64-1	0.26	0.20	0.77
Trichlorofluoromethane	75-69-4	0.27	0.20	0.74
Isopropanol	67-63-0	0.17	0.20	1.18
n-Pentane	109-66-0	0.19	0.20	1.05
1,1-Dichloroethene	75-35-4	0.21	0.20	0.95
Methylene chloride	75-09-2	0.36	0.20	0.56
Tert-butyl alcohol	75-65-0	0.2	0.20	1.00
Allyl Chloride	107-05-1	0.19	0.20	1.05
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	0.25	0.20	0.80
Carbon disulfide	75-15-0	0.22	0.20	0.91
1,2-Dichloroethene (trans)	156-60-5	0.24	0.20	0.83
1,1-Dichloroethane	75-34-3	0.20	0.20	1.00
Methyl tert-butyl ether	1634-04-4	0.17	0.20	1.18
Methyl ethyl ketone	78-93-3	0.2	0.20	1.00
1,2-Dichloroethene (cis)	156-59-2	0.22	0.20	0.91
Ethyl acetate	141-78-6	0.26	0.20	0.77
n-Hexane	110-54-3	0.21	0.20	0.95
Chloroform	67-66-3	0.25	0.20	0.80
Tetrahydrofuran	109-99-9	0.16	0.20	1.25
1,2-Dichloroethane	106-93-4	0.22	0.20	0.91
1,1,1-Trichloroethane	71-55-6	0.24	0.20	0.83
Benzene	71-43-2	0.25	0.20	0.80
Carbon tetrachloride	56-23-5	0.27	0.20	0.74
Cyclohexane	110-82-7	0.17	0.20	1.18
1,2-Dichloropropane	78-87-5	0.24	0.20	0.83
Bromodichloromethane	75-27-4	0.25	0.20	0.80
2,2,4-Trimethylpentane	540-84-1	0.19	0.20	1.05
Trichloroethene	79-01-6	0.18	0.20	1.11
1,4-Dioxane	123-91-1	0.24	0.20	0.83
Methyl methacrylate	80-62-6	0.17	0.20	1.18
n-Heptane	142-82-5	0.20	0.20	1.00
1,3-Dichloropropene (cis)	10061-01-5	0.22	0.20	0.91
Methyl isobutyl ketone	108-10-1	0.22	0.20	0.91
1,3-Dichloropropene (trans)	10061-02-6	0.20	0.20	1.00
1,1,2-Trichloroethane	79-00-5	0.27	0.20	0.74
Toluene	108-88-3	0.20	0.20	1.00
Methyl n-butyl ketone	591-78-6	0.18	0.20	1.11
Dibromochloromethane	75-71-8	0.26	0.20	0.77
1,2-Dibromoethane	107-06-2	0.23	0.20	0.87
Tetrachloroethene	127-18-4	0.21	0.20	0.95

METHOD DETECTION LIMIT VERIFICATION (MDLV) REPORT

Integrated Analytical Laboratories - Randolph, NJ

Analysis Level: 0.20 ppbv, 0.40 for m&p-xylenes
 Matrix: Air
 Column ID: Restek Rtx-1, 60 meter, 0.32mm ID, 1 um
 Instrument Identification: AA
 Date of Verification Study: 4/17/2018
 Study Identification File #: aa6565rllcs
 Analyst: Jeff Schmitt
 Analysis/Processing Method: C:\MSDCHEM\1\METHODS\AA0302.M
 Cylinder ID: CC483586

Lauren Jenkins

Lauren Jenkins
 QA Officer

Date: May 22, 2018

Compound Name	CAS #	MDLV (ppbv)	RL (ppbv)	RL/MDLV Ratio
Chlorobenzene	108-90-7	0.25	0.20	0.80
Ethylbenzene	100-41-4	0.19	0.20	1.05
Xylenes (m&p)	179601-23-1	0.35	0.40	1.14
Bromoform	75-25-2	0.24	0.20	0.83
Styrene	100-42-5	0.15	0.20	1.33
Xylene (o)	95-47-6	0.17	0.20	1.18
1,1,2,2-Tetrachloroethane	79-34-5	0.21	0.20	0.95
n-Nonane	111-84-2	0.15	0.20	1.33
Cumene	98-82-8	0.23	0.20	0.87
2-Chlorotoluene	95-49-8	0.20	0.20	1.00
n-Propyl benzene	103-65-1	0.18	0.20	1.11
4-Ethyltoluene	622-96-8	0.16	0.20	1.25
1,3,5-Trimethylbenzene	108-67-8	0.17	0.20	1.18
1,2,4-Trimethylbenzene	95-63-6	0.14	0.20	1.43
Benzyl chloride	100-44-7	0.20	0.20	1.00
1,3-Dichlorobenzene	541-73-1	0.26	0.20	0.77
1,4-Dichlorobenzene	106-46-7	0.23	0.20	0.87
1,2-Dichlorobenzene	95-50-1	0.22	0.20	0.91
1,2,4-Trichlorobenzene	120-82-1	0.17	0.20	1.18
Naphthalene	91-20-3	0.11	0.20	1.82
1,3-Hexachlorobutadiene	87-68-3	0.40	0.20	0.50

Section V: Quality Control Data Summary

BFB Tune Summary

Method Blank

Laboratory Sample Duplicate

Internal Standard Area Summary

Data Path: C:\DATA\05-18-18\
Data File: AA7071BFB.D
Acq On: 5/18/2018 8:17:00AM
Operator: jls
Sample: BFB
Misc: ALM029426
ALS Vial: 1 **Multiplier:** 1
Integration File: rteint.p
Method: C:\msdchem\1\METHODS\0511.M
Last Update: Mon May 14 11:16:39 2018

Spectrum Information:

Pass/Fail	Target Mass	Rel. to Mass	Lower Limit %	Higher Limit %	Raw Abundance	% Relative Abundance
PASS	50	95	8	40	183818	27.3
PASS	75	95	30	66	378859	56.2
PASS	95	95	100	100	674347	100.0
PASS	96	95	5	9	45051	6.7
PASS	173	174	0.00	2	2912	0.8
PASS	174	95	50	100	387670	57.5
PASS	175	174	4	9	26268	6.8
PASS	176	174	93	101	373555	96.4
PASS	177	176	5	9	24192	6.5

Runs with this BFB:

Lab Sample Number	Date File	Field Sample	Date/Time of Sample/Standard Analysis
BFB	AA7071BFB	NA	5/18/2018 8:17:00 AM
40 PPBV STD	AA7073STD01	NA	5/18/2018 9:46:00 AM
20 PPBV STD	AA7074STD02	NA	5/18/2018 10:19:00 AM
10 PPBV STD	AA7075STD03	NA	5/18/2018 10:53:00 AM
2 PPBV STD	AA7076STD04	NA	5/18/2018 12:20:00 PM
0.2 PPBV STD	AA7077STD05	NA	5/18/2018 1:03:00 PM
10 PPBV ICVSS	AA7078ICVSS	NA	5/18/2018 3:14:00 PM

Data Path: C:\DATA\06-13-18\
Data File: AA7471BFB.D
Acq On: 6/13/2018 8:30:00AM
Operator: jls
Sample: BFB
Misc: ALM029426
ALS Vial: 1 **Multiplier:** 1
Integration File: rteint.p
Method: C:\msdchem\1\METHODS\0518.M
Last Update: Fri May 18 13:51:08 2018

Spectrum Information:

PassFail	Target Mass	Rel. to Mass	Lower Limit %	Higher Limit %	Raw Abundance	% Relative Abundance
PASS	50	95	8	40	200102	30.6
PASS	75	95	30	66	384246	58.8
PASS	95	95	100	100	653589	100.0
PASS	96	95	5	9	42346	6.5
PASS	173	174	0.00	2	3805	1.0
PASS	174	95	50	100	363368	55.6
PASS	175	174	4	9	26616	7.3
PASS	176	174	93	101	348380	95.9
PASS	177	176	5	9	23261	6.7

Runs with this BFB:

Lab Sample Number	Date File	Field Sample	Date/Time of Sample/Standard Analysis
BFB	AA7471BFB	NA	6/13/2018 8:30:00 AM
10 PPBV DCVS	AA7472DCVS	NA	6/13/2018 9:18:00 AM
METHOD BLANK	AA7473BLK	NA	6/13/2018 10:25:00 AM
02 PPBV RLLCS	AA7474RLLCS	NA	6/13/2018 11:14:00 AM
3059	AA7475	NA	6/13/2018 12:20:00 PM
10 PPBV CCCVS	AA7485CCCVS	NA	6/13/2018 8:47:00 PM

Data Path: C:\DATA\07-25-18\
Data File: AA7971BFB.D
Acq On: 7/25/2018 8:35:00AM
Operator: jls
Sample: BFB
Misc: ALM029426
ALS Vial: 1 **Multiplier:** 1
Integration File: rteint.p
Method: C:\msdchem\1\METHODS\0518.M
Last Update: Fri May 18 13:51:08 2018

Spectrum Information:

PassFail	Target Mass	Rel. to Mass	Lower Limit %	Higher Limit %	Raw Abundance	% Relative Abundance
PASS	50	95	8	40	156657	20.1
PASS	75	95	30	66	381141	49.0
PASS	95	95	100	100	778176	100.0
PASS	96	95	5	9	52476	6.7
PASS	173	174	0.00	2	3085	0.7
PASS	174	95	50	100	459712	59.1
PASS	175	174	4	9	31097	6.8
PASS	176	174	93	101	438861	95.5
PASS	177	176	5	9	27982	6.4

Runs with this BFB:

Lab Sample Number	Date File	Field Sample	Date/Time of Sample/Standard Analysis
BFB	AA7971BFB	NA	7/25/2018 8:35:00 AM
40 PPBV STD	AA7972STD01	NA	7/25/2018 9:55:00 AM
20 PPBV STD	AA7973STD02	NA	7/25/2018 10:29:00 AM
10 PPBV STD	AA7974STD03	NA	7/25/2018 11:02:00 AM
2 PPBV STD	AA7975STD04	NA	7/25/2018 12:18:00 PM
0.2 PPBV STD	AA7976STD05	NA	7/25/2018 1:42:00 PM
10 PPBV ICVSS	AA7977ICVSS	NA	7/25/2018 3:05:00 PM

Data Path: C:\DATA\08-06-18\
Data File: AA8191BFB.D
Acq On: 8/6/2018 11:26:00AM
Operator: jls
Sample: BFB
Misc: ALM029426
ALS Vial: 1 **Multiplier:** 1
Integration File: rteint.p
Method: C:\msdchem\1\METHODS\0725.M
Last Update: Wed Jul 25 14:15:57 2018
Spectrum Information:

PassFail	Target Mass	Rel. to Mass	Lower Limit %	Higher Limit %	Raw Abundance	% Relative Abundance
PASS	50	95	8	40	177003	32.6
PASS	75	95	30	66	334482	61.6
PASS	95	95	100	100	542971	100.0
PASS	96	95	5	9	36432	6.7
PASS	173	174	0.00	2	2635	0.9
PASS	174	95	50	100	286464	52.8
PASS	175	174	4	9	21901	7.6
PASS	176	174	93	101	278015	97.1
PASS	177	176	5	9	17336	6.2

Runs with this BFB:

Lab Sample Number	Date File	Field Sample	Date/Time of Sample/Standard Analysis
BFB	AA8191BFB	NA	8/6/2018 11:26:00 AM
10 PPBV DCVS	AA8192DCVS	NA	8/6/2018 1:07:00 PM
METHOD BLANK	AA8193BLK	NA	8/6/2018 1:42:00 PM
02 PPBV RLLCS	AA8194RLLCS	NA	8/6/2018 2:23:00 PM
E18-06141-07	AA8206	SS-101	8/6/2018 10:28:00 PM
E18-06141-08	AA8207	SS-102	8/6/2018 11:02:00 PM
E18-06141-09	AA8208	SS-103	8/6/2018 11:35:00 PM
10 PPBV CCCVS	AA8213CCCVS	NA	8/7/2018 2:22:00 AM

Data Path: C:\DATA\08-07-18\
Data File: AA8221BFB.D
Acq On: 8/7/2018 9:33:00AM
Operator: jls
Sample: BFB
Misc: ALM029426
ALS Vial: 1 **Multiplier:** 1
Integration File: rteint.p
Method: C:\msdchem\1\METHODS\0725.M
Last Update: Wed Jul 25 14:15:57 2018

Spectrum Information:

PassFail	Target Mass	Rel. to Mass	Lower Limit %	Higher Limit %	Raw Abundance	% Relative Abundance
PASS	50	95	8	40	155635	33.4
PASS	75	95	30	66	289323	62.1
PASS	95	95	100	100	466217	100.0
PASS	96	95	5	9	29664	6.4
PASS	173	174	0.00	2	2337	1.0
PASS	174	95	50	100	240853	51.7
PASS	175	174	4	9	17856	7.4
PASS	176	174	93	101	230955	95.9
PASS	177	176	5	9	15117	6.5

Runs with this BFB:

Lab Sample Number	Date File	Field Sample	Date/Time of Sample/Standard Analysis
BFB	AA8221BFB	NA	8/7/2018 9:33:00 AM
10 PPBV DCVS	AA8222DCVS	NA	8/7/2018 10:17:00 AM
METHOD BLANK	AA8223BLK	NA	8/7/2018 11:01:00 AM
02 PPBV RLLCS	AA8224RLLCS	NA	8/7/2018 11:47:00 AM
E18-06141-08	AA8232	SS-102	8/7/2018 4:57:00 PM
E18-06141-09	AA8233	SS-103	8/7/2018 5:31:00 PM
10 PPBV CCCVS	AA8242CCCVS	NA	8/7/2018 10:31:00 PM

Method Blank Report

Lab Sample Name: METHOD BLANK
Field Sample Name: METHOD BLANK
Matrix: Air
Dilution Factor: 1

Data File: AA7473BLK
Date Analyzed: 6/13/2018
Sample Volume: 500ml
GC/MS Column: RTX-1, 0.32 mmID

Runs with this Method Blank:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7471BFB]	06/13/2018 8:30
10 PPBV DCVS [AA7472DCVS]	06/13/2018 9:18
METHOD BLANK [AA7473BLK]	06/13/2018 10:25
02 PPBV RLLCS [AA7474RLLCS]	06/13/2018 11:14
3059 [AA7475]	06/13/2018 12:20
10 PPBV CCCVS [AA7485CCCVS]	06/13/2018 20:47

Compound	CAS #	Reporting Limit (ppbv)	Concentration (ppbv)
Acetone	67-64-1	0.20	ND
Benzene	71-43-2	0.15	ND
Bromodichloromethane	75-27-4	0.20	ND
Bromoform	75-25-2	0.20	ND
Bromomethane	74-83-9	0.20	ND
1,3-Butadiene	106-99-0	0.20	ND
Chlorobenzene	108-90-7	0.20	ND
Chloroethane	75-00-3	0.20	ND
Chloroform	67-66-3	0.20	ND
Chloromethane	74-87-3	0.20	ND
Carbon disulfide	75-15-0	0.20	ND
Carbon tetrachloride	56-23-5	0.20	ND
Cyclohexane	110-82-7	0.12	ND
Dibromochloromethane	124-48-1	0.20	ND
1,2-Dibromoethane	106-93-4	0.20	ND
1,2-Dichlorobenzene	95-50-1	0.20	ND
1,3-Dichlorobenzene	541-73-1	0.20	ND
1,4-Dichlorobenzene	106-46-7	0.20	ND
Dichlorodifluoromethane	75-71-8	0.20	ND
1,1-Dichloroethane	75-34-3	0.20	ND
1,2-Dichloroethane	107-06-2	0.20	ND
1,1-Dichloroethene	75-35-4	0.20	ND
1,2-Dichloroethene (cis)	156-59-2	0.19	ND
1,2-Dichloroethene (trans)	156-60-5	0.20	ND
1,2-Dichloropropane	78-87-5	0.20	ND
1,3-Dichloropropene (cis)	10061-01-5	0.19	ND
1,3-Dichloropropene (trans)	10061-02-6	0.18	ND
1,2-Dichlorotetrafluoroethane	76-14-2	0.20	ND
1,4-Dioxane	123-91-1	0.17	ND
Ethylbenzene	100-41-4	0.20	ND
n-Heptane	142-82-5	0.20	ND
1,3-Hexachlorobutadiene	87-68-3	0.20	ND
n-Hexane	110-54-3	0.20	ND
Methylene chloride	75-09-2	0.20	ND
Methyl ethyl ketone	78-93-3	0.13	ND
Methyl isobutyl ketone	108-10-1	0.20	ND

Method Blank must be less than the Practical Quantitation Limit (PQL).

Method Blank Report

Lab Sample Name: METHOD BLANK
Field Sample Name: METHOD BLANK
Matrix: Air
Dilution Factor: 1

Data File: AA7473BLK
Date Analyzed: 6/13/2018
Sample Volume: 500ml
GC/MS Column: RTX-1, 0.32 mmID

Runs with this Method Blank:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7471BFB]	06/13/2018 8:30
10 PPBV DCVS [AA7472DCVS]	06/13/2018 9:18
METHOD BLANK [AA7473BLK]	06/13/2018 10:25
02 PPBV RLLCS [AA7474RLLCS]	06/13/2018 11:14
3059 [AA7475]	06/13/2018 12:20
10 PPBV CCCVS [AA7485CCCVS]	06/13/2018 20:47

Compound	CAS #	Reporting Limit (ppbv)	Concentration (ppbv)
Methyl tert-butyl ether	1634-04-4	0.11	ND
Styrene	100-42-5	0.14	ND
Tert-butyl alcohol	75-65-0	0.18	ND
1,1,2,2-Tetrachloroethane	79-34-5	0.17	ND
Tetrachloroethene	127-18-4	0.20	ND
Toluene	108-88-3	0.17	ND
1,2,4-Trichlorobenzene	120-82-1	0.20	ND
1,1,1-Trichloroethane	71-55-6	0.20	ND
1,1,2-Trichloroethane	79-00-5	0.20	ND
Trichloroethene	79-01-6	0.20	ND
Trichlorofluoromethane	75-69-4	0.20	ND
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	0.20	ND
1,2,4-Trimethylbenzene	95-63-6	0.12	ND
1,3,5-Trimethylbenzene	108-67-8	0.17	ND
2,2,4-Trimethylpentane	540-84-1	0.20	ND
Vinyl bromide	593-60-2	0.20	ND
Vinyl chloride	75-01-4	0.20	ND
Xylenes (m&p)	179601-23-1	0.20	ND
Xylenes (o)	95-47-6	0.20	ND

Method Blank must be less than the Practical Quantitation Limit (PQL).

Method Blank Report

Lab Sample Name: METHOD BLANK
Field Sample Name: METHOD BLANK
Matrix: Air
Dilution Factor: 1

Data File: AA8193BLK
Date Analyzed: 8/6/2018
Sample Volume: 500ml
GC/MS Column: RTX-1, 0.32 mmID

Runs with this Method Blank:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA8191BFB]	08/06/2018 11:26
10 PPBV DCVS [AA8192DCVS]	08/06/2018 13:07
METHOD BLANK [AA8193BLK]	08/06/2018 13:42
02 PPBV RLLCS [AA8194RLLCS]	08/06/2018 14:23
E18-06141-07 [AA8206]	08/06/2018 22:28
E18-06141-08 [AA8207]	08/06/2018 23:02
E18-06141-09 [AA8208]	08/06/2018 23:35
10 PPBV CCCVS [AA8213CCCVS]	08/07/2018 2:22

Compound	CAS #	Reporting Limit (ppbv)	Concentration (ppbv)
Acetone	67-64-1	0.20	ND
Benzene	71-43-2	0.15	ND
Bromodichloromethane	75-27-4	0.20	ND
Bromoform	75-25-2	0.20	ND
Bromomethane	74-83-9	0.20	ND
1,3-Butadiene	106-99-0	0.20	ND
Chlorobenzene	108-90-7	0.20	ND
Chloroethane	75-00-3	0.20	ND
Chloroform	67-66-3	0.20	ND
Chloromethane	74-87-3	0.20	ND
Carbon disulfide	75-15-0	0.20	ND
Carbon tetrachloride	56-23-5	0.20	ND
Cyclohexane	110-82-7	0.12	ND
Dibromochloromethane	124-48-1	0.20	ND
1,2-Dibromoethane	106-93-4	0.20	ND
1,2-Dichlorobenzene	95-50-1	0.20	ND
1,3-Dichlorobenzene	541-73-1	0.20	ND
1,4-Dichlorobenzene	106-46-7	0.20	ND
Dichlorodifluoromethane	75-71-8	0.20	ND
1,1-Dichloroethane	75-34-3	0.20	ND
1,2-Dichloroethane	107-06-2	0.20	ND
1,1-Dichloroethene	75-35-4	0.20	ND
1,2-Dichloroethene (cis)	156-59-2	0.19	ND
1,2-Dichloroethene (trans)	156-60-5	0.20	ND
1,2-Dichloropropane	78-87-5	0.20	ND
1,3-Dichloropropene (cis)	10061-01-5	0.19	ND
1,3-Dichloropropene (trans)	10061-02-6	0.18	ND
1,2-Dichlorotetrafluoroethane	76-14-2	0.20	ND
1,4-Dioxane	123-91-1	0.17	ND
Ethylbenzene	100-41-4	0.20	ND
n-Heptane	142-82-5	0.20	ND
1,3-Hexachlorobutadiene	87-68-3	0.20	ND
n-Hexane	110-54-3	0.20	ND
Methylene chloride	75-09-2	0.20	ND

Method Blank must be less than the Practical Quantitation Limit (PQL).

Method Blank Report

Lab Sample Name: METHOD BLANK
Field Sample Name: METHOD BLANK
Matrix: Air
Dilution Factor: 1

Data File: AA8193BLK
Date Analyzed: 8/6/2018
Sample Volume: 500ml
GC/MS Column: RTX-1, 0.32 mmID

Runs with this Method Blank:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA8191BFB]	08/06/2018 11:26
10 PPBV DCVS [AA8192DCVS]	08/06/2018 13:07
METHOD BLANK [AA8193BLK]	08/06/2018 13:42
02 PPBV RLLCS [AA8194RLLCS]	08/06/2018 14:23
E18-06141-07 [AA8206]	08/06/2018 22:28
E18-06141-08 [AA8207]	08/06/2018 23:02
E18-06141-09 [AA8208]	08/06/2018 23:35
10 PPBV CCCVS [AA8213CCCVS]	08/07/2018 2:22

Compound	CAS #	Reporting Limit (ppbv)	Concentration (ppbv)
Methyl ethyl ketone	78-93-3	0.13	ND
Methyl isobutyl ketone	108-10-1	0.20	ND
Methyl tert-butyl ether	1634-04-4	0.11	ND
Styrene	100-42-5	0.14	ND
Tert-butyl alcohol	75-65-0	0.18	ND
1,1,2,2-Tetrachloroethane	79-34-5	0.17	ND
Tetrachloroethene	127-18-4	0.20	ND
Toluene	108-88-3	0.17	ND
1,2,4-Trichlorobenzene	120-82-1	0.20	ND
1,1,1-Trichloroethane	71-55-6	0.20	ND
1,1,2-Trichloroethane	79-00-5	0.20	ND
Trichloroethene	79-01-6	0.20	ND
Trichlorofluoromethane	75-69-4	0.20	ND
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	0.20	ND
1,2,4-Trimethylbenzene	95-63-6	0.12	ND
1,3,5-Trimethylbenzene	108-67-8	0.17	ND
2,2,4-Trimethylpentane	540-84-1	0.20	ND
Vinyl bromide	593-60-2	0.20	ND
Vinyl chloride	75-01-4	0.20	ND
Xylenes (m&p)	179601-23-1	0.20	ND
Xylenes (o)	95-47-6	0.20	ND

Method Blank must be less than the Practical Quantitation Limit (PQL).

Method Blank Report

Lab Sample Name: METHOD BLANK
Field Sample Name: METHOD BLANK
Matrix: Air
Dilution Factor: 1

Data File: AA8223BLK
Date Analyzed: 8/7/2018
Sample Volume: 500ml
GC/MS Column: RTX-1, 0.32 mmID

Runs with this Method Blank:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA8221BFB]	08/07/2018 9:33
10 PPBV DCVS [AA8222DCVS]	08/07/2018 10:17
METHOD BLANK [AA8223BLK]	08/07/2018 11:01
02 PPBV RLLCS [AA8224RLLCS]	08/07/2018 11:47
E18-06141-08 [AA8232]	08/07/2018 16:57
E18-06141-09 [AA8233]	08/07/2018 17:31
10 PPBV CCCVS [AA8242CCCVS]	08/07/2018 22:31

Compound	CAS #	Reporting Limit (ppbv)	Concentration (ppbv)
Acetone	67-64-1	0.20	ND
Benzene	71-43-2	0.15	ND
Bromodichloromethane	75-27-4	0.20	ND
Bromoform	75-25-2	0.20	ND
Bromomethane	74-83-9	0.20	ND
1,3-Butadiene	106-99-0	0.20	ND
Chlorobenzene	108-90-7	0.20	ND
Chloroethane	75-00-3	0.20	ND
Chloroform	67-66-3	0.20	ND
Chloromethane	74-87-3	0.20	ND
Carbon disulfide	75-15-0	0.20	ND
Carbon tetrachloride	56-23-5	0.20	ND
Cyclohexane	110-82-7	0.12	ND
Dibromochloromethane	124-48-1	0.20	ND
1,2-Dibromoethane	106-93-4	0.20	ND
1,2-Dichlorobenzene	95-50-1	0.20	ND
1,3-Dichlorobenzene	541-73-1	0.20	ND
1,4-Dichlorobenzene	106-46-7	0.20	ND
Dichlorodifluoromethane	75-71-8	0.20	ND
1,1-Dichloroethane	75-34-3	0.20	ND
1,2-Dichloroethane	107-06-2	0.20	ND
1,1-Dichloroethene	75-35-4	0.20	ND
1,2-Dichloroethene (cis)	156-59-2	0.19	ND
1,2-Dichloroethene (trans)	156-60-5	0.20	ND
1,2-Dichloropropane	78-87-5	0.20	ND
1,3-Dichloropropene (cis)	10061-01-5	0.19	ND
1,3-Dichloropropene (trans)	10061-02-6	0.18	ND
1,2-Dichlorotetrafluoroethane	76-14-2	0.20	ND
1,4-Dioxane	123-91-1	0.17	ND
Ethylbenzene	100-41-4	0.20	ND
n-Heptane	142-82-5	0.20	ND
1,3-Hexachlorobutadiene	87-68-3	0.20	ND
n-Hexane	110-54-3	0.20	ND
Methylene chloride	75-09-2	0.20	ND
Methyl ethyl ketone	78-93-3	0.13	ND

Method Blank must be less than the Practical Quantitation Limit (PQL).

Method Blank Report

Lab Sample Name: METHOD BLANK
Field Sample Name: METHOD BLANK
Matrix: Air
Dilution Factor: 1

Data File: AA8223BLK
Date Analyzed: 8/7/2018
Sample Volume: 500ml
GC/MS Column: RTX-1, 0.32 mmID

Runs with this Method Blank:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA8221BFB]	08/07/2018 9:33
10 PPBV DCVS [AA8222DCVS]	08/07/2018 10:17
METHOD BLANK [AA8223BLK]	08/07/2018 11:01
02 PPBV RLLCS [AA8224RLLCS]	08/07/2018 11:47
E18-06141-08 [AA8232]	08/07/2018 16:57
E18-06141-09 [AA8233]	08/07/2018 17:31
10 PPBV CCCVS [AA8242CCCVS]	08/07/2018 22:31

Compound	CAS #	Reporting Limit (ppbv)	Concentration (ppbv)
Methyl isobutyl ketone	108-10-1	0.20	ND
Methyl tert-butyl ether	1634-04-4	0.11	ND
Styrene	100-42-5	0.14	ND
Tert-butyl alcohol	75-65-0	0.18	ND
1,1,2,2-Tetrachloroethane	79-34-5	0.17	ND
Tetrachloroethene	127-18-4	0.20	ND
Toluene	108-88-3	0.17	ND
1,2,4-Trichlorobenzene	120-82-1	0.20	ND
1,1,1-Trichloroethane	71-55-6	0.20	ND
1,1,2-Trichloroethane	79-00-5	0.20	ND
Trichloroethene	79-01-6	0.20	ND
Trichlorofluoromethane	75-69-4	0.20	ND
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	0.20	ND
1,2,4-Trimethylbenzene	95-63-6	0.12	ND
1,3,5-Trimethylbenzene	108-67-8	0.17	ND
2,2,4-Trimethylpentane	540-84-1	0.20	ND
Vinyl bromide	593-60-2	0.20	ND
Vinyl chloride	75-01-4	0.20	ND
Xylenes (m&p)	179601-23-1	0.20	ND
Xylenes (o)	95-47-6	0.20	ND

Method Blank must be less than the Practical Quantitation Limit (PQL).

Integrated Analytical Laboratories, LLC

Volatile Organic Compounds by EPA Method TO-15

Laboratory Sample Duplicate Report

SDG Number: E18-04189
 IAL Sample ID: E18-04189-01
 Matrix: Air
 Summa ID: 5096

Date Received: 6/8/18
 Date Analyzed: 6/13/18,6/13/18
 Lab Data File#: AA7478,AA7479
 Dilution Factor: 1
 Injection Volume: 500ml
 GC/MS Column: RTX-1, 0.32 mmID

Compound	CAS #	Sample E18-04189-01		Sample Dup E18-04189-21		Reporting Limits ppbv	RPD
		Concentration Reported ppbv	Q	Concentration Reported ppbv	Q		
Acetone	67-64-1	3.1		3.1		0.40	0.00%
Allyl Chloride	107-05-1		0.40 U		0.40 U	0.40	0.00%
Benzene	71-43-2		0.40 U		0.40 U	0.40	0.00%
Bromodichloromethane	75-27-4		0.40 U		0.40 U	0.40	0.00%
Bromoform	75-25-2		0.40 U		0.40 U	0.40	0.00%
Bromomethane	74-83-9		0.40 U		0.40 U	0.40	0.00%
1,3-Butadiene	106-99-0		0.40 U		0.40 U	0.40	0.00%
Chlorobenzene	108-90-7		0.40 U		0.40 U	0.40	0.00%
Chloroethane	75-00-3		0.40 U		0.40 U	0.40	0.00%
Chloroform	67-66-3		0.40 U		0.40 U	0.40	0.00%
Chloromethane	74-87-3		0.40 U		0.40 U	0.40	0.00%
Carbon disulfide	75-15-0		0.40 U		0.40 U	0.40	0.00%
Carbon tetrachloride	56-23-5		0.40 U		0.40 U	0.40	0.00%
2-Chlorotoluene	95-49-8		0.40 U		0.40 U	0.40	0.00%
Cyclohexane	110-82-7		0.40 U		0.40 U	0.40	0.00%
Dibromochloromethane	124-48-1		0.40 U		0.40 U	0.40	0.00%
1,2-Dibromoethane	106-93-4		0.40 U		0.40 U	0.40	0.00%
1,2-Dichlorobenzene	95-50-1		0.40 U		0.40 U	0.40	0.00%
1,3-Dichlorobenzene	541-73-1		0.40 U		0.40 U	0.40	0.00%
1,4-Dichlorobenzene	106-46-7		0.40 U		0.40 U	0.40	0.00%
Dichlorodifluoromethane	75-71-8		0.40 U		0.40 U	0.40	0.00%
1,1-Dichloroethane	75-34-3		0.40 U		0.40 U	0.40	0.00%
1,2-Dichloroethane	107-06-2		0.40 U		0.40 U	0.40	0.00%
1,1-Dichloroethene	75-35-4		0.40 U		0.40 U	0.40	0.00%
1,2-Dichloroethene (cis)	156-59-2		0.40 U		0.40 U	0.40	0.00%
1,2-Dichloroethene (trans)	156-60-5		0.40 U		0.40 U	0.40	0.00%
1,2-Dichloropropane	78-87-5		0.40 U		0.40 U	0.40	0.00%
1,3-Dichloropropene (cis)	10061-01-5		0.40 U		0.40 U	0.40	0.00%
1,3-Dichloropropene (trans)	10061-02-6		0.40 U		0.40 U	0.40	0.00%
1,2-Dichlorotetrafluoroethane	76-14-2		0.40 U		0.40 U	0.40	0.00%
Ethylbenzene	100-41-4		0.40 U		0.40 U	0.40	0.00%
4-Ethyltoluene	622-96-8		0.40 U		0.40 U	0.40	0.00%
n-Heptane	142-82-5		0.40 U		0.40 U	0.40	0.00%
1,3-Hexachlorobutadiene	87-68-3		0.40 U		0.40 U	0.40	0.00%
n-Hexane	110-54-3		0.40 U		0.40 U	0.40	0.00%
Methylene chloride	75-09-2	2.3		2.6		0.40	-12.24%
Methyl ethyl ketone	78-93-3		0.40 U		0.40 U	0.40	0.00%
Methyl isobutyl ketone	108-10-1		0.40 U		0.40 U	0.40	0.00%

Qualifir:

E=Concentration exceeds upper level of calibration range for instrument.

D=Extra dilution required for this compound. J=Duplicate samples do not met RPD criteria.

U=Compound ND or under reporting limit.

Integrated Analytical Laboratories, LLC

Volatile Organic Compounds by EPA Method TO-15

Laboratory Sample Duplicate Report

SDG Number: E18-04189
 IAL Sample ID: E18-04189-01
 Matrix: Air
 Summa ID: 5096

Date Received: 6/8/18
 Date Analyzed: 6/13/18,6/13/18
 Lab Data File#: AA7478,AA7479
 Dilution Factor: 1
 Injection Volume: 500ml
 GC/MS Column: RTX-1, 0.32 mmID

Compound	CAS #	Sample E18-04189-01 Concentration Reported		Sample Dup E18-04189-21 Concentration Reported		Reporting Limits ppbv	RPD
		ppbv	Q	ppbv	Q		
Methyl tert-butyl ether	1634-04-4	0.40	U	0.40	U	0.40	0.00%
Styrene	100-42-5	0.40	U	0.40	U	0.40	0.00%
Tert-butyl alcohol	75-65-0	0.40	U	0.40	U	0.40	0.00%
1,1,2,2-Tetrachloroethane	79-34-5	0.40	U	0.40	U	0.40	0.00%
Tetrachloroethene	127-18-4	0.40	U	0.40	U	0.40	0.00%
Toluene	108-88-3	0.40	U	0.40	U	0.40	0.00%
1,2,4-Trichlorobenzene	120-82-1	0.40	U	0.40	U	0.40	0.00%
1,1,1-Trichloroethane	71-55-6	0.40	U	0.40	U	0.40	0.00%
1,1,2-Trichloroethane	79-00-5	0.40	U	0.40	U	0.40	0.00%
Trichloroethene	79-01-6	0.40	U	0.40	U	0.40	0.00%
Trichlorofluoromethane	75-69-4	0.40	U	0.40	U	0.40	0.00%
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	0.40	U	0.40	U	0.40	0.00%
1,2,4-Trimethylbenzene	95-63-6	0.40	U	0.40	U	0.40	0.00%
1,3,5-Trimethylbenzene	108-67-8	0.40	U	0.40	U	0.40	0.00%
2,2,4-Trimethylpentane	540-84-1	0.40	U	0.40	U	0.40	0.00%
Vinyl bromide	593-60-2	0.40	U	0.40	U	0.40	0.00%
Vinyl chloride	75-01-4	0.40	U	0.40	U	0.40	0.00%
Xylenes (m&p)	179601-23-1	0.40	U	0.40	U	0.40	0.00%
Xylenes (o)	95-47-6	0.40	U	0.40	U	0.40	0.00%

RPD must be <25% for all laboratory duplicate samples. Laboratory duplicate samples are run once daily.
NC = The RPD could not be calculated since the compound was only detected in either the parent or duplicate sample.

Qualifir:

E=Concentration exceeds upper level of calibration range for instrument.
 D=Extra dilution required for this compound. J=Duplicate samples do not met RPD criteria.
 U=Compound ND or under reporting limit.

Integrated Analytical Laboratories, LLC

Volatile Organic Compounds by EPA Method TO-15

Laboratory Sample Duplicate Report

SDG Number: E18-06204
 IAL Sample ID: E18-06204-01
 Matrix: Air
 Summa ID: 3280

Date Received: 8/6/18
 Date Analyzed: 8/6/18,8/6/18
 Lab Data File#: AA8195,AA8196
 Dilution Factor: 1
 Injection Volume: 500ml
 GC/MS Column: RTX-1, 0.32 mmID

Compound	CAS #	Sample E18-06204-01 Concentration Reported		Sample Dup E18-06204-21 Concentration Reported		Reporting Limits ppbv	RPD
		ppbv	Q	ppbv	Q		
Acetone	67-64-1	10		10.0		0.20	0.00%
Allyl Chloride	107-05-1		0.20 U		0.20 U	0.20	0.00%
Benzene	71-43-2	0.26		0.24		0.20	8.00%
Bromodichloromethane	75-27-4		0.20 U		0.20 U	0.20	0.00%
Bromoform	75-25-2		0.20 U		0.20 U	0.20	0.00%
Bromomethane	74-83-9		0.20 U		0.20 U	0.20	0.00%
1,3-Butadiene	106-99-0		0.20 U		0.20 U	0.20	0.00%
Chlorobenzene	108-90-7		0.20 U		0.20 U	0.20	0.00%
Chloroethane	75-00-3		0.20 U		0.20 U	0.20	0.00%
Chloroform	67-66-3		0.20 U		0.20 U	0.20	0.00%
Chloromethane	74-87-3		0.20 U		0.20 U	0.20	0.00%
Carbon disulfide	75-15-0		0.20 U		0.20 U	0.20	0.00%
Carbon tetrachloride	56-23-5		0.20 U		0.20 U	0.20	0.00%
2-Chlorotoluene	95-49-8		0.20 U		0.20 U	0.20	0.00%
Cyclohexane	110-82-7		0.20 U		0.20 U	0.20	0.00%
Dibromochloromethane	124-48-1		0.20 U		0.20 U	0.20	0.00%
1,2-Dibromoethane	106-93-4		0.20 U		0.20 U	0.20	0.00%
1,2-Dichlorobenzene	95-50-1		0.20 U		0.20 U	0.20	0.00%
1,3-Dichlorobenzene	541-73-1		0.20 U		0.20 U	0.20	0.00%
1,4-Dichlorobenzene	106-46-7		0.20 U		0.20 U	0.20	0.00%
Dichlorodifluoromethane	75-71-8		0.20 U		0.20 U	0.20	0.00%
1,1-Dichloroethane	75-34-3		0.20 U		0.20 U	0.20	0.00%
1,2-Dichloroethane	107-06-2		0.20 U		0.20 U	0.20	0.00%
1,1-Dichloroethene	75-35-4		0.20 U		0.20 U	0.20	0.00%
1,2-Dichloroethene (cis)	156-59-2		0.20 U		0.20 U	0.20	0.00%
1,2-Dichloroethene (trans)	156-60-5		0.20 U		0.20 U	0.20	0.00%
1,2-Dichloropropane	78-87-5		0.20 U		0.20 U	0.20	0.00%
1,3-Dichloropropene (cis)	10061-01-5		0.20 U		0.20 U	0.20	0.00%
1,3-Dichloropropene (trans)	10061-02-6		0.20 U		0.20 U	0.20	0.00%
1,2-Dichlorotetrafluoroethane	76-14-2		0.20 U		0.20 U	0.20	0.00%
1,4-Dioxane	123-91-1		0.20 U		0.20 U	0.20	0.00%
Ethanol	64-17-5	84	E	80	E	0.20	4.88%
Ethylbenzene	100-41-4		0.20 U		0.20 U	0.20	0.00%
4-Ethyltoluene	622-96-8		0.20 U		0.20 U	0.20	0.00%
n-Heptane	142-82-5		0.20 U		0.20 U	0.20	0.00%
1,3-Hexachlorobutadiene	87-68-3		0.20 U		0.20 U	0.20	0.00%
n-Hexane	110-54-3	0.20		0.21		0.20	-4.88%
Isopropanol	67-63-0	1.2		1.2		0.20	0.00%

Qualifir:

E=Concentration exceeds upper level of calibration range for instrument.
 D=Extra dilution required for this compound. J=Duplicate samples do not met RPD criteria.
 U=Compound ND or under reporting limit.

Integrated Analytical Laboratories, LLC

Volatile Organic Compounds by EPA Method TO-15

Laboratory Sample Duplicate Report

SDG Number: E18-06204
 IAL Sample ID: E18-06204-01
 Matrix: Air
 Summa ID: 3280

Date Received: 8/6/18
 Date Analyzed: 8/6/18,8/6/18
 Lab Data File#: AA8195,AA8196
 Dilution Factor: 1
 Injection Volume: 500ml
 GC/MS Column: RTX-1, 0.32 mmID

Compound	CAS #	Sample E18-06204-01 Concentration Reported		Sample Dup E18-06204-21 Concentration Reported		Reporting Limits ppbv	RPD
		ppbv	Q	ppbv	Q		
Methylene chloride	75-09-2	1.5		1.4		0.20	6.90%
Methyl ethyl ketone	78-93-3	1.7		1.8		0.20	-5.71%
Methyl isobutyl ketone	108-10-1		0.20 U		0.20 U	0.20	0.00%
Methyl methacrylate	80-62-6		0.20 U		0.20 U	0.20	0.00%
Methyl tert-butyl ether	1634-04-4		0.20 U		0.20 U	0.20	0.00%
Styrene	100-42-5		0.20 U		0.20 U	0.20	0.00%
Tert-butyl alcohol	75-65-0		0.20 U		0.20 U	0.20	0.00%
1,1,2,2-Tetrachloroethane	79-34-5		0.20 U		0.20 U	0.20	0.00%
Tetrachloroethene	127-18-4		0.20 U		0.20 U	0.20	0.00%
Tetrahydrofuran	109-99-9		0.20 U		0.20 U	0.20	0.00%
Toluene	108-88-3	0.23		0.23		0.20	0.00%
1,2,4-Trichlorobenzene	120-82-1		0.20 U		0.20 U	0.20	0.00%
1,1,1-Trichloroethane	71-55-6		0.20 U		0.20 U	0.20	0.00%
1,1,2-Trichloroethane	79-00-5		0.20 U		0.20 U	0.20	0.00%
Trichloroethene	79-01-6		0.20 U		0.20 U	0.20	0.00%
Trichlorofluoromethane	75-69-4	0.57		0.55		0.20	3.57%
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1		0.20 U		0.20 U	0.20	0.00%
1,2,4-Trimethylbenzene	95-63-6		0.20 U		0.20 U	0.20	0.00%
1,3,5-Trimethylbenzene	108-67-8		0.20 U		0.20 U	0.20	0.00%
2,2,4-Trimethylpentane	540-84-1		0.20 U		0.20 U	0.20	0.00%
Vinyl bromide	593-60-2		0.20 U		0.20 U	0.20	0.00%
Vinyl chloride	75-01-4		0.20 U		0.20 U	0.20	0.00%
Xylenes (m&p)	179601-23-1		0.40 U		0.40 U	0.40	0.00%
Xylenes (o)	95-47-6		0.20 U		0.20 U	0.20	0.00%

RPD must be <25% for all laboratory duplicate samples. Laboratory duplicate samples are run once daily.

NC = The RPD could not be calculated since the compound was only detected in either the parent or duplicate sample.

Qualifir:

E=Concentration exceeds upper level of calibration range for instrument.

D=Extra dilution required for this compound. J=Duplicate samples do not met RPD criteria.

U=Compound ND or under reporting limit.

Integrated Analytical Laboratories, LLC

Volatile Organic Compounds by EPA Method TO-15

Laboratory Sample Duplicate Report

SDG Number: E18-06173
 IAL Sample ID: E18-06173-01
 Matrix: Air
 Summa ID: 1070

Date Received: 8/3/18
 Date Analyzed: 8/7/18,8/7/18
 Lab Data File#: AA8228,AA8229
 Dilution Factor: 1
 Injection Volume: 50ml
 GC/MS Column: RTX-1, 0.32 mmID

Compound	CAS #	Sample E18-06173-01 Concentration Reported		Sample Dup E18-06173-21 Concentration Reported		Reporting Limits ppbv	RPD
		ppbv	Q	ppbv	Q		
Acetone	67-64-1	30		27		2.0	10.53%
Allyl Chloride	107-05-1		2.0 U		2.0 U	2.0	0.00%
Benzene	71-43-2		2.0 U		2.0 U	2.0	0.00%
Bromodichloromethane	75-27-4		2.0 U		2.0 U	2.0	0.00%
Bromoform	75-25-2		2.0 U		2.0 U	2.0	0.00%
Bromomethane	74-83-9		2.0 U		2.0 U	2.0	0.00%
1,3-Butadiene	106-99-0		2.0 U		2.0 U	2.0	0.00%
Chlorobenzene	108-90-7		2.0 U		2.0 U	2.0	0.00%
Chloroethane	75-00-3		2.0 U		2.0 U	2.0	0.00%
Chloroform	67-66-3		2.0 U		2.0 U	2.0	0.00%
Chloromethane	74-87-3		2.0 U		2.0 U	2.0	0.00%
Carbon disulfide	75-15-0		2.0 U		2.0 U	2.0	0.00%
Carbon tetrachloride	56-23-5		2.0 U		2.0 U	2.0	0.00%
2-Chlorotoluene	95-49-8		2.0 U		2.0 U	2.0	0.00%
Cyclohexane	110-82-7		2.0 U		2.0 U	2.0	0.00%
Dibromochloromethane	124-48-1		2.0 U		2.0 U	2.0	0.00%
1,2-Dibromoethane	106-93-4		2.0 U		2.0 U	2.0	0.00%
1,2-Dichlorobenzene	95-50-1		2.0 U		2.0 U	2.0	0.00%
1,3-Dichlorobenzene	541-73-1		2.0 U		2.0 U	2.0	0.00%
1,4-Dichlorobenzene	106-46-7		2.0 U		2.0 U	2.0	0.00%
Dichlorodifluoromethane	75-71-8		2.0 U		2.0 U	2.0	0.00%
1,1-Dichloroethane	75-34-3		2.0 U		2.0 U	2.0	0.00%
1,2-Dichloroethane	107-06-2		2.0 U		2.0 U	2.0	0.00%
1,1-Dichloroethene	75-35-4		2.0 U		2.0 U	2.0	0.00%
1,2-Dichloroethene (cis)	156-59-2		2.0 U		2.0 U	2.0	0.00%
1,2-Dichloroethene (trans)	156-60-5		2.0 U		2.0 U	2.0	0.00%
1,2-Dichloropropane	78-87-5		2.0 U		2.0 U	2.0	0.00%
1,3-Dichloropropene (cis)	10061-01-5		2.0 U		2.0 U	2.0	0.00%
1,3-Dichloropropene (trans)	10061-02-6		2.0 U		2.0 U	2.0	0.00%
1,2-Dichlorotetrafluoroethane	76-14-2		2.0 U		2.0 U	2.0	0.00%
1,4-Dioxane	123-91-1		2.0 U		2.0 U	2.0	0.00%
Ethanol	64-17-5	380		370		2.0	2.67%
Ethylbenzene	100-41-4		2.0 U		2.0 U	2.0	0.00%
4-Ethyltoluene	622-96-8		2.0 U		2.0 U	2.0	0.00%
n-Heptane	142-82-5		2.0 U		2.0 U	2.0	0.00%
1,3-Hexachlorobutadiene	87-68-3		2.0 U		2.0 U	2.0	0.00%
n-Hexane	110-54-3		2.0 U		2.0 U	2.0	0.00%
Isopropanol	67-63-0	5.2		5.5		2.0	-5.61%

Qualifir:

E=Concentration exceeds upper level of calibration range for instrument.

D=Extra dilution required for this compound. J=Duplicate samples do not met RPD criteria.

U=Compound ND or under reporting limit.

Integrated Analytical Laboratories, LLC

Volatile Organic Compounds by EPA Method TO-15

Laboratory Sample Duplicate Report

SDG Number: E18-06173
 IAL Sample ID: E18-06173-01
 Matrix: Air
 Summa ID: 1070

Date Received: 8/3/18
 Date Analyzed: 8/7/18,8/7/18
 Lab Data File#: AA8228,AA8229
 Dilution Factor: 1
 Injection Volume: 50ml
 GC/MS Column: RTX-1, 0.32 mmID

Compound	CAS #	Sample E18-06173-01 Concentration Reported		Sample Dup E18-06173-21 Concentration Reported		Reporting Limits ppbv	RPD
		ppbv	Q	ppbv	Q		
Methylene chloride	75-09-2	99		100		2.0	-1.01%
Methyl ethyl ketone	78-93-3	2.8		2.2		2.0	24.00%
Methyl isobutyl ketone	108-10-1		2.0 U		2.0 U	2.0	0.00%
Methyl methacrylate	80-62-6		2.0 U		2.0 U	2.0	0.00%
Methyl tert-butyl ether	1634-04-4		2.0 U		2.0 U	2.0	0.00%
Styrene	100-42-5		2.0 U		2.0 U	2.0	0.00%
Tert-butyl alcohol	75-65-0	3.1		3.0		2.0	3.28%
1,1,2,2-Tetrachloroethane	79-34-5		2.0 U		2.0 U	2.0	0.00%
Tetrachloroethene	127-18-4		2.0 U		2.0 U	2.0	0.00%
Tetrahydrofuran	109-99-9		2.0 U		2.0 U	2.0	0.00%
Toluene	108-88-3	2.7		2.7		2.0	0.00%
1,2,4-Trichlorobenzene	120-82-1		2.0 U		2.0 U	2.0	0.00%
1,1,1-Trichloroethane	71-55-6		2.0 U		2.0 U	2.0	0.00%
1,1,2-Trichloroethane	79-00-5		2.0 U		2.0 U	2.0	0.00%
Trichloroethene	79-01-6		2.0 U		2.0 U	2.0	0.00%
Trichlorofluoromethane	75-69-4		2.0 U		2.0 U	2.0	0.00%
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1		2.0 U		2.0 U	2.0	0.00%
1,2,4-Trimethylbenzene	95-63-6		2.0 U		2.0 U	2.0	0.00%
1,3,5-Trimethylbenzene	108-67-8		2.0 U		2.0 U	2.0	0.00%
2,2,4-Trimethylpentane	540-84-1		2.0 U		2.0 U	2.0	0.00%
Vinyl bromide	593-60-2		2.0 U		2.0 U	2.0	0.00%
Vinyl chloride	75-01-4		2.0 U		2.0 U	2.0	0.00%
Xylenes (m&p)	179601-23-1		4.0 U		4.0 U	4.0	0.00%
Xylenes (o)	95-47-6		2.0 U		2.0 U	2.0	0.00%

RPD must be <25% for all laboratory duplicate samples. Laboratory duplicate samples are run once daily.

NC = The RPD could not be calculated since the compound was only detected in either the parent or duplicate sample.

Qualifir:

E=Concentration exceeds upper level of calibration range for instrument.
 D=Extra dilution required for this compound. J=Duplicate samples do not met RPD criteria.
 U=Compound ND or under reporting limit.

Initial Calibration Curve Internal Standard Area and Retention Time Summary

Instrument: AA

ICAL Date: 5/18/2018

		BROMOCHLOROMETHANE				1,4-DIFLUOROBENZENE				D-5 CHLOROBENZENE			
		Area #	RT			Area #	RT			Area #	RT		
AVERAGE OF CALIBRATION STANDARDS		534003	7.733			2218829	9.782			1956425	15.111		
UPPER LIMIT		747604	8.063			3106360	10.112			2738995	15.441		
LOWER LIMIT		320402	7.403			1331297	9.452			1173855	14.781		
Lab ID		Area #	%	RT	+/-	Area #	%	RT	+/-	Area #	%	RT	+/-
40 PPBV STD	AA 7073 STD01	497626	6.81	7.759	0.03	2152735	2.98	9.788	0.01	1991801	1.81	15.113	0.00
2 PPBV STD	AA 7074 STD04	531970	0.38	7.740	0.01	2238046	0.87	9.785	0.00	2008933	2.68	15.112	0.00
10 PPBV STD	AA 7075 STD03	543623	1.80	7.724	0.01	2258745	1.80	9.782	0.00	1967114	0.55	15.109	0.00
2 PPBV STD	AA 7076 STD04	531646	0.44	7.717	0.02	2168184	2.28	9.775	0.01	1841018	5.90	15.112	0.00
0.2 PPBV STD	AA 7077 STD05	565151	5.83	7.724	0.01	2276433	2.60	9.778	0.00	1973260	0.86	15.109	0.00
ICVSS	AA 7078 ICVSS	522049	2.24	7.724	0.01	2190215	1.29	9.782	0.00	1913238	2.21	15.113	0.00

Difference of Internal Area must be within +/- 40%; Retention Times must be within +/- 0.33 minute.

* Values outside QC limits.

Internal Standard Area and Retention Time Summary

Lab File ID (Standard): AA7472DCVS

Date Analyzed: 6/13/2018

Instrument: AA

ICAL Date: 5/18/2018

			BROMOCHLOROMETHANE				1,4-DIFLUOROBENZENE				D-5 CHLORO BENZENE				
CALIBRATION STANDARD			Area #	RT				Area #	RT				Area #	RT	
UPPER LIMIT			539798	7.724				2150341	9.785				1773674	15.116	
LOWER LIMIT			755717	8.05				3010477	10.12				2483144	15.45	
			323879	7.39				1290205	9.46				1064204	14.79	
Lab ID	DF		Area #	%	RT	+/-	Area #	%	RT	+/-	Area #	%	RT	+/-	
Method Blank	AA7473BLK	1.0	510475	-5.43	7.724	0.00	1863647	-13.33	9.785	0.00	1558649	-12.12	15.116	0.00	
Reporting Limit Laboratory Control Standard	AA7474RLLCS	1.0	558319	3.43	7.720	0.00	2000895	-6.95	9.781	0.00	1704889	-3.88	15.116	0.00	
3059	AA7475	1.0	486608	-9.85	7.724	0.00	1675394	-22.09	9.785	0.00	1426411	-19.58	15.116	0.00	
Closing Calibration	AA7485CCCVS	1.0	512140	-5.12	7.727	0.00	1948511	-9.39	9.785	0.00	1565782	-11.72	15.116	0.00	

Difference of Internal Area must be within +/- 40%; Retention Times must be within +/- 0.33 minute.

* Values outside QC limits.

Initial Calibration Curve Internal Standard Area and Retention Time Summary

Instrument: AA

ICAL Date: 7/25/2018

		BROMOCHLOROMETHANE				1,4-DIFLUOROBENZENE				D-5 CHLOROBENZENE			
		Area #	RT			Area #	RT			Area #	RT		
AVERAGE OF CALIBRATION STANDARDS		476007	7.689			2043575	9.740			1757787	15.057		
UPPER LIMIT		666410	8.019			2861004	10.070			2460901	15.387		
LOWER LIMIT		285604	7.359			1226145	9.410			1054672	14.727		
Lab ID		Area #	%	RT	+/-	Area #	%	RT	+/-	Area #	%	RT	+/-
40 PPBV STD	AA 7972 STD01	459428	3.48	7.702	0.01	2108198	3.16	9.743	0.00	1916771	9.04	15.058	0.00
20 PPBV STD	AA 7973 STD04	486694	2.25	7.695	0.01	2128550	4.16	9.744	0.00	1865915	6.15	15.058	0.00
10 PPBV STD	AA 7974 STD03	485778	2.05	7.689	0.00	2080600	1.81	9.740	0.00	1771224	0.76	15.058	0.00
2 PPBV STD	AA 7975 STD04	442159	7.11	7.679	0.01	1836044	10.16	9.737	0.00	1531011	12.90	15.055	0.00
0.2 PPBV STD	AA 7976 STD05	505975	6.30	7.679	0.01	2064481	1.02	9.734	0.01	1704012	3.06	15.055	0.00
ICVSS	AA 7977 ICVSS	434360	8.75	7.686	0.00	1814996	11.19	9.734	0.01	1540277	12.37	15.055	0.00

Difference of Internal Area must be within +/- 40%; Retention Times must be within +/- 0.33 minute.

* Values outside QC limits.

Internal Standard Area and Retention Time Summary

Lab File ID (Standard): AA8192DCVS

Date Analyzed: 8/6/2018

Instrument: AA

ICAL Date: 7/25/2018

			BROMOCHLOROMETHANE				1,4-DIFLUOROBENZENE				D-5 CHLORO BENZENE				
CALIBRATION STANDARD			Area #	RT				Area #	RT				Area #	RT	
UPPER LIMIT			438428	7.686				1771164	9.734				1475100	15.049	
LOWER LIMIT			613799	8.02				2479630	10.06				2065140	15.38	
			263057	7.36				1062698	9.40				885060	14.72	
Lab ID	DF		Area #	%	RT	+/-	Area #	%	RT	+/-	Area #	%	RT	+/-	
Method Blank	AA8193BLK	1.0	418368	-4.58	7.679	-0.01	1550382	-12.47	9.731	0.00	1251713	-15.14	15.052	0.00	
Reporting Limit Laboratory Control Standard	AA8194RLLCS	1.0	423253	-3.46	7.679	-0.01	1605091	-9.38	9.731	0.00	1341071	-9.09	15.052	0.00	
E18-06141-07	AA8206	1.0	352575	-19.58	7.724	0.04	1467744	-17.13	9.747	0.01	1402127	-4.95	15.058	0.01	
E18-06141-08	AA8207	1.0	485982	10.85	7.731	0.05	1975065	11.51	9.747	0.01	1713875	16.19	15.058	0.01	
E18-06141-09	AA8208	1.0	516268	17.75	7.721	0.04	2306557	30.23	9.747	0.01	1996784	35.37	15.058	0.01	
Closing Calibration	AA8213CCCVS	1.0	386051	-11.95	7.682	0.00	1532831	-13.46	9.737	0.00	1278044	-13.36	15.052	0.00	

Difference of Internal Area must be within +/- 40%; Retention Times must be within +/- 0.33 minute.

* Values outside QC limits.

Internal Standard Area and Retention Time Summary

Lab File ID (Standard): AA8222DCVS

Date Analyzed: 8/7/2018

Instrument: AA

ICAL Date: 7/25/2018

			BROMOCHLOROMETHANE				1,4-DIFLUOROBENZENE				D-5 CHLORO BENZENE				
CALIBRATION STANDARD			Area #	RT				Area #	RT				Area #	RT	
UPPER LIMIT			404488	7.682				1591121	9.734				1304098	15.049	
LOWER LIMIT			566283	8.01				2227569	10.06				1825737	15.38	
			242693	7.35				954673	9.40				782459	14.72	
Lab ID	DF		Area #	%	RT	+/-	Area #	%	RT	+/-	Area #	%	RT	+/-	
Method Blank	AA8223BLK	1.0	386755	-4.38	7.682	0.00	1385734	-12.91	9.734	0.00	1124467	-13.77	15.052	0.00	
Reporting Limit Laboratory Control Standard	AA8224RLLCS	1.0	398730	-1.42	7.676	-0.01	1455932	-8.50	9.730	0.00	1218171	-6.59	15.055	0.01	
E18-06141-08	AA8232	100.0	332191	-17.87	7.682	0.00	1080070	-32.12	9.737	0.00	922301	-29.28	15.052	0.00	
E18-06141-09	AA8233	10.0	305198	-24.55	7.692	0.01	1178297	-25.95	9.743	0.01	1041852	-20.11	15.055	0.01	
Closing Calibration	AA8242CCCVS	1.0	316732	-21.70	7.682	0.00	1419983	-10.76	9.734	0.00	1112022	-14.73	15.055	0.01	

Difference of Internal Area must be within +/- 40%; Retention Times must be within +/- 0.33 minute.

* Values outside QC limits.

Section VI: Sample Data Summary

Certificate of Analysis

Summary of Results

**Quantitation Reports, Chromatograms,
and Peak Integration Reports**

CERTIFICATE OF ANALYSIS

**ANALYTICAL DATA PACKAGE FOR THE
NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION
ALBANY NEW YORK 12233**

Integrated Analytical Laboratories, LLC
Project#: Congers CP / 060141
SDG #: E18-06141
Date of first sample receipt: 8/2/2018

Randolph, NJ 07869
NY ELAP Certification#: 11402
NJDEP (Primary AB) Certification#: 14751
Date of last sample receipt: 8/2/2018

Client: Brennan Environmental
19 Chatham Road
Summit, NJ 07901

Attention: Attention: Jeff McCurdy

Project/Site: Congers CP / 060141/NY

Analysis conducted at: Integrated Analytical Laboratories, LLC
273 Franklin Road
Randolph, NJ 07869

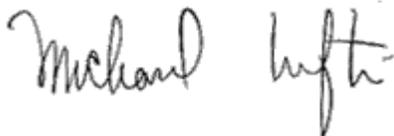
Contact: Michael H. Leftin, Ph.D.

Sample(s):

E18-06141-07
E18-06141-08
E18-06141-09

Samples for this analysis were received in good condition with a chain of custody.

All work recorded herein has been done in accordance with normal professional standards using accepted testing methodologies, quality assurance and quality control procedures except where otherwise agreed to by the client and testing company in writing. Once analysis has been performed on canisters that meets regulatory criteria, samples are recycled for future use, unless other provisions have been made by the client.



Michael H. Leftin, Ph.D.
Laboratory Director

Date: August 27, 2018



Integrated Analytical Laboratories LLC

Summary of Results

Brennan Environmental
 19 Chatham Road
 Summit, NJ 07901
 Attn: Jeff McCurdy
 Project: Congers CP / 060141
 Site: NY

Report Date: 08/13/18
 SDG Number: E18-06141
 Date Sampled: 08/01/18
 Date Received: 08/02/18
 Date Analyzed: 08/06/18
 Data File: AA8206
 Summa ID: 3014A
 DF: 1

Analysis: Volatile Organic Compounds by EPA Method TO-15

Compound	CAS #	Sample Name: SS-101		Reporting Limits	
		IAL ID: E18-06141-07		ppbv	ug/m3
1,1-Dichloroethane	75-34-3	ND	ND	0.20	0.81
1,2-Dichloroethane	107-06-2	ND	ND	0.20	0.81
1,1-Dichloroethene	75-35-4	ND	ND	0.20	0.79
1,2-Dichloroethene (cis)	156-59-2	ND	ND	0.20	0.79
1,2-Dichloroethene (trans)	156-60-5	ND	ND	0.20	0.79
1,1,2,2-Tetrachloroethane	79-34-5	ND	ND	0.20	1.4
Tetrachloroethene	127-18-4	1.5	10	0.20	1.4
1,1,1-Trichloroethane	71-55-6	ND	ND	0.20	1.1
1,1,2-Trichloroethane	79-00-5	ND	ND	0.20	1.1
Trichloroethene	79-01-6	ND	ND	0.05	0.25
Vinyl chloride	75-01-4	ND	ND	0.20	0.51

Data Path : C:\DATA\08-06-18\
 Data File : aa8206.D
 Acq On : 6 Aug 2018 10:28 pm
 Operator : jls
 Sample : E18-06141-07
 Misc : 3014A, 500cc
 ALS Vial : 16 Sample Multiplier: 1

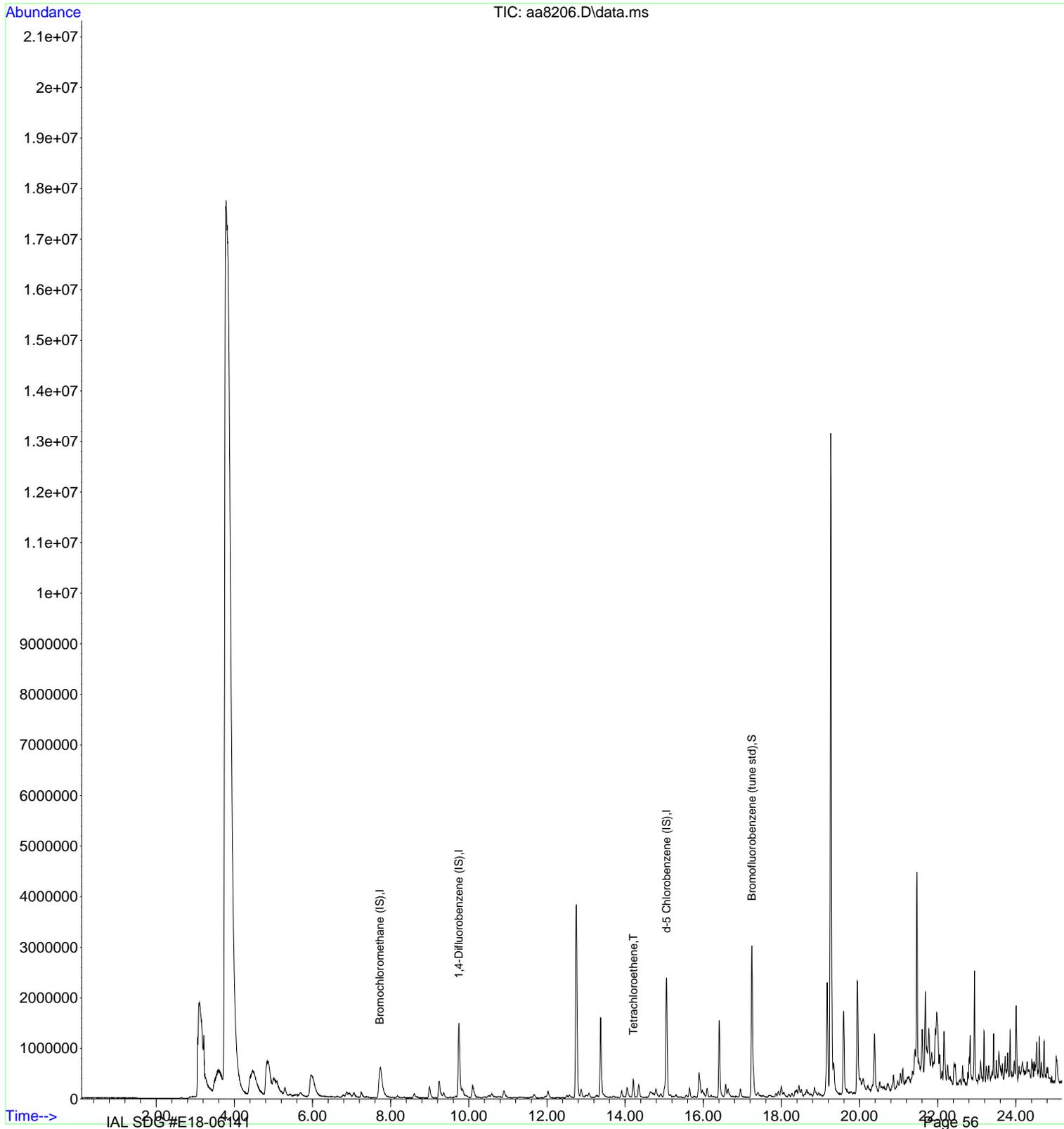
Quant Time: Aug 07 11:27:02 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

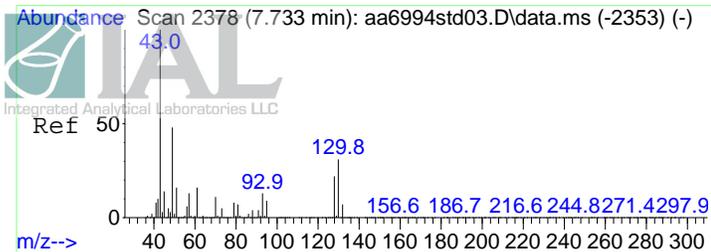
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Bromochloromethane (IS)	7.724	130	352575	10.00	ppbV	0.04
38) 1,4-Difluorobenzene (IS)	9.747	114	1467744	10.00	ppbV #	0.00
55) d-5 Chlorobenzene (IS)	15.058	117	1402127	10.00	ppbV	0.00
System Monitoring Compounds						
64) Bromofluorobenzene (tu...	17.245	95	1320113	10.32	ppbV	0.00
Target Compounds						
54) Tetrachloroethene	14.213	166	102997	1.53	ppbV	Qvalue 99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\08-06-18\
 Data File : aa8206.D
 Acq On : 6 Aug 2018 10:28 pm
 Operator : jls
 Sample : E18-06141-07
 Misc : 3014A, 500cc
 ALS Vial : 16 Sample Multiplier: 1

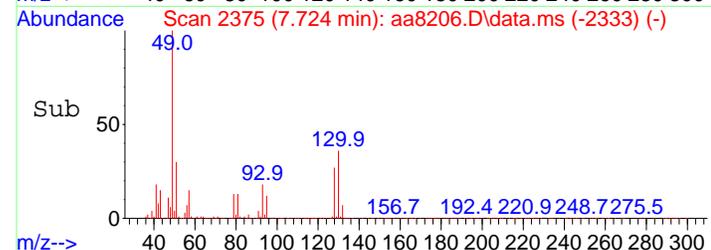
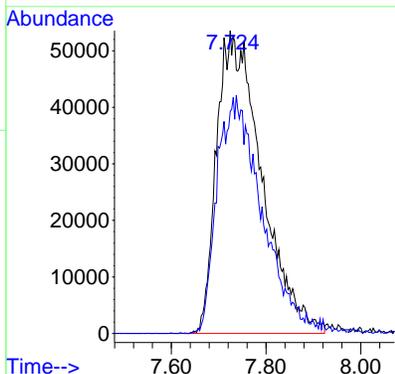
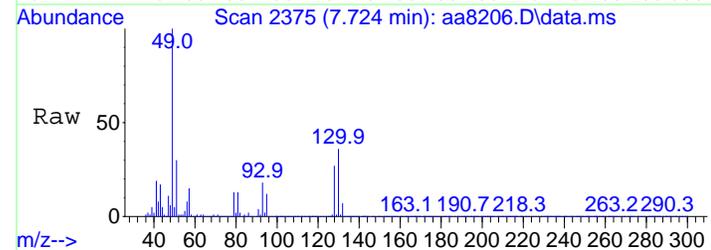
Quant Time: Aug 07 11:27:02 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration





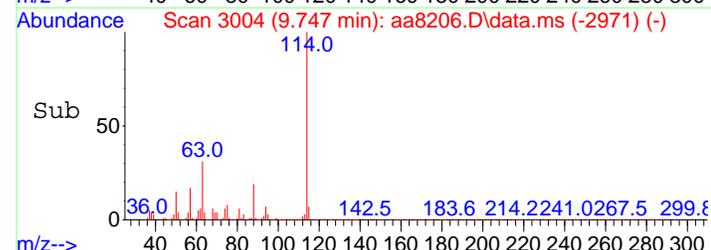
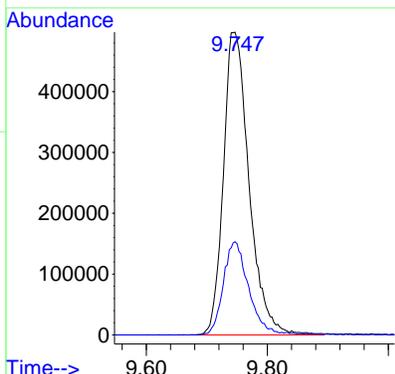
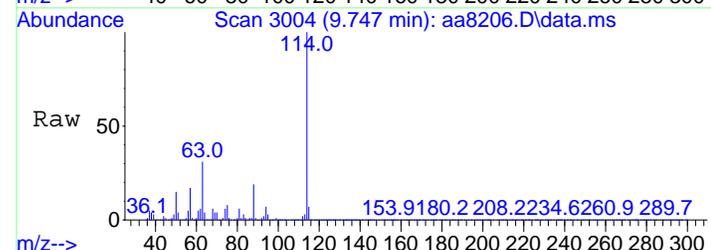
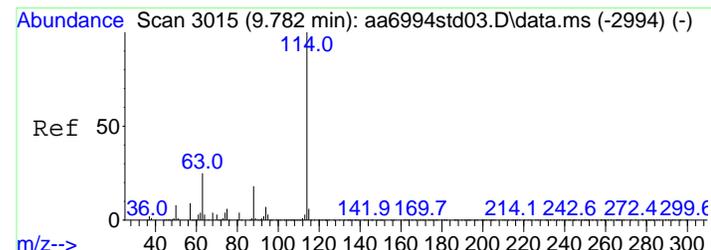
#1
 Bromochloromethane (IS)
 Concen: 10.00 ppbV
 RT: 7.724 min Scan# 2375
 Delta R.T. 0.035 min
 Lab File: aa8206.D
 Acq: 6 Aug 2018 10:28 pm

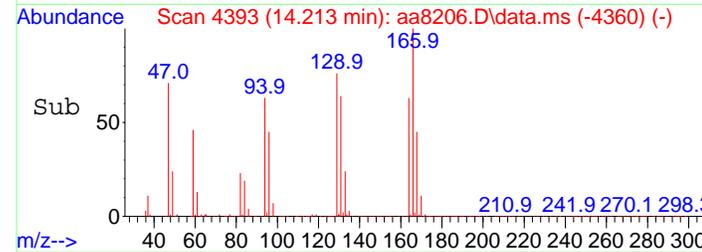
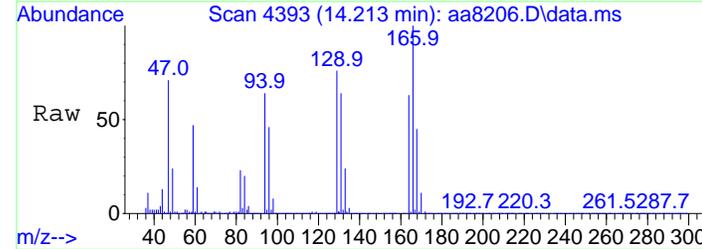
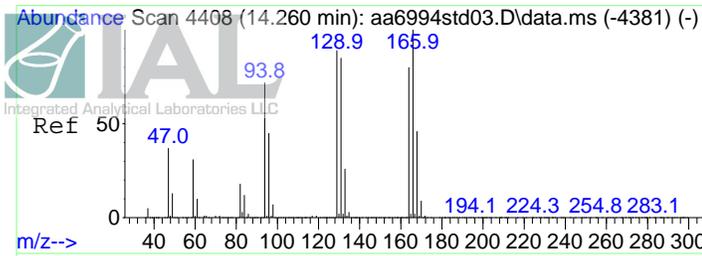
Tgt Ion:130 Resp: 352575
 Ion Ratio Lower Upper
 130 100
 128 77.9 62.6 94.0



#38
 1,4-Difluorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 9.747 min Scan# 3004
 Delta R.T. 0.006 min
 Lab File: aa8206.D
 Acq: 6 Aug 2018 10:28 pm

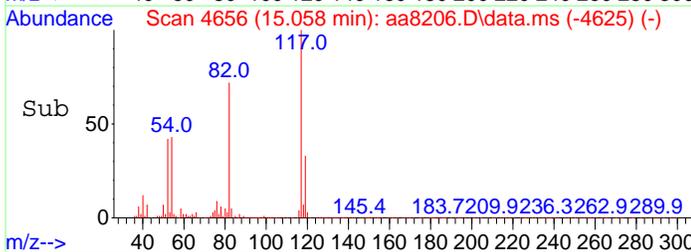
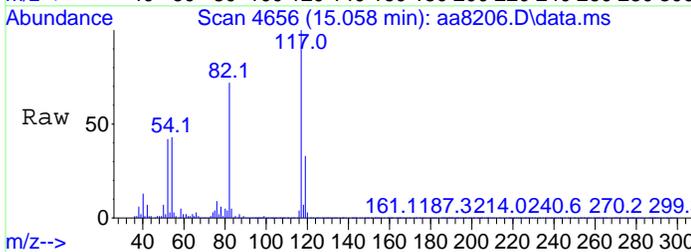
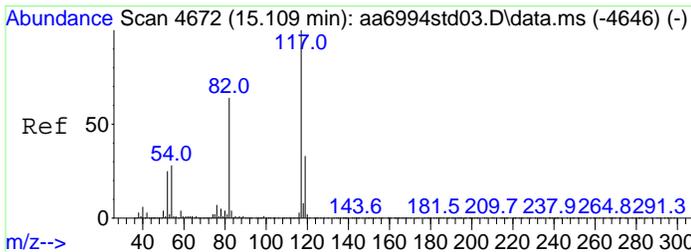
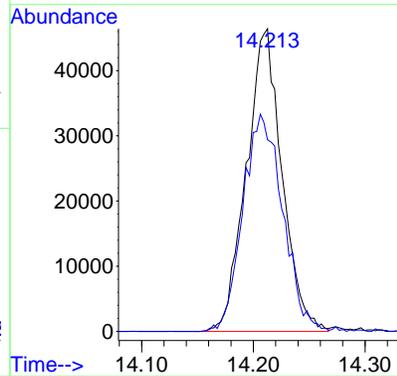
Tgt Ion:114 Resp: 1467744
 Ion Ratio Lower Upper
 114 100
 63 30.5 20.0 30.0#





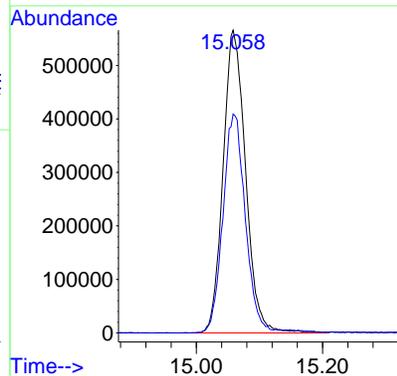
#54
 Tetrachloroethene
 Concen: 1.53 ppbV
 RT: 14.213 min Scan# 4393
 Delta R.T. 0.006 min
 Lab File: aa8206.D
 Acq: 6 Aug 2018 10:28 pm

Tgt Ion:166 Resp: 102997
 Ion Ratio Lower Upper
 166 100
 164 80.0 64.6 96.8

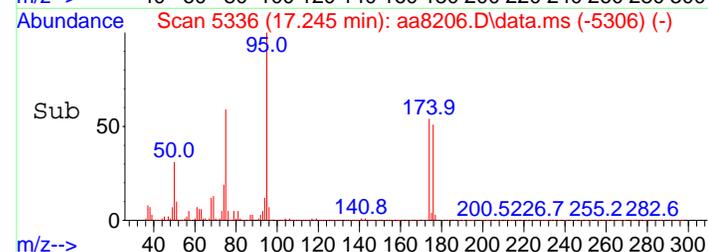
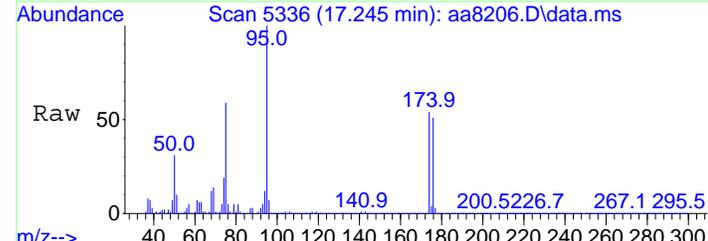
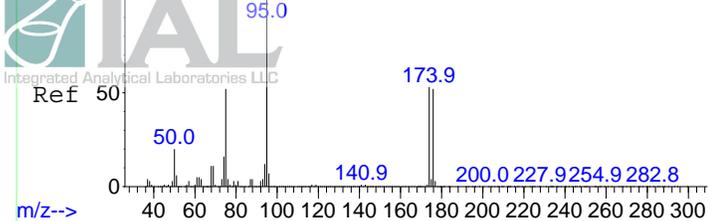


#55
 d-5 Chlorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 15.058 min Scan# 4656
 Delta R.T. 0.000 min
 Lab File: aa8206.D
 Acq: 6 Aug 2018 10:28 pm

Tgt Ion:117 Resp: 1402127
 Ion Ratio Lower Upper
 117 100
 82 71.6 56.0 84.0

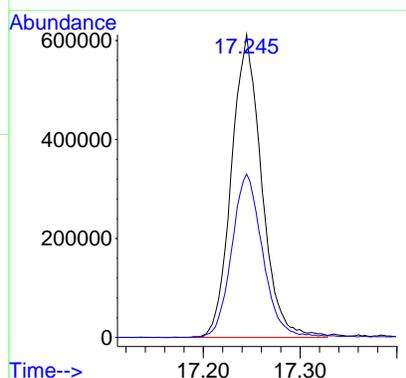


Abundance Scan 5355 (17.305 min): aa6994std03.D\data.ms (-5334) (-)



#64
Bromofluorobenzene (tune std)
Concen: 10.32 ppbV
RT: 17.245 min Scan# 5336
Delta R.T. -0.003 min
Lab File: aa8206.D
Acq: 6 Aug 2018 10:28 pm

Tgt Ion	Resp	Lower	Upper
95	1320113		
95	100		
174	55.7	61.5	92.3#





Integrated Analytical Laboratories LLC

Summary of Results

Brennan Environmental
 19 Chatham Road
 Summit, NJ 07901
 Attn: Jeff McCurdy
 Project: Congers CP / 060141
 Site: NY

Report Date: 08/13/18
 SDG Number: E18-06141
 Date Sampled: 08/01/18
 Date Received: 08/02/18
 Date Analyzed: 08/06/18, 08/07/18
 Data File: AA8207, AA8232
 Summa ID: 2155
 DF: 1, 100

Analysis: Volatile Organic Compounds by EPA Method TO-15

Compound	Sample Name:		SS-102		Reporting Limits	
	IAL ID:	CAS #	ppbv	ug/m3	ppbv	ug/m3
1,1-Dichloroethane		75-34-3	ND	ND	0.20	0.81
1,2-Dichloroethane		107-06-2	ND	ND	0.20	0.81
1,1-Dichloroethene		75-35-4	0.50	2.0	0.20	0.79
1,2-Dichloroethene (cis)	D	156-59-2	240	950	20	79
1,2-Dichloroethene (trans)		156-60-5	1.8	6.9	0.20	0.79
1,1,2,2-Tetrachloroethane		79-34-5	ND	ND	0.20	1.4
Tetrachloroethene	D	127-18-4	1500	10000	20	136
1,1,1-Trichloroethane		71-55-6	ND	ND	0.20	1.1
1,1,2-Trichloroethane		79-00-5	ND	ND	0.20	1.1
Trichloroethene	D	79-01-6	86	460	4.6	25
Vinyl chloride		75-01-4	ND	ND	0.20	0.51

Data Path : C:\DATA\08-06-18\
 Data File : aa8207.D
 Acq On : 6 Aug 2018 11:02 pm
 Operator : jls
 Sample : E18-06141-08
 Misc : 2155, 500cc
 ALS Vial : 17 Sample Multiplier: 1

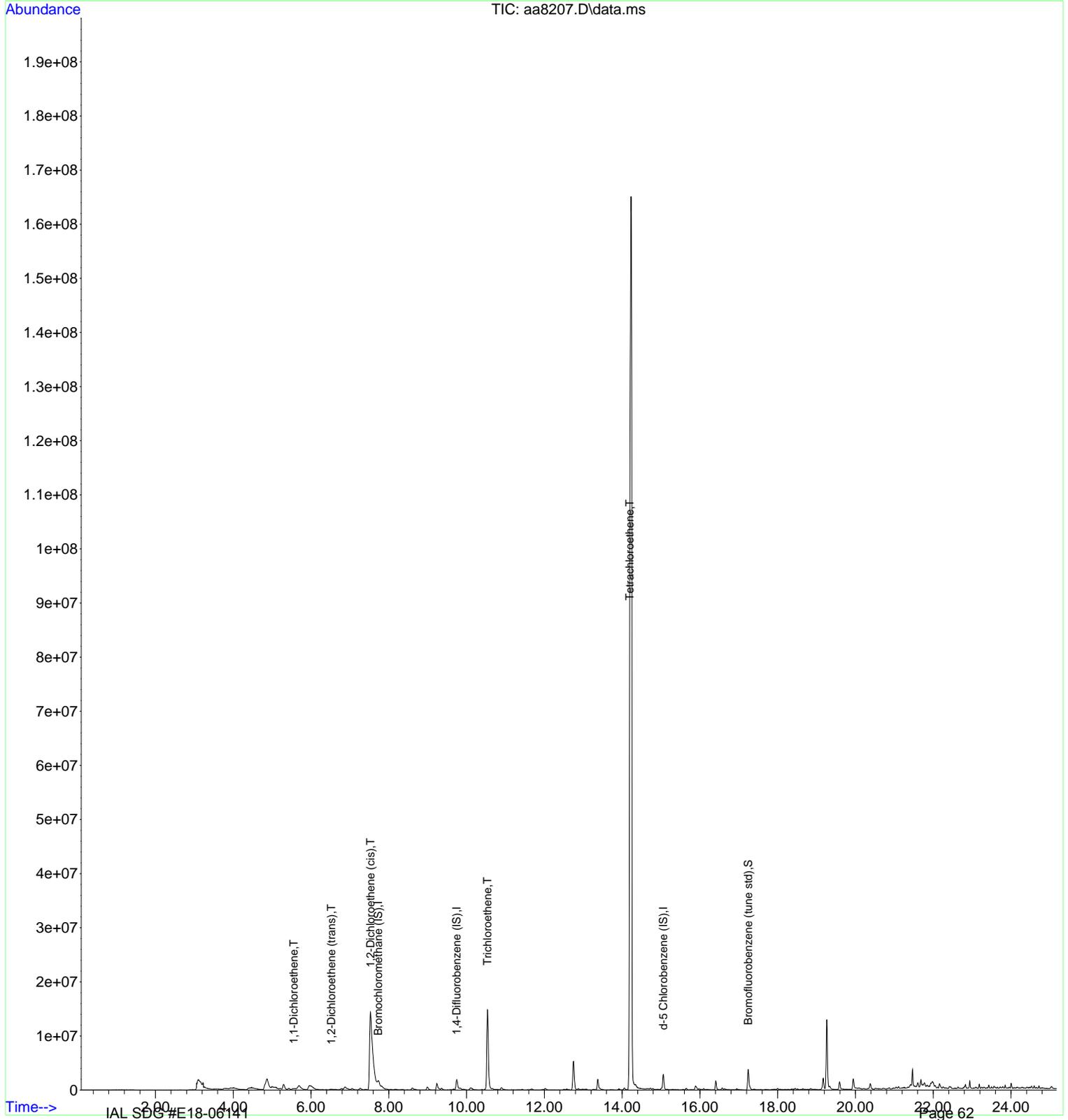
Quant Time: Aug 07 11:33:40 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

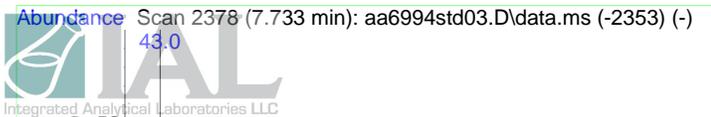
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Bromochloromethane (IS)	7.731	130	485982	10.00	ppbV	0.04
38) 1,4-Difluorobenzene (IS)	9.747	114	1975065	10.00	ppbV	0.00
55) d-5 Chlorobenzene (IS)	15.058	117	1713875	10.00	ppbV	0.00
System Monitoring Compounds						
64) Bromofluorobenzene (tu...	17.245	95	1735320	11.10	ppbV	0.00
Target Compounds						
18) 1,1-Dichloroethene	5.554	61	57333	0.50	ppbV #	62
24) 1,2-Dichloroethene (tr...	6.525	61	181309	1.75	ppbV	96
28) 1,2-Dichloroethene (cis)	7.528	61	27917034	278.07	ppbV	91
42) Trichloroethene	10.538	130	5206629	72.49	ppbV	100
54) Tetrachloroethene	14.194	166	37556129	414.72	ppbV	83

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\08-06-18\
 Data File : aa8207.D
 Acq On : 6 Aug 2018 11:02 pm
 Operator : jls
 Sample : E18-06141-08
 Misc : 2155, 500cc
 ALS Vial : 17 Sample Multiplier: 1

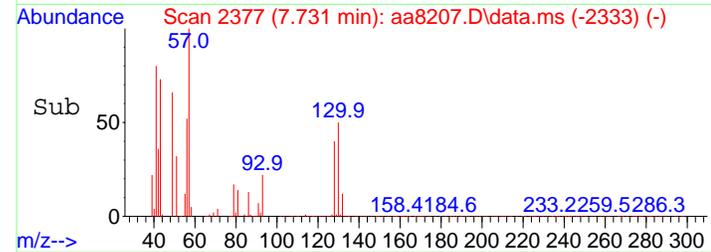
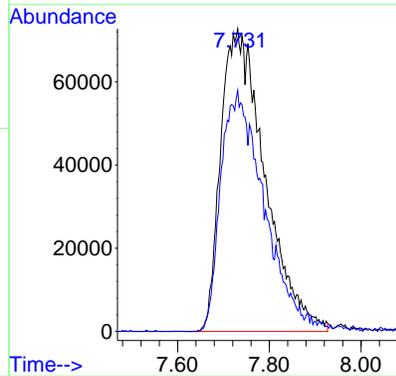
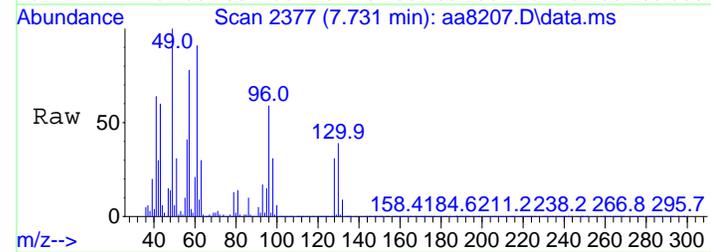
Quant Time: Aug 07 11:33:40 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration





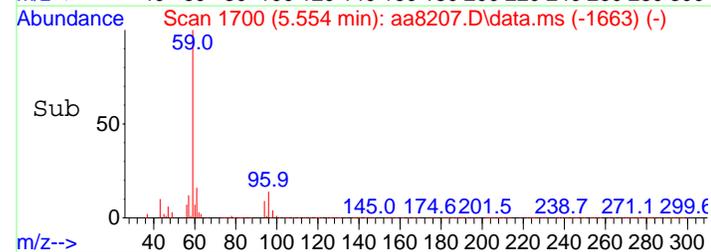
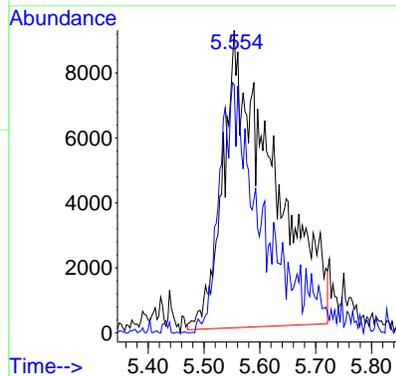
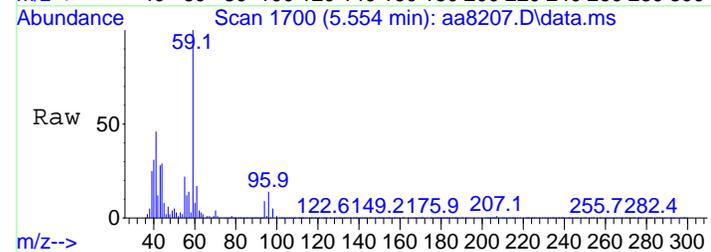
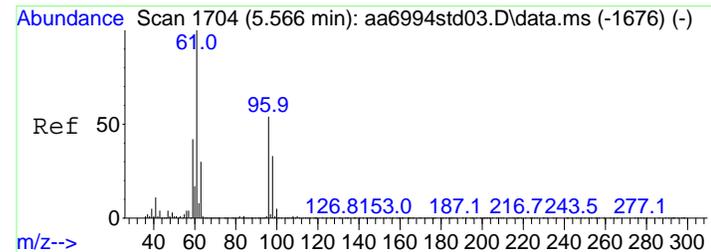
#1
 Bromochloromethane (IS)
 Concen: 10.00 ppbV
 RT: 7.731 min Scan# 2377
 Delta R.T. 0.042 min
 Lab File: aa8207.D
 Acq: 6 Aug 2018 11:02 pm

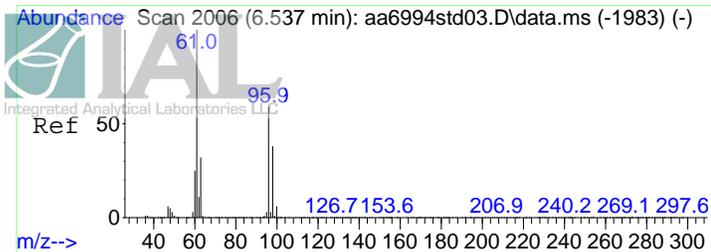
Tgt Ion: 130 Resp: 485982
 Ion Ratio Lower Upper
 130 100
 128 77.6 62.6 94.0



#18
 1,1-Dichloroethene
 Concen: 0.50 ppbV
 RT: 5.554 min Scan# 1700
 Delta R.T. 0.019 min
 Lab File: aa8207.D
 Acq: 6 Aug 2018 11:02 pm

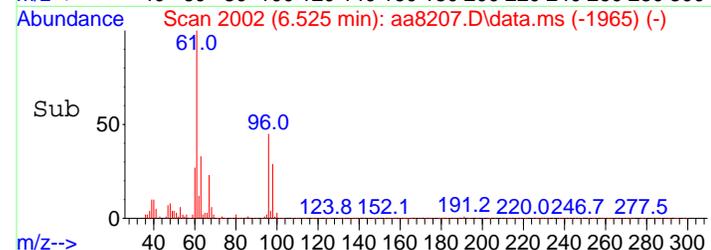
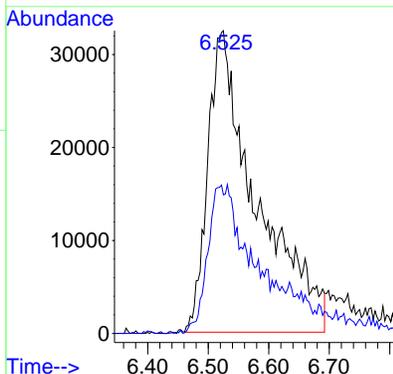
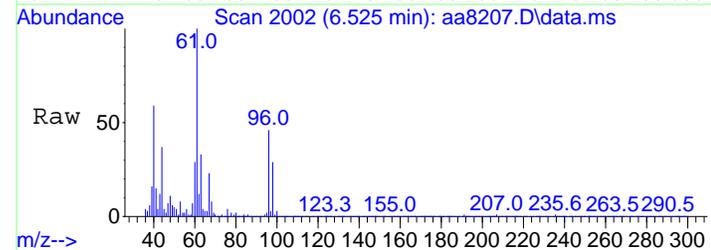
Tgt Ion: 61 Resp: 57333
 Ion Ratio Lower Upper
 61 100
 96 25.3 41.6 62.4#





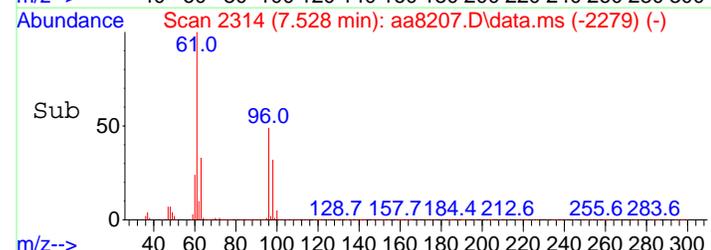
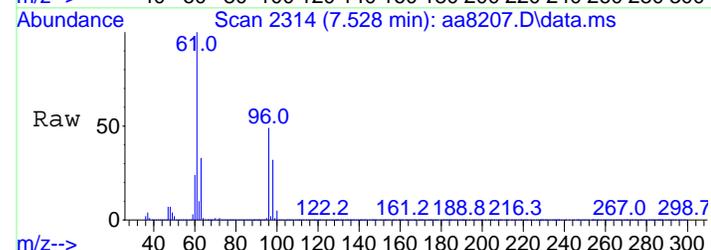
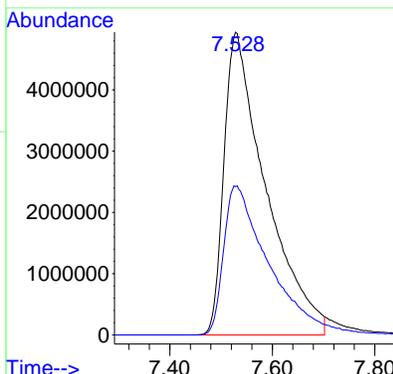
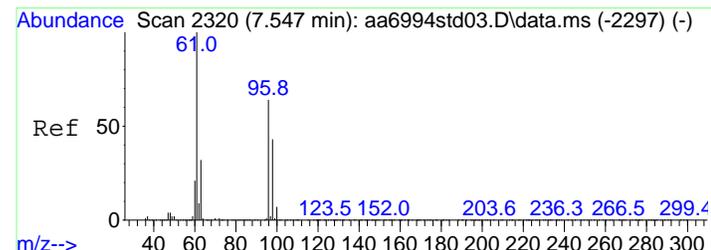
#24
 1,2-Dichloroethene (trans)
 Concen: 1.75 ppbV
 RT: 6.525 min Scan# 2002
 Delta R.T. 0.019 min
 Lab File: aa8207.D
 Acq: 6 Aug 2018 11:02 pm

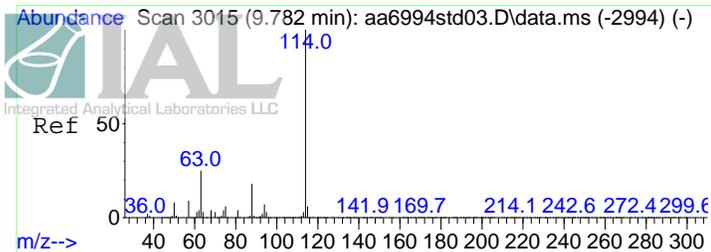
Tgt Ion: 61 Resp: 181309
 Ion Ratio Lower Upper
 61 100
 96 51.4 43.2 64.8



#28
 1,2-Dichloroethene (cis)
 Concen: 278.07 ppbV
 RT: 7.528 min Scan# 2314
 Delta R.T. 0.013 min
 Lab File: aa8207.D
 Acq: 6 Aug 2018 11:02 pm

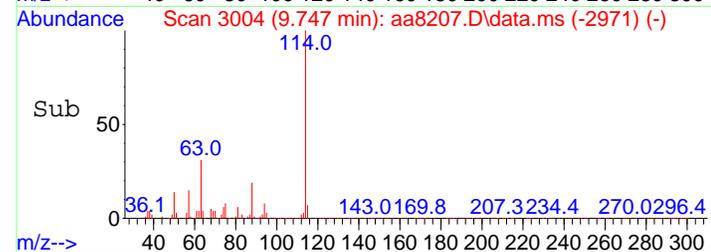
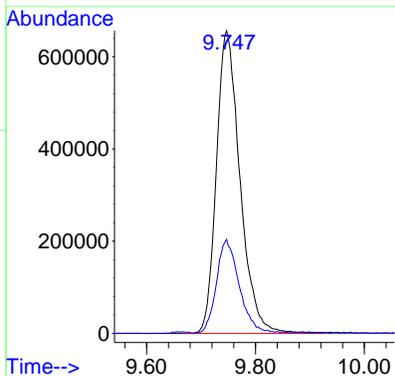
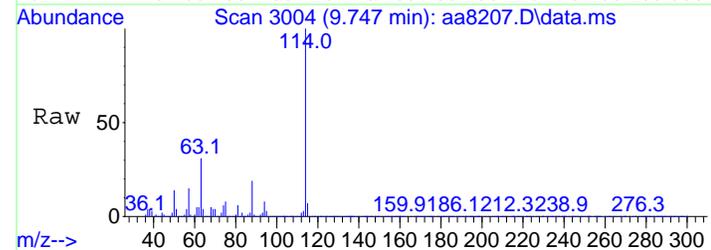
Tgt Ion: 61 Resp: 27917034
 Ion Ratio Lower Upper
 61 100
 96 52.0 47.2 70.8





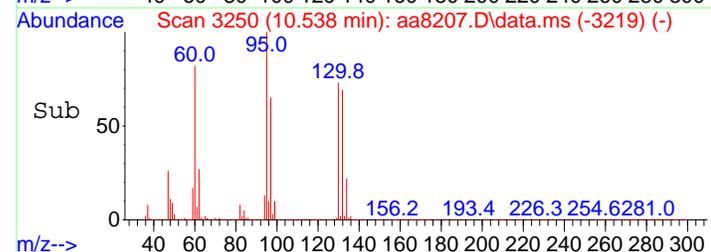
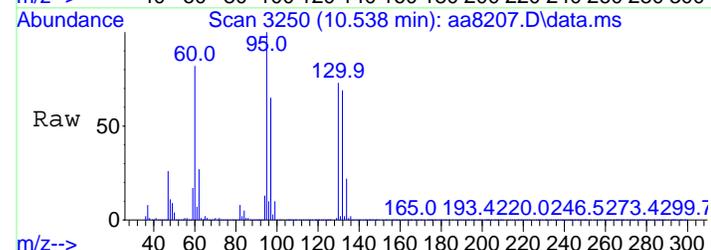
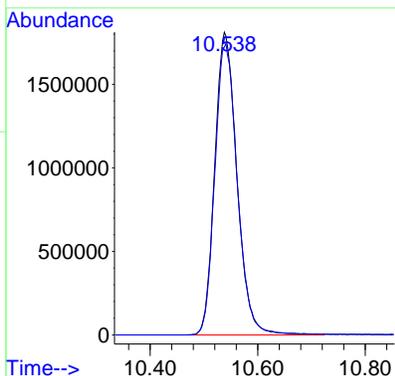
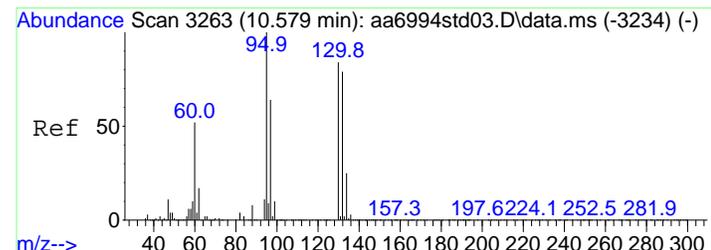
#38
 1,4-Difluorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 9.747 min Scan# 3004
 Delta R.T. 0.007 min
 Lab File: aa8207.D
 Acq: 6 Aug 2018 11:02 pm

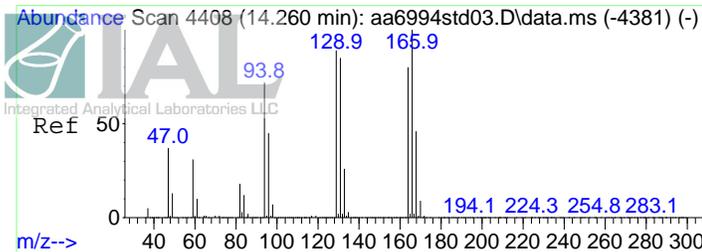
Tgt Ion:114 Resp: 1975065
 Ion Ratio Lower Upper
 114 100
 63 29.9 20.0 30.0



#42
 Trichloroethene
 Concen: 72.49 ppbV
 RT: 10.538 min Scan# 3250
 Delta R.T. 0.000 min
 Lab File: aa8207.D
 Acq: 6 Aug 2018 11:02 pm

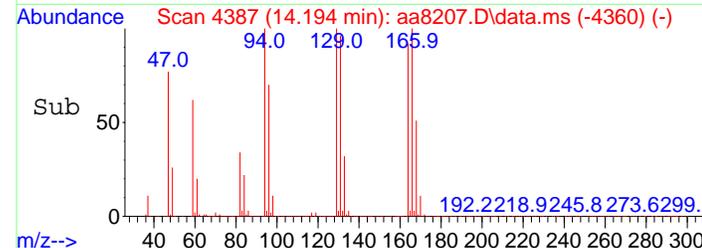
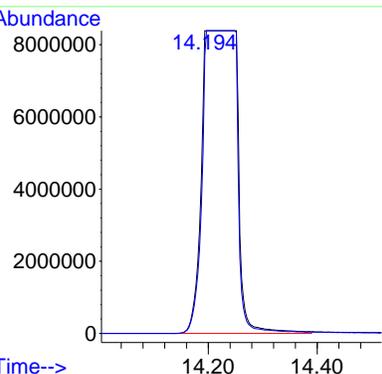
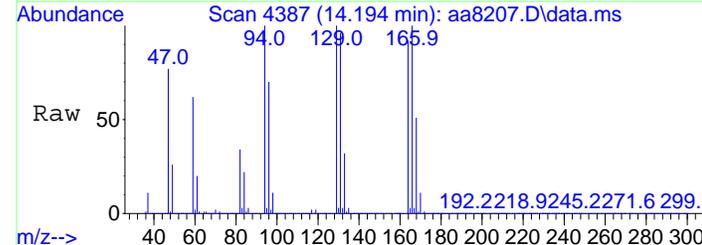
Tgt Ion:130 Resp: 5206629
 Ion Ratio Lower Upper
 130 100
 132 97.3 77.8 116.6





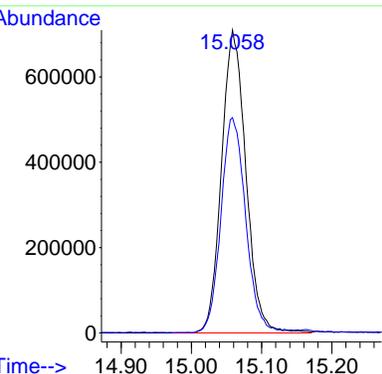
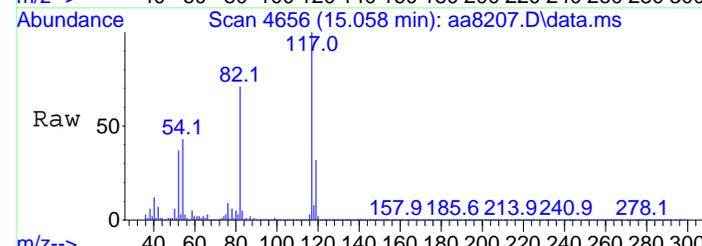
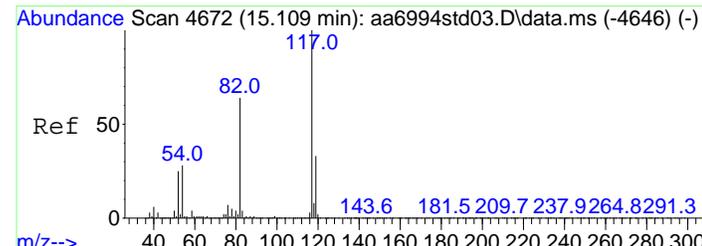
#54
 Tetrachloroethene
 Concen: 414.72 ppbV
 RT: 14.194 min Scan# 4387
 Delta R.T. -0.013 min
 Lab File: aa8207.D
 Acq: 6 Aug 2018 11:02 pm

Tgt Ion:166 Resp:37556129
 Ion Ratio Lower Upper
 166 100
 164 95.9 64.6 96.8

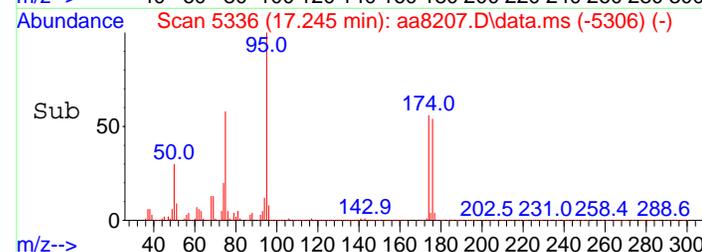
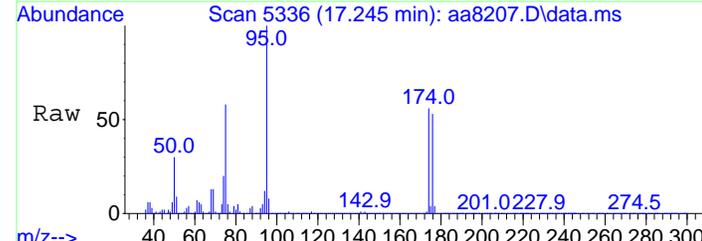
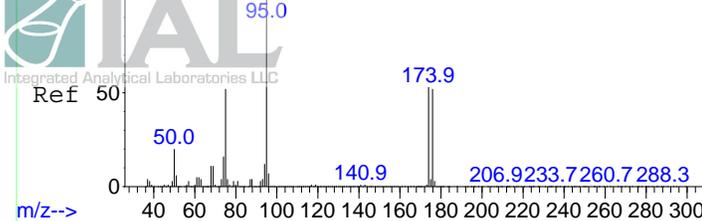


#55
 d-5 Chlorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 15.058 min Scan# 4656
 Delta R.T. 0.000 min
 Lab File: aa8207.D
 Acq: 6 Aug 2018 11:02 pm

Tgt Ion:117 Resp: 1713875
 Ion Ratio Lower Upper
 117 100
 82 73.0 56.0 84.0

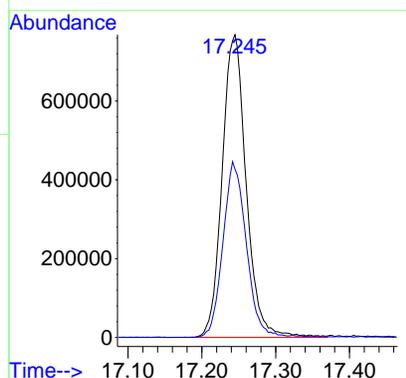


Abundance Scan 5355 (17.305 min): aa6994std03.D\data.ms (-5334) (-)



#64
Bromofluorobenzene (tune std)
Concen: 11.10 ppbV
RT: 17.245 min Scan# 5336
Delta R.T. -0.003 min
Lab File: aa8207.D
Acq: 6 Aug 2018 11:02 pm

Tgt Ion	Resp	Lower	Upper
95	1735320		
95	100		
174	57.0	61.5	92.3#



Data Path : C:\DATA\08-07-18\
 Data File : aa8232.D
 Acq On : 7 Aug 2018 4:57 pm
 Operator : jjw
 Sample : E18-06141-08 x 100 dil
 Misc : 2155, 5cc
 ALS Vial : 12 Sample Multiplier: 1

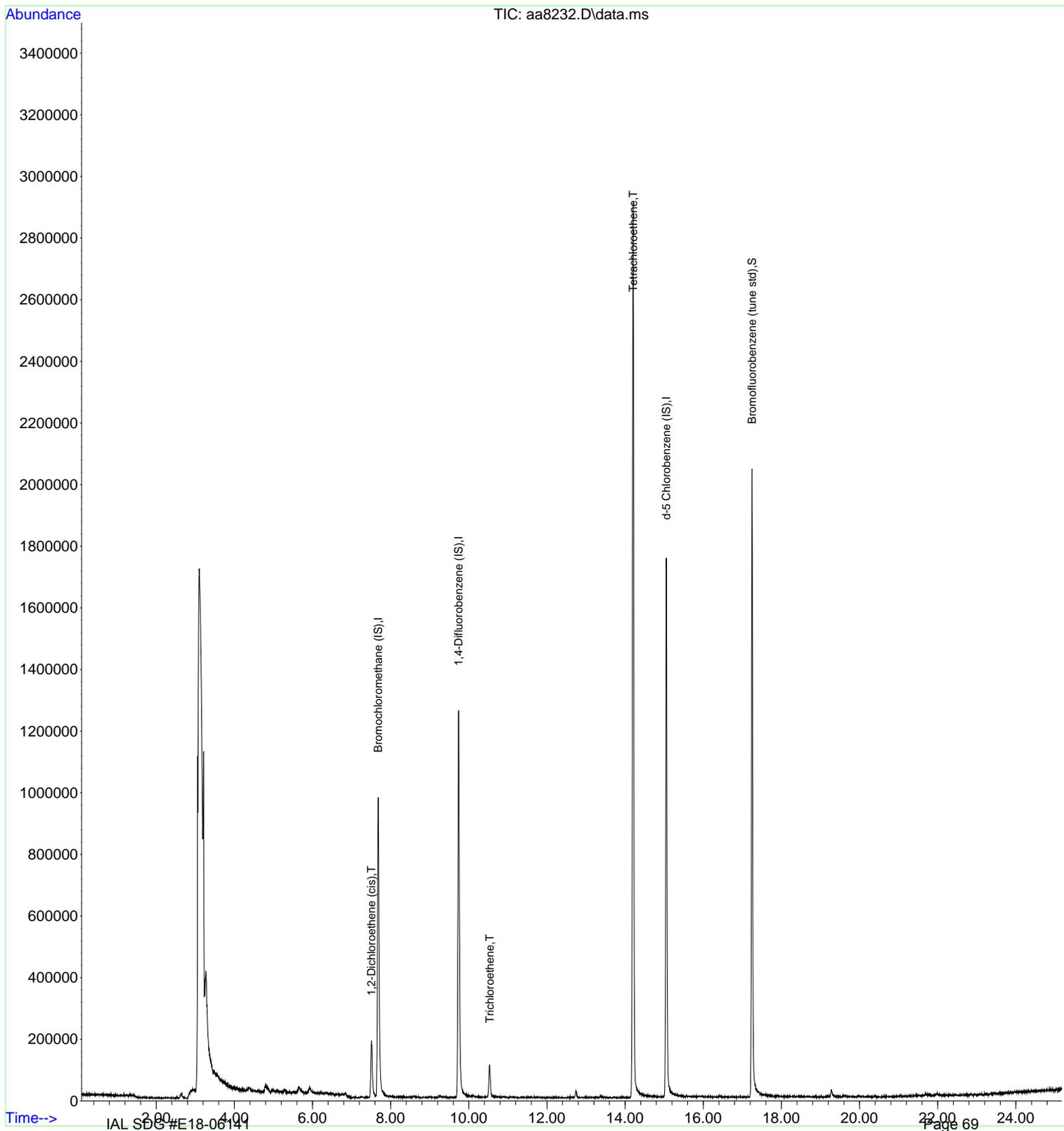
Quant Time: Aug 09 11:20:50 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

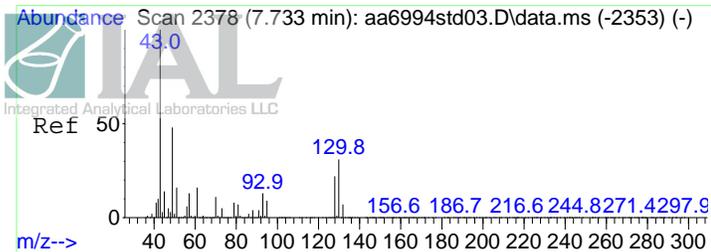
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Bromochloromethane (IS)	7.682	130	332191	10.00	ppbV	0.00
38) 1,4-Difluorobenzene (IS)	9.737	114	1080070	10.00	ppbV #	0.00
55) d-5 Chlorobenzene (IS)	15.052	117	922301	10.00	ppbV	0.00
System Monitoring Compounds						
64) Bromofluorobenzene (tu...	17.248	95	817274	9.72	ppbV	0.00
Target Compounds						
28) 1,2-Dichloroethene (cis)	7.509	61	164078	2.39	ppbV	98
42) Trichloroethene	10.538	130	33853	0.86	ppbV	98
54) Tetrachloroethene	14.206	166	753398	15.21	ppbV	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\08-07-18\
Data File : aa8232.D
Acq On : 7 Aug 2018 4:57 pm
Operator : jjw
Sample : E18-06141-08 x 100 dil
Misc : 2155, 5cc
ALS Vial : 12 Sample Multiplier: 1

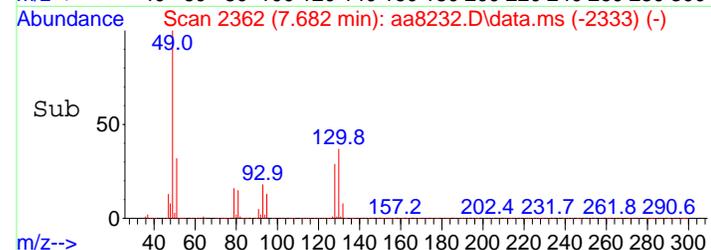
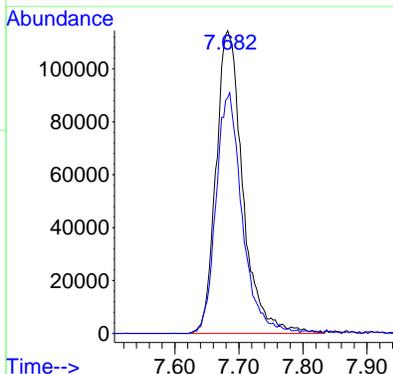
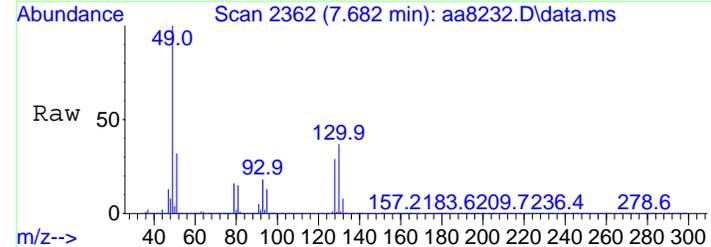
Quant Time: Aug 09 11:20:50 2018
Quant Method : C:\msdchem\1\METHODS\0725.M
Quant Title : TO-15 on the Agilent 7890A / 5975C
QLast Update : Wed Jul 25 14:15:57 2018
Response via : Initial Calibration





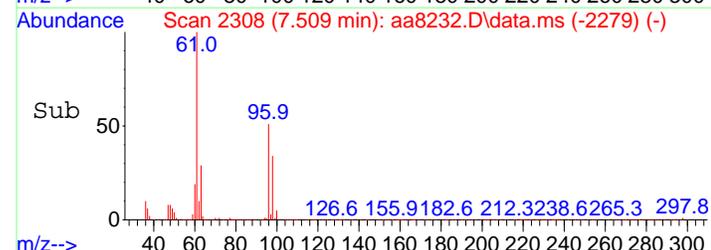
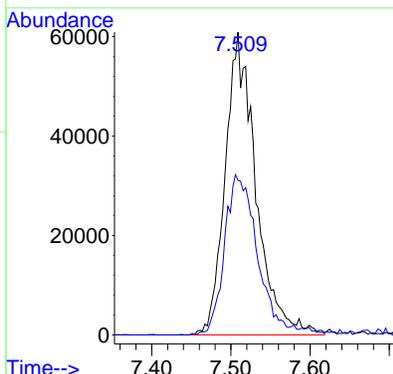
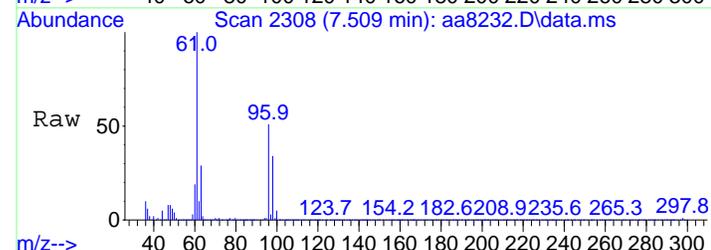
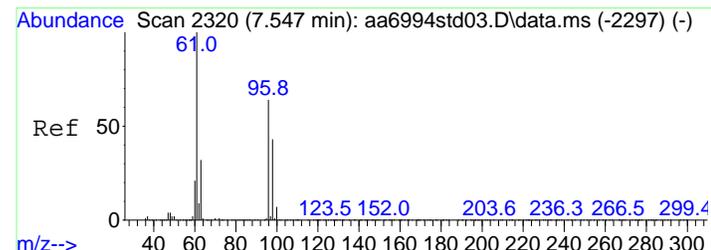
#1
 Bromochloromethane (IS)
 Concen: 10.00 ppbV
 RT: 7.682 min Scan# 2362
 Delta R.T. -0.007 min
 Lab File: aa8232.D
 Acq: 7 Aug 2018 4:57 pm

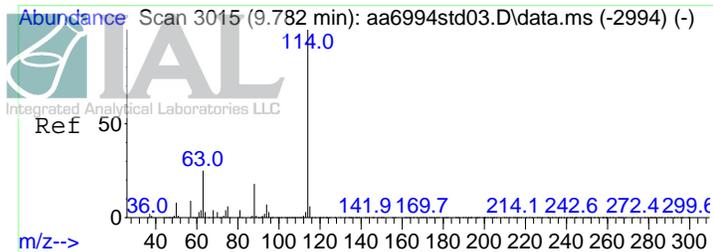
Tgt Ion:130 Resp: 332191
 Ion Ratio Lower Upper
 130 100
 128 77.7 62.6 94.0



#28
 1,2-Dichloroethene (cis)
 Concen: 2.39 ppbV
 RT: 7.509 min Scan# 2308
 Delta R.T. -0.006 min
 Lab File: aa8232.D
 Acq: 7 Aug 2018 4:57 pm

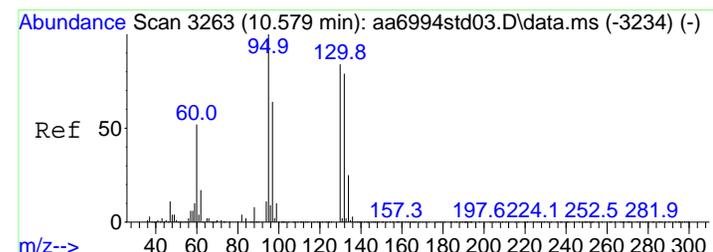
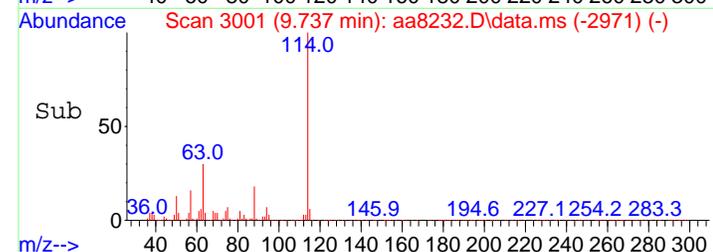
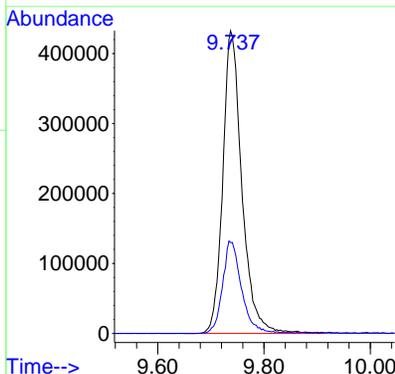
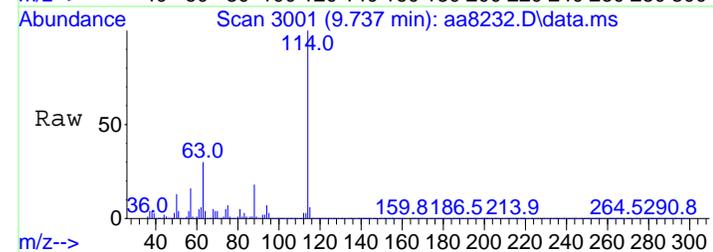
Tgt Ion: 61 Resp: 164078
 Ion Ratio Lower Upper
 61 100
 96 57.5 47.2 70.8





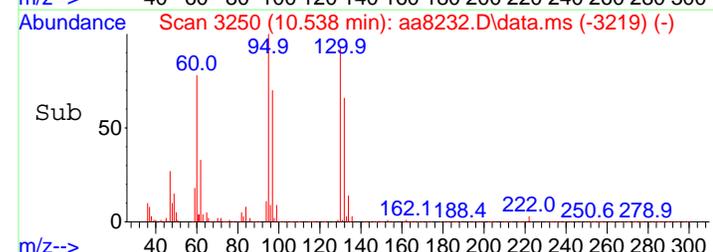
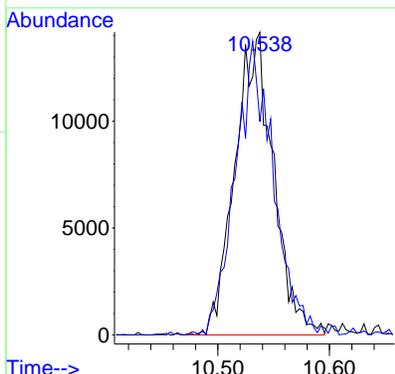
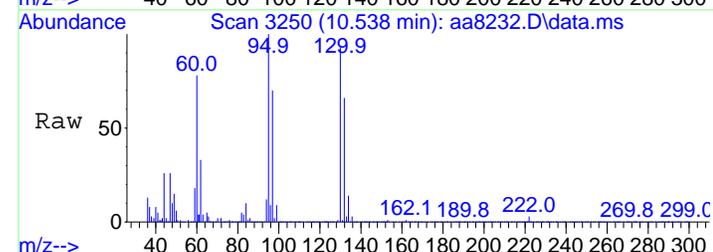
#38
 1,4-Difluorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 9.737 min Scan# 3001
 Delta R.T. -0.003 min
 Lab File: aa8232.D
 Acq: 7 Aug 2018 4:57 pm

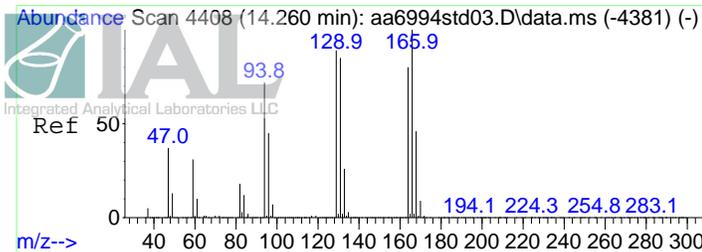
Tgt Ion:114 Resp: 1080070
 Ion Ratio Lower Upper
 114 100
 63 30.2 20.0 30.0#



#42
 Trichloroethene
 Concen: 0.86 ppbV
 RT: 10.538 min Scan# 3250
 Delta R.T. 0.000 min
 Lab File: aa8232.D
 Acq: 7 Aug 2018 4:57 pm

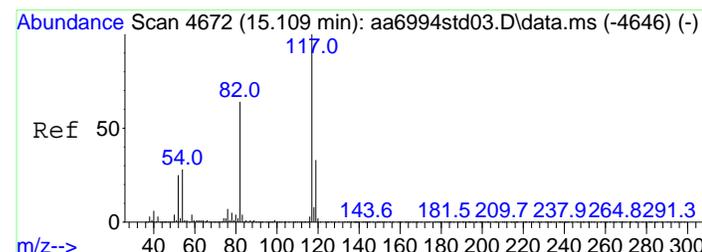
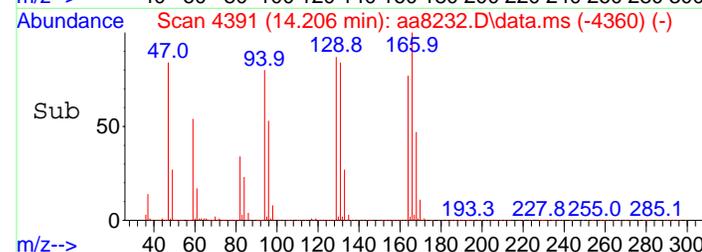
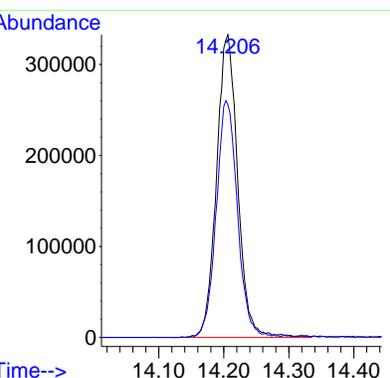
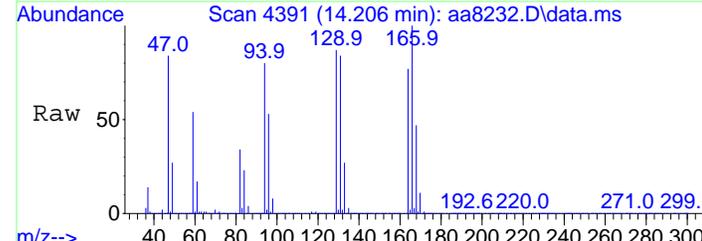
Tgt Ion:130 Resp: 33853
 Ion Ratio Lower Upper
 130 100
 132 95.3 77.8 116.6





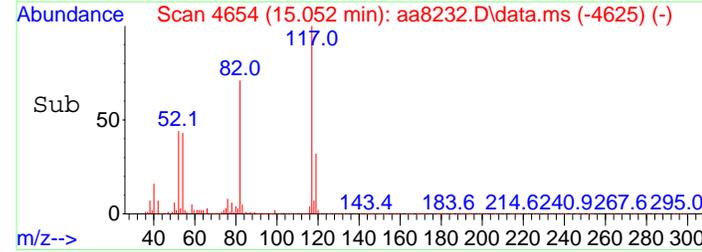
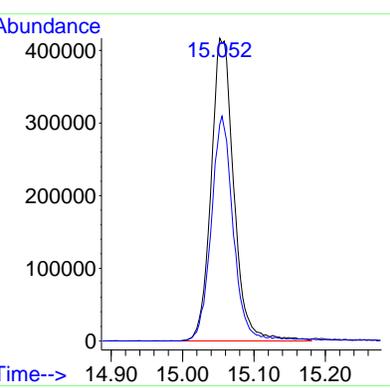
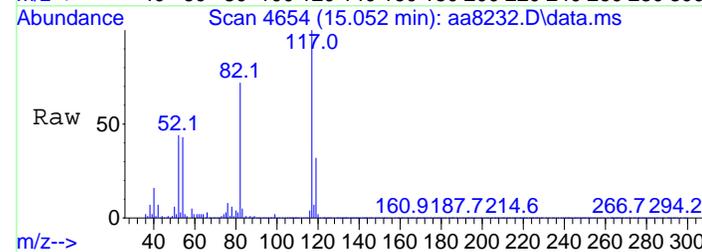
#54
 Tetrachloroethene
 Concen: 15.21 ppbV
 RT: 14.206 min Scan# 4391
 Delta R.T. 0.000 min
 Lab File: aa8232.D
 Acq: 7 Aug 2018 4:57 pm

Tgt Ion	Resp	Lower	Upper
166	753398		
166	100		
164	78.3	64.6	96.8

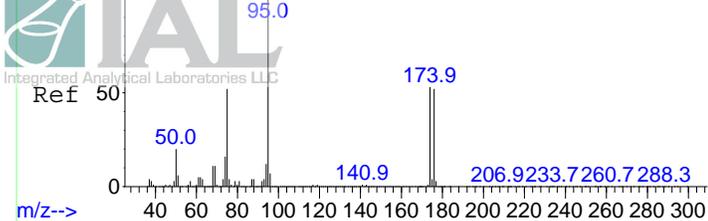


#55
 d-5 Chlorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 15.052 min Scan# 4654
 Delta R.T. -0.006 min
 Lab File: aa8232.D
 Acq: 7 Aug 2018 4:57 pm

Tgt Ion	Resp	Lower	Upper
117	922301		
117	100		
82	71.7	56.0	84.0

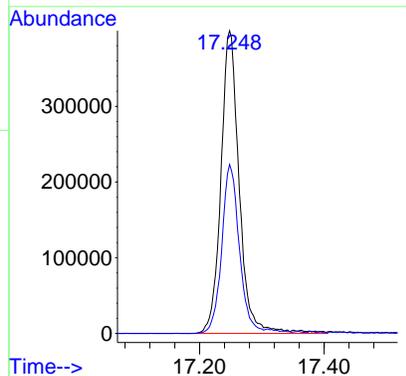
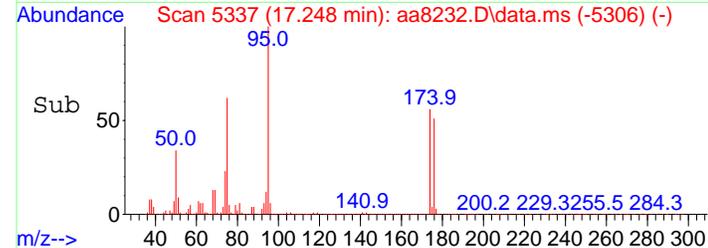
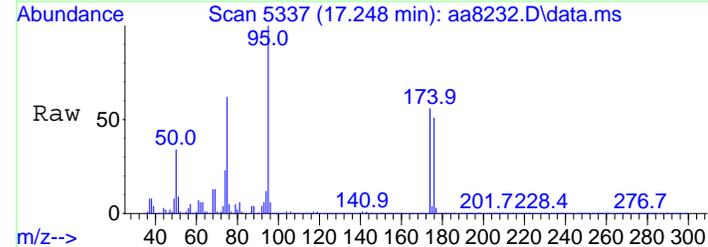


Abundance Scan 5355 (17.305 min): aa6994std03.D\data.ms (-5334) (-)



#64
Bromofluorobenzene (tune std)
Concen: 9.72 ppbV
RT: 17.248 min Scan# 5337
Delta R.T. 0.000 min
Lab File: aa8232.D
Acq: 7 Aug 2018 4:57 pm

Tgt Ion: 95 Resp: 817274
Ion Ratio Lower Upper
95 100
174 54.5 61.5 92.3#





Integrated Analytical Laboratories LLC

Summary of Results

Brennan Environmental
 19 Chatham Road
 Summit, NJ 07901
 Attn: Jeff McCurdy
 Project: Congers CP / 060141
 Site: NY

Report Date: 08/13/18
 SDG Number: E18-06141
 Date Sampled: 08/01/18
 Date Received: 08/02/18
 Date Analyzed: 08/06/18, 08/07/18
 Data File: AA8208, AA8233
 Summa ID: 5089
 DF: 1, 10

Analysis: Volatile Organic Compounds by EPA Method TO-15

Compound	Sample Name:	SS-103		Reporting Limits	
	IAL ID:	E18-06141-09			
	CAS #	ppbv	ug/m3	ppbv	ug/m3
1,1-Dichloroethane	75-34-3	ND	ND	0.20	0.81
1,2-Dichloroethane	107-06-2	ND	ND	0.20	0.81
1,1-Dichloroethene	75-35-4	ND	ND	0.20	0.79
1,2-Dichloroethene (cis)	156-59-2	ND	ND	0.20	0.79
1,2-Dichloroethene (trans)	156-60-5	ND	ND	0.20	0.79
1,1,2,2-Tetrachloroethane	79-34-5	ND	ND	0.20	1.4
Tetrachloroethene	127-18-4	D	140	950	2.0
1,1,1-Trichloroethane	71-55-6	ND	ND	0.20	1.1
1,1,2-Trichloroethane	79-00-5	ND	ND	0.20	1.1
Trichloroethene	79-01-6	0.17	0.93	0.05	0.25
Vinyl chloride	75-01-4	ND	ND	0.20	0.51

Data Path : C:\DATA\08-06-18\
 Data File : aa8208.D
 Acq On : 6 Aug 2018 11:35 pm
 Operator : jls
 Sample : E18-06141-09
 Misc : 5089, 500cc
 ALS Vial : 18 Sample Multiplier: 1

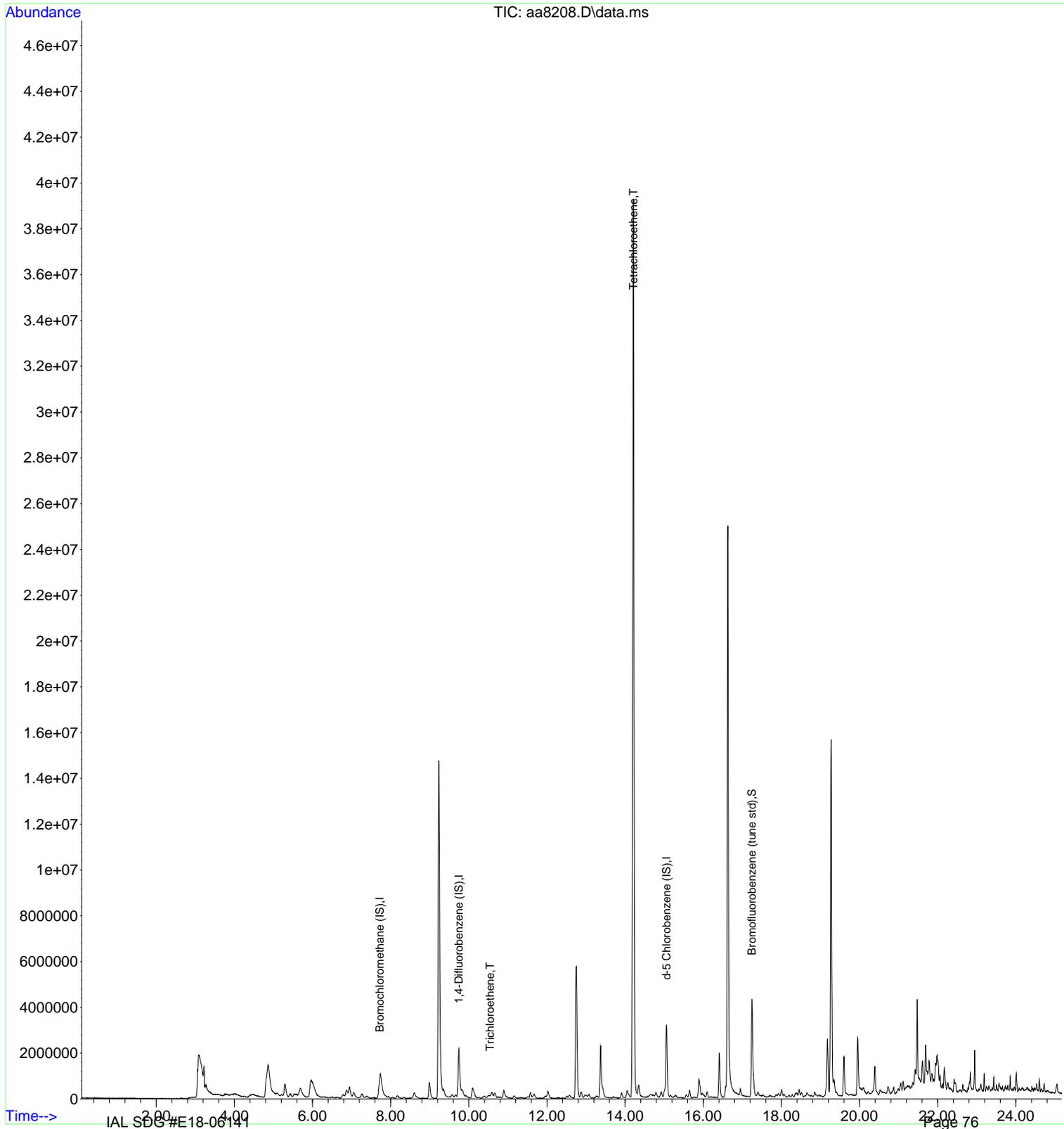
Quant Time: Aug 07 11:41:23 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

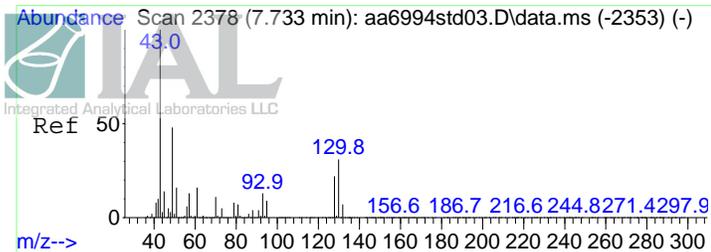
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.721	130	516268	10.00	ppbV	0.03	
38) 1,4-Difluorobenzene (IS)	9.747	114	2306557	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.058	117	1996784	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...	17.245	95	2065867	11.34	ppbV	0.00	
Target Compounds							
42) Trichloroethene	10.541	130	14488	0.17	ppbV	99	Qvalue
54) Tetrachloroethene	14.210	166	10904647	103.11	ppbV	98	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\08-06-18\
Data File : aa8208.D
Acq On : 6 Aug 2018 11:35 pm
Operator : jls
Sample : E18-06141-09
Misc : 5089, 500cc
ALS Vial : 18 Sample Multiplier: 1

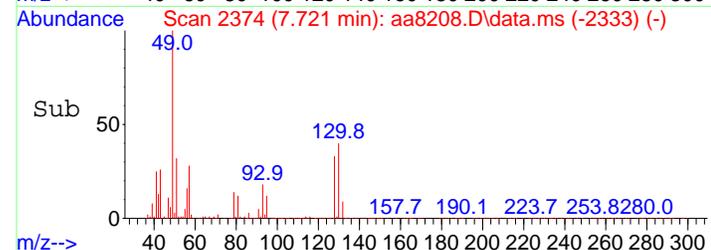
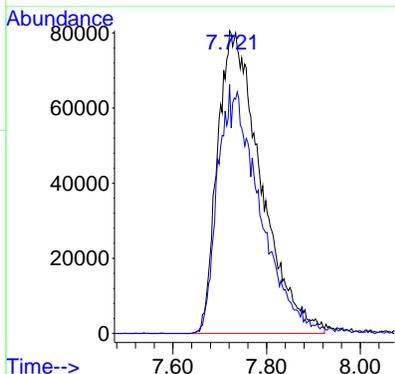
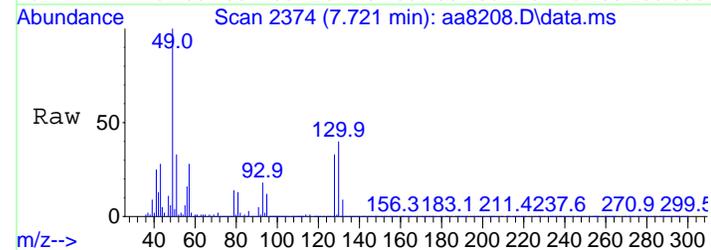
Quant Time: Aug 07 11:41:23 2018
Quant Method : C:\msdchem\1\METHODS\0725.M
Quant Title : TO-15 on the Agilent 7890A / 5975C
QLast Update : Wed Jul 25 14:15:57 2018
Response via : Initial Calibration





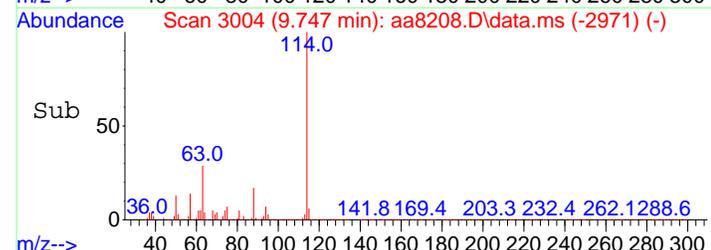
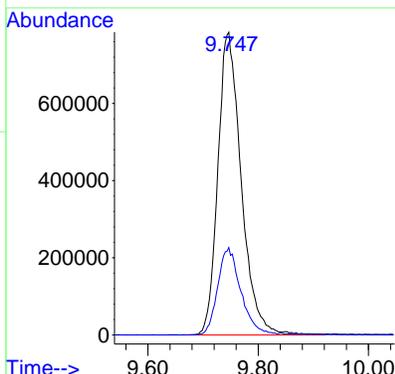
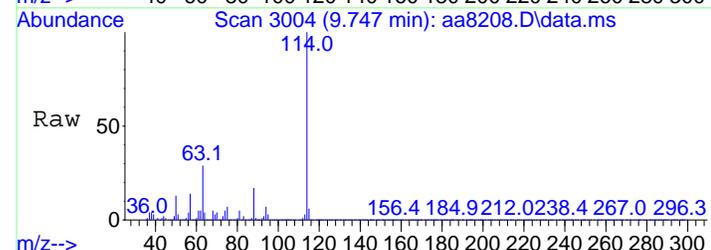
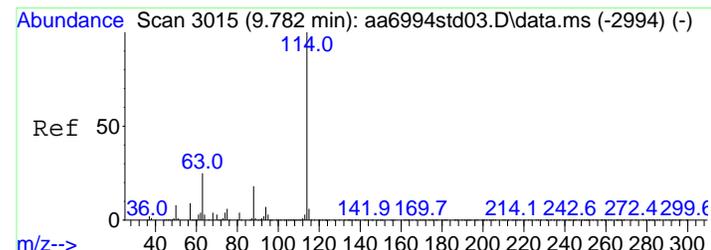
#1
 Bromochloromethane (IS)
 Concen: 10.00 ppbV
 RT: 7.721 min Scan# 2374
 Delta R.T. 0.032 min
 Lab File: aa8208.D
 Acq: 6 Aug 2018 11:35 pm

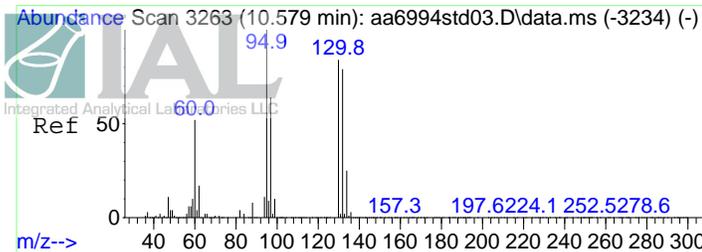
Tgt Ion:130 Resp: 516268
 Ion Ratio Lower Upper
 130 100
 128 79.1 62.6 94.0



#38
 1,4-Difluorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 9.747 min Scan# 3004
 Delta R.T. 0.007 min
 Lab File: aa8208.D
 Acq: 6 Aug 2018 11:35 pm

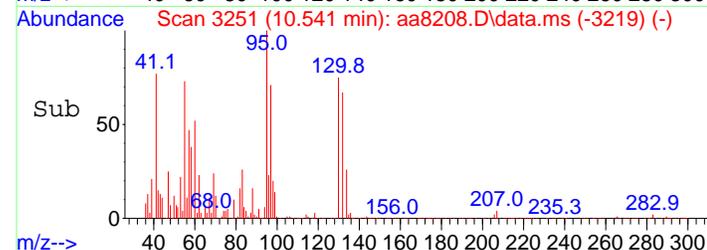
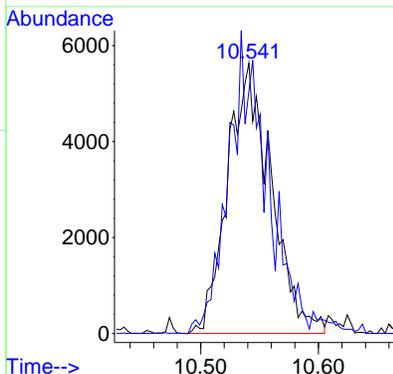
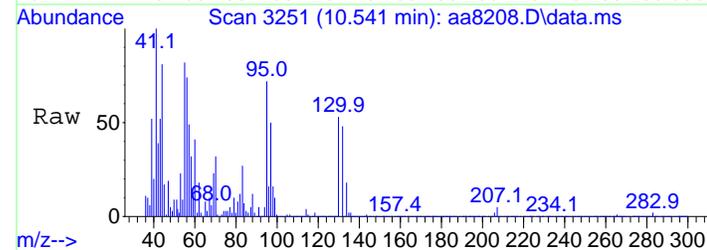
Tgt Ion:114 Resp: 2306557
 Ion Ratio Lower Upper
 114 100
 63 28.4 20.0 30.0





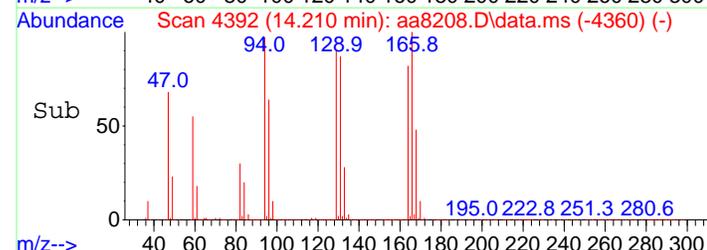
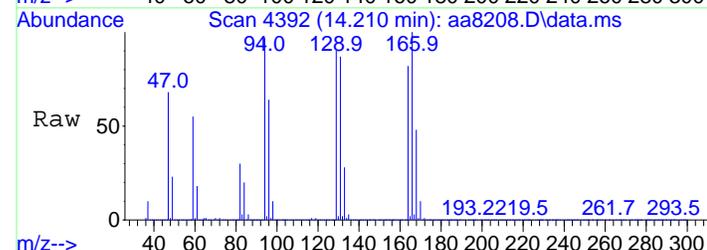
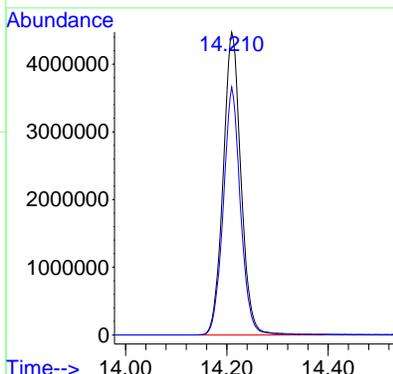
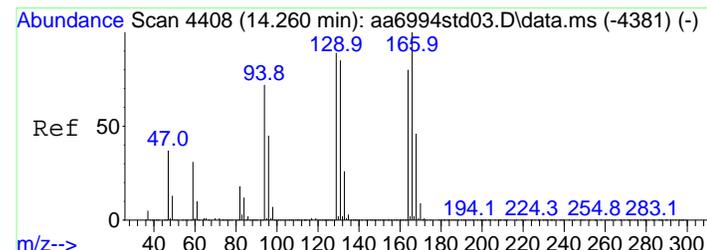
#42
 Trichloroethene
 Concen: 0.17 ppbV
 RT: 10.541 min Scan# 3251
 Delta R.T. 0.003 min
 Lab File: aa8208.D
 Acq: 6 Aug 2018 11:35 pm

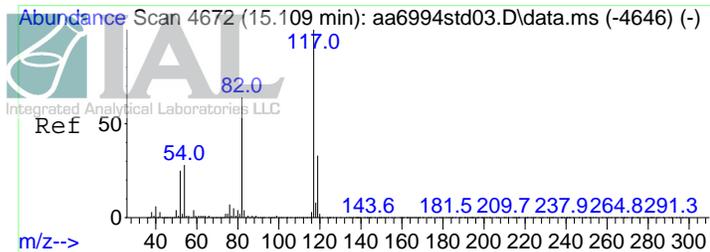
Tgt Ion:130 Resp: 14488
 Ion Ratio Lower Upper
 130 100
 132 97.7 77.8 116.6



#54
 Tetrachloroethene
 Concen: 103.11 ppbV
 RT: 14.210 min Scan# 4392
 Delta R.T. 0.003 min
 Lab File: aa8208.D
 Acq: 6 Aug 2018 11:35 pm

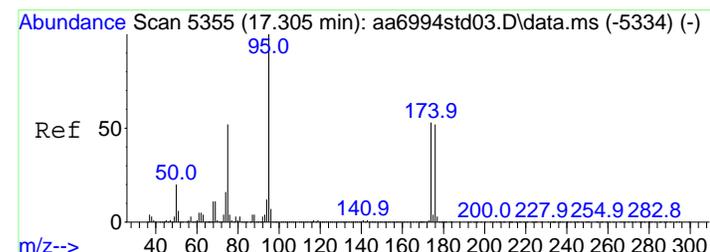
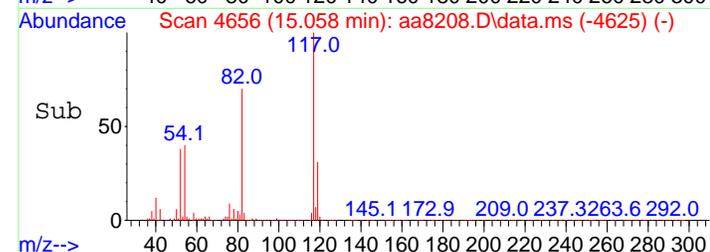
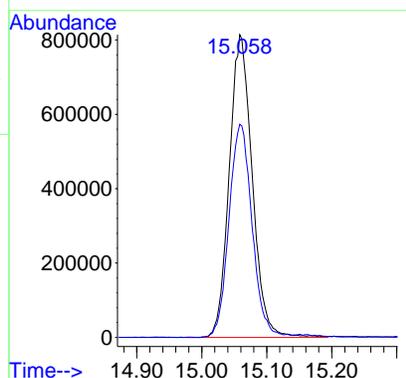
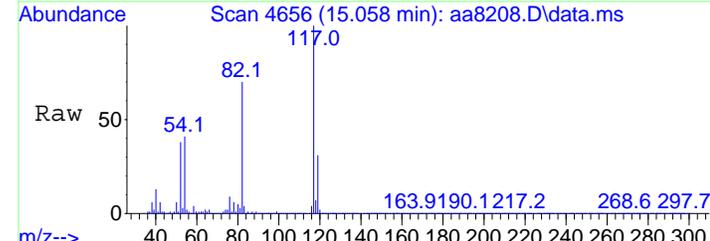
Tgt Ion:166 Resp:10904647
 Ion Ratio Lower Upper
 166 100
 164 79.3 64.6 96.8





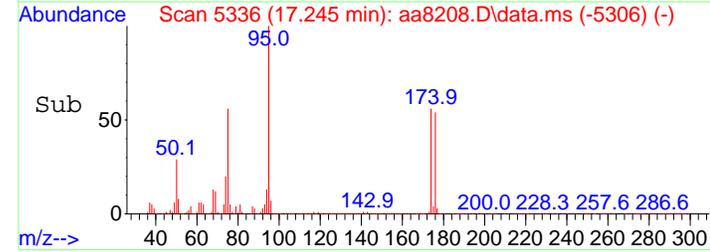
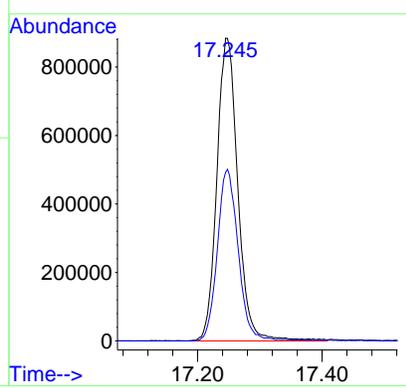
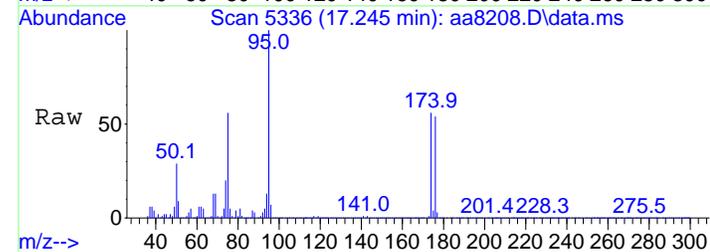
#55
 d-5 Chlorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 15.058 min Scan# 4656
 Delta R.T. 0.000 min
 Lab File: aa8208.D
 Acq: 6 Aug 2018 11:35 pm

Tgt Ion	Resp	Lower	Upper
117	100		
82	71.8	56.0	84.0



#64
 Bromofluorobenzene (tune std)
 Concen: 11.34 ppbV
 RT: 17.245 min Scan# 5336
 Delta R.T. -0.003 min
 Lab File: aa8208.D
 Acq: 6 Aug 2018 11:35 pm

Tgt Ion	Resp	Lower	Upper
95	100		
174	56.4	61.5	92.3#



Data Path : C:\DATA\08-07-18\
 Data File : aa8233.D
 Acq On : 7 Aug 2018 5:31 pm
 Operator : jjw
 Sample : E18-06141-09 x 10 dil
 Misc : 5089, 50cc
 ALS Vial : 13 Sample Multiplier: 1

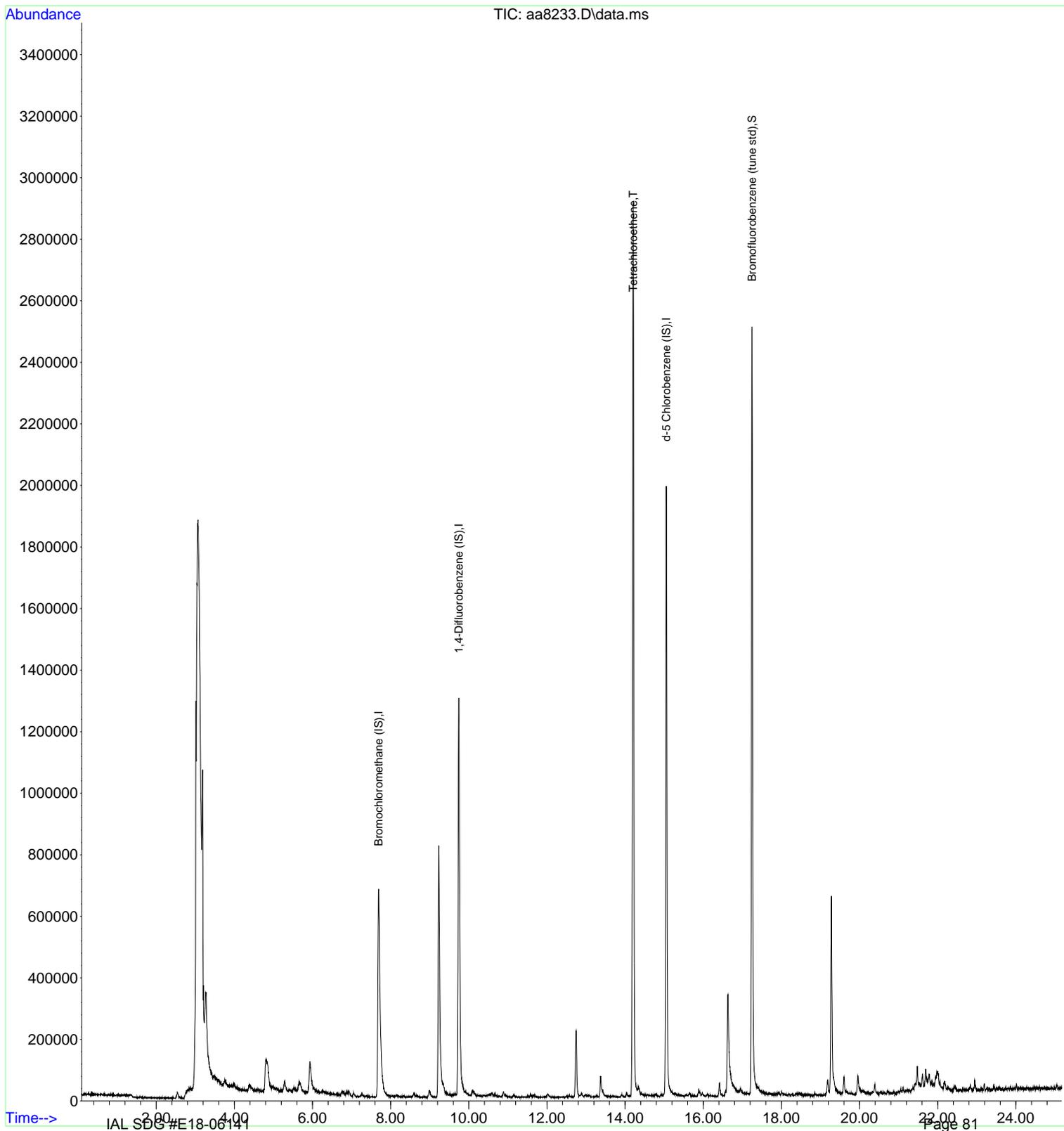
Quant Time: Aug 09 11:24:30 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

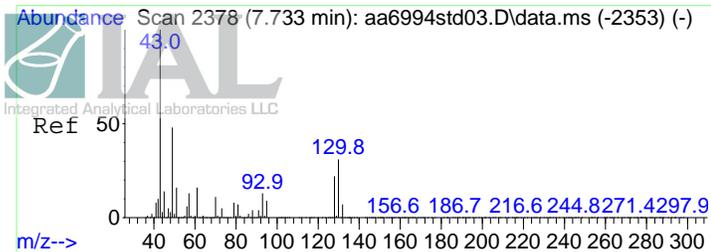
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Bromochloromethane (IS)	7.692	130	305198	10.00	ppbV	0.00
38) 1,4-Difluorobenzene (IS)	9.743	114	1178297	10.00	ppbV #	0.00
55) d-5 Chlorobenzene (IS)	15.055	117	1041852	10.00	ppbV	0.00
System Monitoring Compounds						
64) Bromofluorobenzene (tu...	17.248	95	1043585	10.98	ppbV	0.00
Target Compounds						
54) Tetrachloroethene	14.206	166	755419	13.98	ppbV	Qvalue 98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\08-07-18\
Data File : aa8233.D
Acq On : 7 Aug 2018 5:31 pm
Operator : jjw
Sample : E18-06141-09 x 10 dil
Misc : 5089, 50cc
ALS Vial : 13 Sample Multiplier: 1

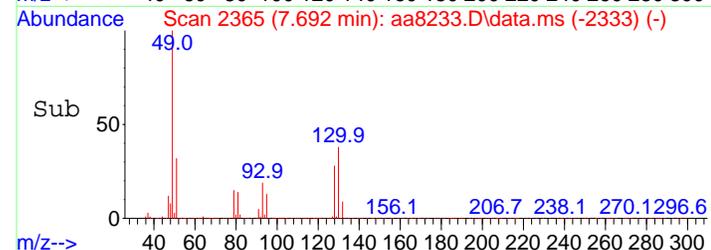
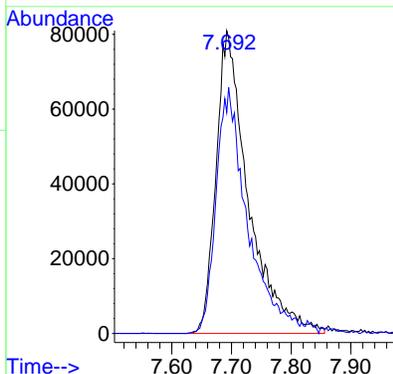
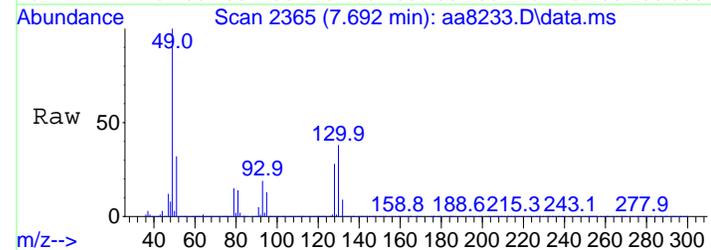
Quant Time: Aug 09 11:24:30 2018
Quant Method : C:\msdchem\1\METHODS\0725.M
Quant Title : TO-15 on the Agilent 7890A / 5975C
QLast Update : Wed Jul 25 14:15:57 2018
Response via : Initial Calibration





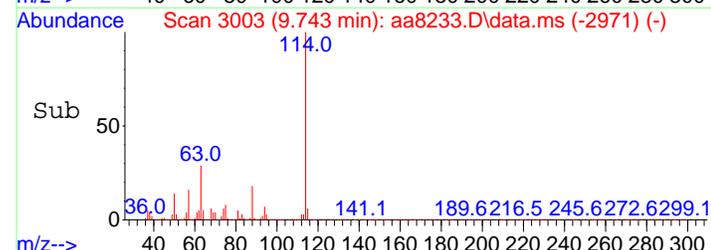
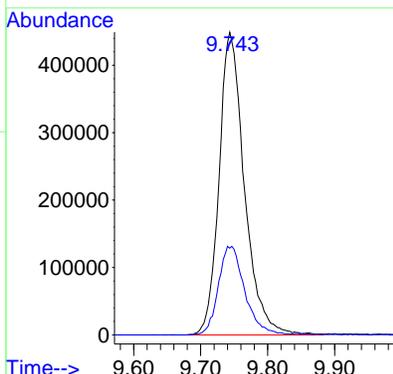
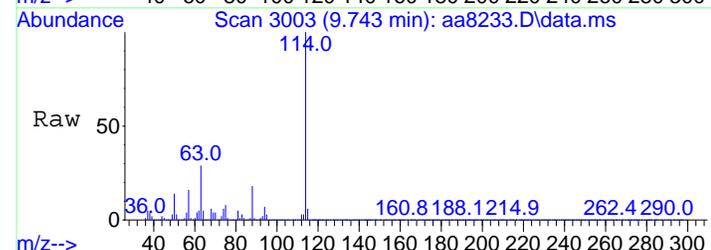
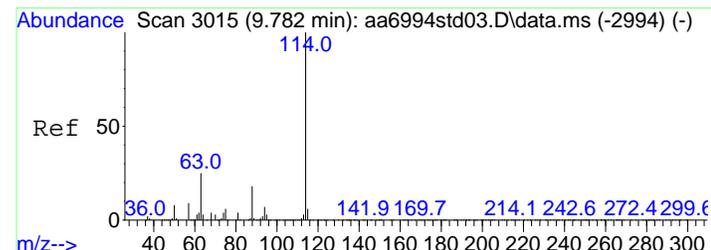
#1
 Bromochloromethane (IS)
 Concen: 10.00 ppbV
 RT: 7.692 min Scan# 2365
 Delta R.T. 0.003 min
 Lab File: aa8233.D
 Acq: 7 Aug 2018 5:31 pm

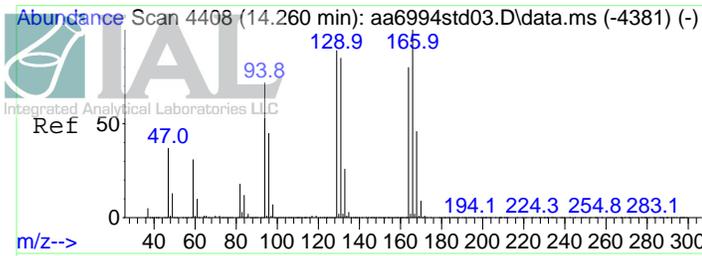
Tgt Ion:130 Resp: 305198
 Ion Ratio Lower Upper
 130 100
 128 79.0 62.6 94.0



#38
 1,4-Difluorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 9.743 min Scan# 3003
 Delta R.T. 0.003 min
 Lab File: aa8233.D
 Acq: 7 Aug 2018 5:31 pm

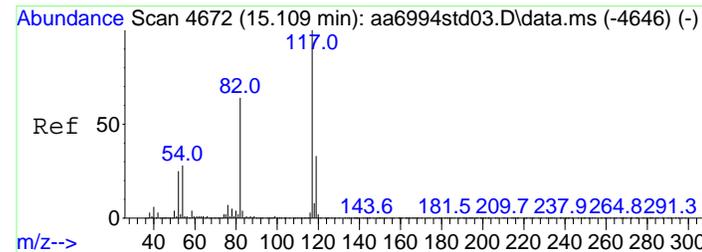
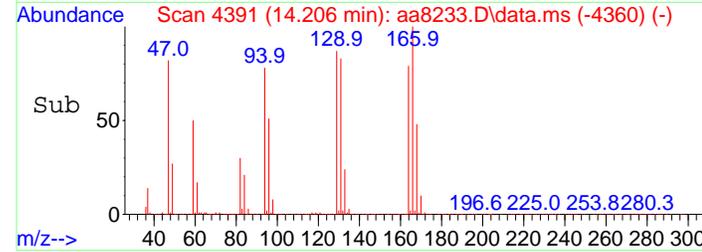
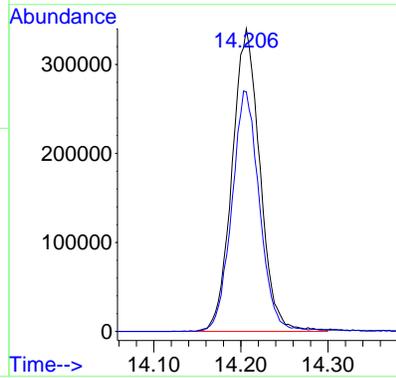
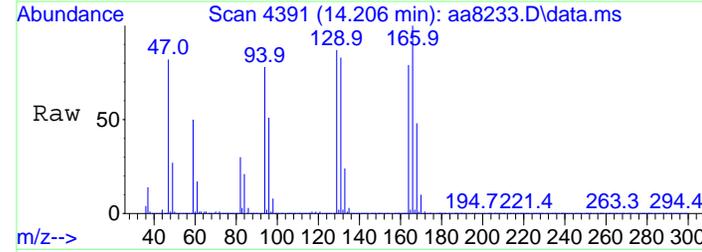
Tgt Ion:114 Resp: 1178297
 Ion Ratio Lower Upper
 114 100
 63 30.5 20.0 30.0#





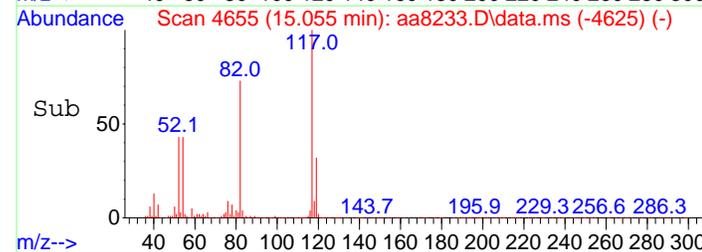
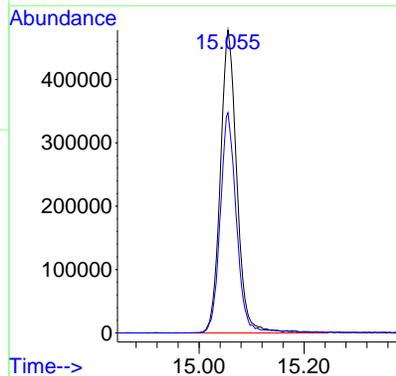
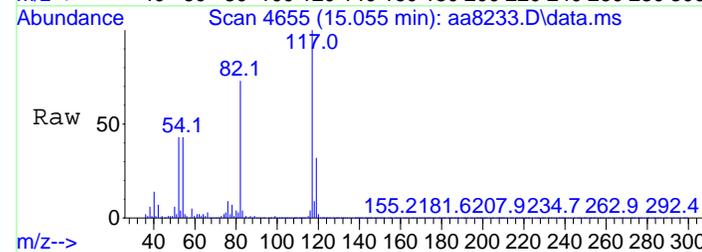
#54
 Tetrachloroethene
 Concen: 13.98 ppbV
 RT: 14.206 min Scan# 4391
 Delta R.T. 0.000 min
 Lab File: aa8233.D
 Acq: 7 Aug 2018 5:31 pm

Tgt Ion:166 Resp: 755419
 Ion Ratio Lower Upper
 166 100
 164 78.8 64.6 96.8

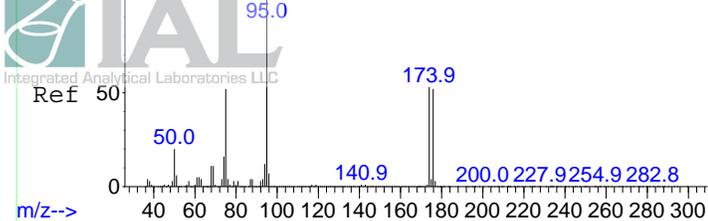


#55
 d-5 Chlorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 15.055 min Scan# 4655
 Delta R.T. -0.003 min
 Lab File: aa8233.D
 Acq: 7 Aug 2018 5:31 pm

Tgt Ion:117 Resp: 1041852
 Ion Ratio Lower Upper
 117 100
 82 71.7 56.0 84.0



Abundance Scan 5355 (17.305 min): aa6994std03.D\data.ms (-5334) (-)



#64

Bromofluorobenzene (tune std)

Concen: 10.98 ppbV

RT: 17.248 min Scan# 5337

Delta R.T. 0.000 min

Lab File: aa8233.D

Acq: 7 Aug 2018 5:31 pm

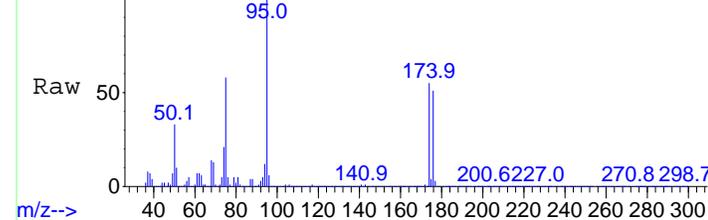
Tgt Ion: 95 Resp: 1043585

Ion Ratio Lower Upper

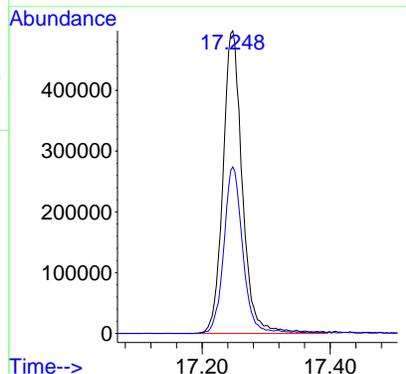
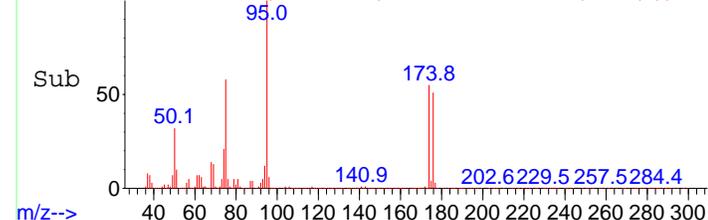
95 100

174 54.0 61.5 92.3#

Abundance Scan 5337 (17.248 min): aa8233.D\data.ms



Abundance Scan 5337 (17.248 min): aa8233.D\data.ms (-5306) (-)



Section VII: Standards Data

Initial Calibration Data

Initial Calibration Verification Data

Continuing Calibration Data

Initial Calibration Data Summary Report

Initial Calibration Curve: 5/18/2018
 Instrument: AA

Method ID: 0518.M

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7071BFB]	05/18/2018 08:17
40 PPBV STD [AA7073STD01]	05/18/2018 09:46
20 PPBV STD [AA7074STD02]	05/18/2018 10:19
10 PPBV STD [AA7075STD03]	05/18/2018 10:53
2 PPBV STD [AA7076STD04]	05/18/2018 12:20
0.2 PPBV STD [AA7077STD05]	05/18/2018 13:03
10 PPBV ICVSS [AA7078ICVSS]	05/18/2018 15:14

RParameter	RRF 0.2ppbv	RRF 2ppbv	RRF 10ppbv	RRF 20ppbv	RRF 40ppbv	Avg ppbv	% RSD
Bromochloromethane	-----ISTD-----						
1,4-Difluorobenzene	-----ISTD-----						
d-5 Chlorobenzene	-----ISTD-----						
Acetone	0.92	0.64	0.84	0.82	0.82	0.81	13
Acrolein	0.48	0.41	0.53	0.56	0.57	0.51	13
Allyl Chloride	0.60	0.44	0.59	0.62	0.59	0.57	12
Benzene	4.9	3.7	4.2	4.4	4.6	4.4	10
Benzyl chloride	0.79	0.83	1.2	1.3	1.4	1.1	26
Bromodichloromethane	0.97	0.72	0.79	0.82	0.83	0.82	11
Bromoform	0.71	0.63	0.60	0.59	0.58	0.62	8.0
Bromomethane	1.2	0.92	1.1	1.1	1.0	1.1	10
1,3-Butadiene	1.2	1.1	1.3	1.3	1.2	1.2	6.3
Chlorobenzene	1.4	1.1	1.00	0.99	0.97	1.1	18
Chloroethane	1.1	0.66	0.78	0.80	0.81	0.82	18
Chloroform	3.6	2.7	3.1	3.2	3.3	3.2	11
Chloromethane	0.64	0.44	0.47	0.47	0.44	0.49	18
Carbon disulfide	4.1	3.3	3.9	3.9	3.8	3.8	8.4
Carbon tetrachloride	3.5	2.8	2.8	2.9	3.1	3.0	10
2-Chlorotoluene	1.6	1.6	1.5	1.5	1.5	1.6	5.0
Cumene	2.1	2.1	1.9	1.8	1.7	1.9	7.9
Cyclohexane	2.6	2.4	2.6	2.6	2.7	2.6	3.9
Dibromochloromethane	0.81	0.63	0.64	0.67	0.69	0.69	10
1,2-Dibromoethane	0.79	0.62	0.64	0.66	0.67	0.68	9.7
1,2-Dichlorobenzene	1.2	1.1	0.98	0.94	0.90	1.0	12
1,3-Dichlorobenzene	1.3	1.1	0.97	0.92	0.85	1.0	18
1,4-Dichlorobenzene	1.3	1.1	0.99	0.95	0.90	1.0	14
Dichlorodifluoromethane	4.1	3.0	3.5	3.5	3.6	3.5	10
1,1-Dichloroethane	3.5	2.4	2.9	3.0	3.1	3.0	13
1,2-Dichloroethane	3.2	2.2	2.5	2.6	2.7	2.7	14
1,1-Dichloroethene	2.8	2.1	2.6	2.7	2.7	2.6	10
1,2-Dichloroethene (trans)	2.3	1.7	2.2	2.3	2.4	2.2	12
1,2-Dichloroethene (trans)	2.5	1.8	2.3	2.4	2.5	2.3	12
1,2-Dichloropropane	0.59	0.44	0.46	0.47	0.48	0.49	12
1,3-Dichloropropene (cis)	0.60	0.49	0.61	0.65	0.67	0.60	11
1,3-Dichloropropene (trans)	0.54	0.46	0.60	0.66	0.70	0.59	17
1,2-Dichlorotetrafluoroethane	4.5	3.5	3.4	3.7	3.6	3.7	12
1,4-Dioxane	0.22	0.17	0.21	0.22	0.22	0.21	9.6
Ethanol	0.91	0.92	0.69	0.72	0.73	0.79	14
Ethylbenzene	1.9	1.9	1.8	1.8	1.7	1.8	4.7
4-Ethyltoluene	1.7	2.0	1.9	1.8	1.7	1.8	6.7
n-Heptane	0.87	0.84	0.82	0.81	0.81	0.83	3.0
1,3-Hexachlorobutadiene	0.89	0.65	0.55	0.49	0.42	0.60	30

*% RSD (Relative Standard Deviation) must be within 30%
 **An exception is made for 2 compounds that must be within 40%
 RRF - Relative Response Factor

Initial Calibration Data Summary Report

Initial Calibration Curve: 5/18/2018
Instrument: AA

Method ID: 0518.M

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7071BFB]	05/18/2018 08:17
40 PPBV STD [AA7073STD01]	05/18/2018 09:46
20 PPBV STD [AA7074STD02]	05/18/2018 10:19
10 PPBV STD [AA7075STD03]	05/18/2018 10:53
2 PPBV STD [AA7076STD04]	05/18/2018 12:20
0.2 PPBV STD [AA7077STD05]	05/18/2018 13:03
10 PPBV ICVSS [AA7078ICVSS]	05/18/2018 15:14

RParameter	RRF 0.2ppbv	RRF 2ppbv	RRF 10ppbv	RRF 20ppbv	RRF 40ppbv	Avg ppbv	% RSD
n-Hexane	2.8	2.2	2.5	2.6	2.7	2.6	9.1
Isopropanol	2.6	2.0	2.7	2.9	2.8	2.6	13
Methylene chloride	3.3	1.9	2.1	2.2	2.2	2.3	25
Methyl ethyl ketone	3.7	2.9	3.7	3.9	4.1	3.7	13
Methyl isobutyl ketone	1.0	1.1	1.1	1.2	1.2	1.1	5.1
Methyl methacrylate	0.54	0.50	0.63	0.68	0.68	0.61	13
Methyl n-butyl ketone	0.85	0.88	1.0	1.1	1.1	1.0	13
Methyl tert-butyl ether	3.3	3.0	3.3	3.5	3.6	3.3	6.8
Naphthalene	0.23	0.21	0.25	0.25	0.24	0.24	8.4
Propene	1.4	1.1	1.2	1.3	1.3	1.3	6.8
Styrene	0.80	0.93	1.0	1.0	0.99	0.94	9.4
Tert-butyl alcohol	2.3	1.9	2.3	2.5	2.6	2.3	12
1,1,2,2-Tetrachloroethane	1.7	1.5	1.2	1.2	1.1	1.3	19
Tetrachloroethene	0.63	0.47	0.45	0.45	0.44	0.49	17
Tetrahydrofuran	1.7	1.3	1.8	1.9	2.0	1.7	17
Toluene	1.2	1.1	1.2	1.2	1.2	1.2	3.9
1,2,4-Trichlorobenzene	1.1	0.72	0.76	0.72	0.65	0.79	22
1,1,1-Trichloroethane	3.5	2.8	2.9	2.9	3.1	3.0	9.3
1,1,2-Trichloroethane	0.56	0.43	0.42	0.43	0.44	0.45	13
Trichloroethene	0.46	0.34	0.37	0.38	0.38	0.38	11
Trichlorofluoromethane	4.7	3.5	3.7	3.7	3.8	3.9	12
1,1,2-Trichloro-1,2,2-trifluoroethane	3.8	3.1	2.7	2.7	2.8	3.0	16
1,2,4-Trimethylbenzene	1.4	1.7	1.7	1.6	1.6	1.6	9.0
1,3,5-Trimethylbenzene	1.5	1.8	1.7	1.6	1.5	1.6	7.9
2,2,4-Trimethylpentane	1.6	1.6	1.9	1.9	1.9	1.8	9.2
Vinyl bromide	0.93	0.87	1.1	1.1	1.2	1.0	12
Vinyl chloride	1.6	1.2	1.5	1.5	1.5	1.5	8.3
Xylenes (m&p)	1.5	1.6	1.4	1.4	1.3	1.4	6.4
Xylenes (o)	1.6	1.7	1.5	1.5	1.4	1.5	7.3

*% RSD (Relative Standard Deviation) must be within 30%

**An exception is made for 2 compounds that must be within 40%

RRF - Relative Response Factor

Method Path : C:\msdchem\1\METHODS\
 Method File : 0518.M
 Title : TO-15 on the Agilent 7890A / 5975C
 Last Update : Fri May 18 13:51:08 2018
 Response Via : Initial Calibration

Calibration Files

0.2 =aa7077std05.D 2 =aa7076std04.D 10 =aa7075std03.D 20 =aa7074std02.D 40 =aa7073std01.D

Compound	0.2	2	10	20	40	Avg	%RSD
-----ISTD-----							
1) I Bromochloromethane...							
2) T Propene	1.359	1.134	1.238	1.269	1.326	1.265	6.89
3) T Dichlorodifluoro...	4.055	3.029	3.478	3.540	3.604	3.541	10.32
4) T Chloromethane	0.643	0.436	0.466	0.471	0.436	0.490	17.76
5) T 1,2-Dichlorotetr...	4.503	3.528	3.375	3.675	3.554	3.727	11.99
6) T Vinyl chloride	1.570	1.246	1.460	1.485	1.488	1.450	8.36
7) T 1,3-Butadiene	1.216	1.110	1.297	1.301	1.229	1.231	6.32
8) T n-Butane	3.596	2.551	2.767	2.725	2.591	2.846	15.07
9) T Bromomethane	1.223	0.923	1.066	1.072	1.022	1.061	10.20
10) T Chloroethane	1.058	0.661	0.778	0.796	0.808	0.820	17.68
11) T Ethanol	0.905	0.917	0.692	0.716	0.732	0.792	13.79
12) T Vinyl bromide	0.934	0.872	1.090	1.122	1.155	1.035	12.04
13) T Acrolein	0.477	0.414	0.534	0.558	0.567	0.510	12.61
14) T Acetone	0.919	0.643	0.838	0.823	0.820	0.808	12.51
15) T Trichlorofluorom...	4.726	3.516	3.676	3.730	3.782	3.886	12.35
16) T Isopropanol	2.617	2.047	2.741	2.877	2.777	2.612	12.59
17) T n-Pentane	3.233	2.582	3.095	3.157	3.239	3.061	8.97
18) T 1,1-Dichloroethene	2.773	2.111	2.608	2.678	2.675	2.569	10.21
19) T Methylene chloride	3.342	1.870	2.114	2.150	2.197	2.334	24.73
20) T Tert-butyl alcohol	2.258	1.905	2.337	2.549	2.647	2.339	12.36
21) T Allyl chloride	0.600	0.443	0.592	0.616	0.589	0.568	12.48
22) T 1,1,2-Trichloro-...	3.820	3.060	2.700	2.679	2.770	3.006	15.97
23) T Carbon disulfide	4.116	3.252	3.859	3.853	3.780	3.772	8.42
24) T 1,2-Dichloroethe...	2.487	1.818	2.292	2.382	2.461	2.288	11.94
25) T 1,1-Dichloroethane	3.471	2.434	2.867	2.974	3.060	2.961	12.60
26) T Methyl tert-butyl...	3.277	2.996	3.322	3.471	3.600	3.333	6.82
27) T Methyl ethyl ketone	3.732	2.882	3.689	3.914	4.100	3.663	12.73
28) T 1,2-Dichloroethe...	2.299	1.732	2.228	2.331	2.418	2.201	12.33
29) T Ethyl acetate	0.565	0.462	0.606	0.650	0.680	0.593	14.34
30) T n-Hexane	2.784	2.199	2.503	2.620	2.738	2.569	9.11
31) T Chloroform	3.641	2.658	3.074	3.165	3.279	3.163	11.23
32) T Tetrahydrofuran	1.712	1.270	1.777	1.914	2.015	1.738	16.51
33) T 1,2-Dichloroethane	3.197	2.159	2.531	2.630	2.742	2.652	14.15
34) T 1,1,1-Trichloroe...	3.526	2.849	2.852	2.923	3.058	3.042	9.33
35) T Benzene	4.941	3.722	4.240	4.420	4.583	4.381	10.27
36) T Carbon tetrachlo...	3.503	2.750	2.789	2.895	3.066	3.001	10.21
37) T Cyclohexane	2.640	2.432	2.600	2.638	2.708	2.604	3.98
-----ISTD-----							
38) I 1,4-Difluorobenzen...							
39) T 1,2-Dichloropropane	0.588	0.442	0.455	0.470	0.477	0.487	11.97
40) T Bromodichloromet...	0.967	0.717	0.792	0.816	0.829	0.824	11.00
41) T 2,2,4-Trimethylp...	1.629	1.601	1.894	1.923	1.935	1.796	9.28
42) T Trichloroethene	0.458	0.339	0.372	0.377	0.376	0.384	11.43
43) T 1,4-Dioxane	0.220	0.174	0.205	0.217	0.222	0.208	9.64
44) T Methyl methacrylate	0.542	0.502	0.631	0.675	0.684	0.607	13.40
45) T n-Heptane	0.869	0.835	0.821	0.809	0.808	0.828	3.04
46) T cis-1,3-Dichloro...	0.599	0.492	0.608	0.649	0.670	0.604	11.40
47) T Methyl isobutyl ...	1.048	1.075	1.141	1.178	1.170	1.122	5.16
48) T trans-1,3-Dichlo...	0.539	0.458	0.604	0.664	0.704	0.594	16.56
49) T 1,1,2-Trichloroe...	0.560	0.425	0.418	0.429	0.436	0.454	13.20
50) T Toluene	1.242	1.117	1.154	1.192	1.196	1.180	3.99
51) T Methyl n-butyl k...	0.845	0.884	1.041	1.111	1.130	1.002	13.03
52) T Dibromochloromet...	0.807	0.633	0.642	0.666	0.686	0.687	10.24
53) T 1,2-Dibromoethane	0.787	0.617	0.639	0.660	0.672	0.675	9.76
54) T Tetrachloroethene	0.633	0.474	0.446	0.448	0.440	0.488	16.80
-----ISTD-----							
55) I d-5 Chlorobenzene ...							
56) T Chlorobenzene	1.435	1.082	0.996	0.985	0.969	1.094	17.92
57) T Ethylbenzene	1.941	1.912	1.836	1.796	1.723	1.841	4.77
58) T Xylenes (m&p)	1.461	1.554	1.433	1.383	1.305	1.427	6.47
59) T Bromoform	0.706	0.630	0.597	0.594	0.584	0.622	8.01
60) T Styrene	0.795	0.926	1.000	1.006	0.990	0.943	9.41

ALSDG #E18-06141

Method Path : C:\msdchem\1\METHODS\
Method File : 0518.M

Peak #	Retention Time (min)	Response Factor	Area	Height	Width	Height	Area	Height
61)	Xylene (o)	1.567	1.670	1.521	1.450	1.380	1.518	7.32
62)	1,1,2,2-Tetrachl...	1.692	1.454	1.220	1.160	1.099	1.325	18.50
63)	n-Nonane	1.365	1.687	1.537	1.453	1.351	1.479	9.36
64)	Bromofluorobenze...	0.915	0.929	0.920	0.914	0.924	0.921	0.67
65)	Cumene	2.056	2.051	1.893	1.812	1.710	1.904	7.92
66)	2-Chlorotoluene	1.646	1.633	1.546	1.520	1.460	1.561	5.02
67)	n-Propyl benzene	2.546	2.954	2.616	2.467	2.278	2.572	9.63
68)	4-Ethyltoluene	1.683	1.957	1.855	1.772	1.672	1.788	6.73
69)	1,3,5-Trimethylb...	1.522	1.829	1.697	1.607	1.522	1.635	7.97
70)	1,2,4-Trimethylb...	1.362	1.725	1.692	1.641	1.565	1.597	9.04
71)	Benzyl chloride	0.787	0.826	1.164	1.321	1.397	1.099	25.53
72)	1,3-Dichlorobenzene	1.318	1.087	0.968	0.919	0.853	1.029	17.76
73)	1,4-Dichlorobenzene	1.264	1.094	0.993	0.954	0.904	1.042	13.66
74)	1,2-Dichlorobenzene	1.208	1.058	0.975	0.942	0.897	1.016	12.04
75)	1,2,4-Trichlorob...	1.094	0.719	0.760	0.723	0.651	0.789	22.12
76)	Naphthalene	0.227	0.206	0.253	0.253	0.238	0.235	8.44
77)	1,3-Hexachlorobu...	0.887	0.652	0.546	0.487	0.423	0.599	30.35

(#) = Out of Range

Data Path : C:\DATA\05-18-18\
 Data File : aa7073std01.D
 Acq On : 18 May 2018 9:46 am
 Operator : jls
 Sample : 40 ppbv Std
 Misc : CC483586
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: May 18 11:37:38 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 11:37:26 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.759	130	497626	10.00	ppbV	0.04	
38) 1,4-Difluorobenzene (IS)	9.788	114	2152735	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.113	117	1991801	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...	17.309	95	1841417	10.08	ppbV	0.00	
Target Compounds							
							Qvalue
2) Propene	3.541	41	2903235	46.55	ppbV		82
3) Dichlorodifluoromethane	3.592	85	7890221	45.18	ppbV		97
4) Chloromethane	3.795	52	953559	40.90	ppbV #		1
5) 1,2-Dichlorotetrafluor...	3.760	85	7073452m	44.67	ppbV		
6) Vinyl chloride	3.882	62	3257860	44.46	ppbV #		50
7) 1,3-Butadiene	3.972	54	2690988	41.62	ppbV		100
8) n-Butane	4.023	43	5672952	41.52	ppbV		93
9) Bromomethane	4.277	94	2014205	37.86	ppbV		97
10) Chloroethane	4.374	64	1608730	41.06	ppbV		92
11) Ethanol	4.399	45	1369985	39.11	ppbV		96
12) Vinyl bromide	4.673	106	2529945	45.96	ppbV		99
13) Acrolein	4.737	56	1355476	49.86	ppbV		96
14) Acetone	4.827	58	1794562	43.43	ppbV		85
15) Trichlorofluoromethane	5.017	101	8281215	44.94	ppbV		100
16) Isopropanol	5.014	45	5527707	39.55	ppbV #		53
17) n-Pentane	5.313	43	7090927	45.58	ppbV		95
18) 1,1-Dichloroethene	5.573	61	5856653	44.53	ppbV		90
19) Methylene chloride	5.689	49	4810194	45.35	ppbV		100
20) Tert-butyl alcohol	5.538	59	6323555	52.01	ppbV		100
21) Allyl chloride	5.788	76	1289380	42.90	ppbV		100
22) 1,1,2-Trichloro-1,2,2-...	5.911	101	6064592	45.32	ppbV		87
23) Carbon disulfide	5.975	76	8276326	43.13	ppbV		99
24) 1,2-Dichloroethene (tr...	6.544	61	5387714	46.32	ppbV		100
25) 1,1-Dichloroethane	6.737	63	6699550	46.10	ppbV		98
26) Methyl tert-butyl ether	6.759	73	7881704	46.63	ppbV		96
27) Methyl ethyl ketone	7.075	43	8976908	47.45	ppbV		100
28) 1,2-Dichloroethene (cis)	7.554	61	5293956	46.67	ppbV		99
29) Ethyl acetate	7.743	45	1489617	47.68	ppbV		100
30) n-Hexane	7.772	57	5996014	47.04	ppbV		100
31) Chloroform	7.853	83	7180003	46.26	ppbV		100
32) Tetrahydrofuran	8.238	42	4814098	52.43	ppbV		92
33) 1,2-Dichloroethane	8.621	62	6004482	46.77	ppbV		99
34) 1,1,1-Trichloroethane	8.901	97	6696335	46.61	ppbV		100
35) Benzene	9.396	78	10034352	46.57	ppbV		99
36) Carbon tetrachloride	9.563	117	6712610	47.47	ppbV		99
37) Cyclohexane	9.708	56	5929762	45.50	ppbV		90
39) 1,2-Dichloropropane	10.306	63	4520946	45.38	ppbV		100
40) Bromodichloromethane	10.528	83	7848057	45.36	ppbV		100
41) 2,2,4-Trimethylpentane	10.640	57	18330470	44.62	ppbV		96
42) Trichloroethene	10.582	130	3561175	44.20	ppbV		96
43) 1,4-Dioxane	10.544	88	1909558	42.01	ppbV		98
44) Methyl methacrylate	10.795	41	6478787	46.07	ppbV		98
45) n-Heptane	10.936	43	7654967	43.62	ppbV		96
46) cis-1,3-Dichloropropene	11.595	75	6350305	46.94	ppbV		100
47) Methyl isobutyl ketone	11.611	43	11084002	44.42	ppbV		97
48) trans-1,3-Dichloropropene	12.229	75	6063832	44.44	ppbV		98
49) 1,1,2-Trichloroethane	12.441	97	4128347	45.27	ppbV		90
50) Toluene	12.798	91	11327617	44.86	ppbV		99
51) Methyl n-butyl ketone	13.113	43	10706608	46.24	ppbV		97

AL SDG #E18-06141

Data Path : C:\DATA\05-18-18\
 Data File : aa7073std01.D
 Acq On : 18 May 2018 9:46 am
 Operator : jls
 Sample : 40 ppbv Std
 Misc : CC483586
 ALS Vial : 3 Sample Multiplier: 1

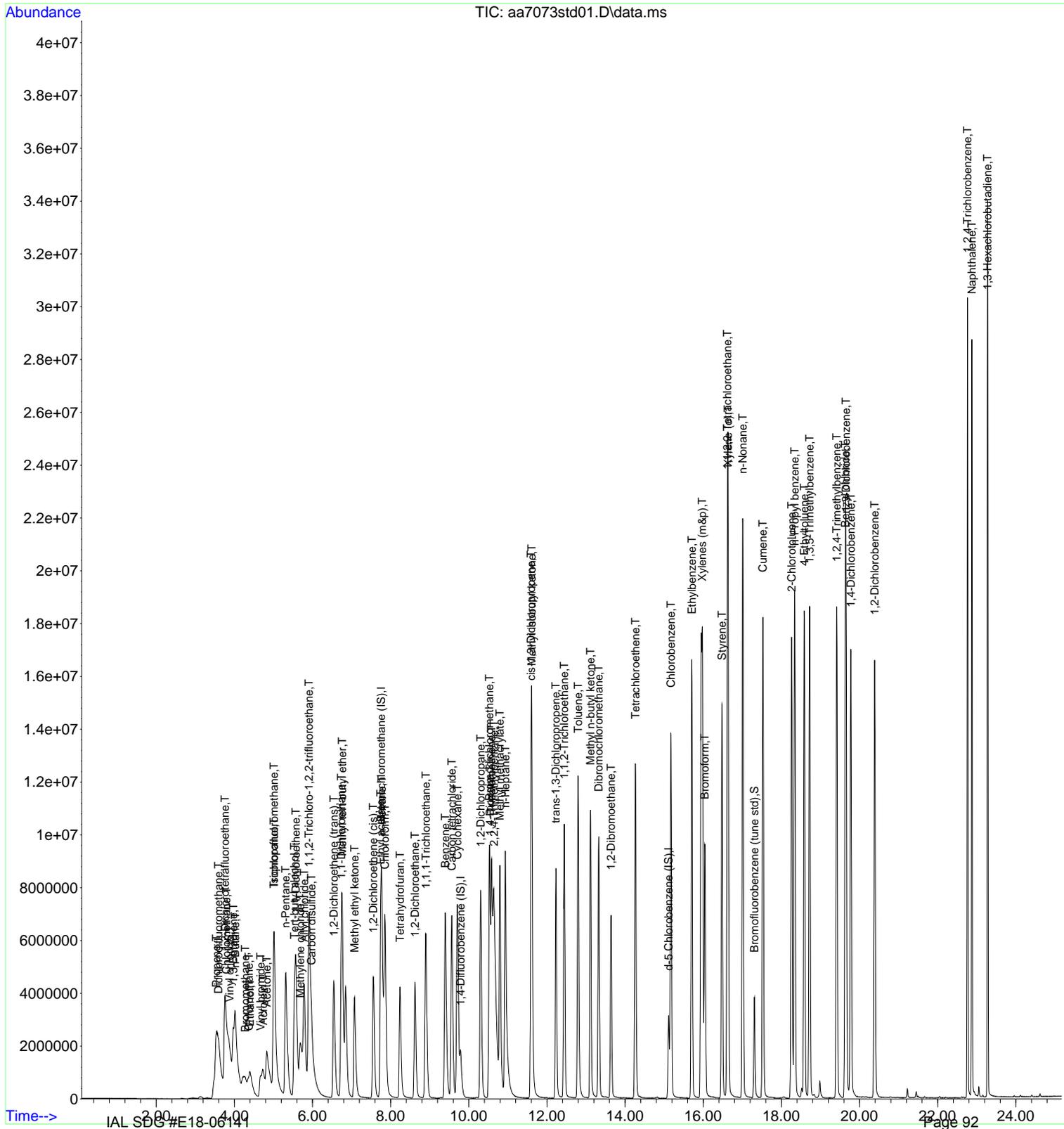
Quant Time: May 18 11:37:38 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 11:37:26 2018
 Response via : Initial Calibration

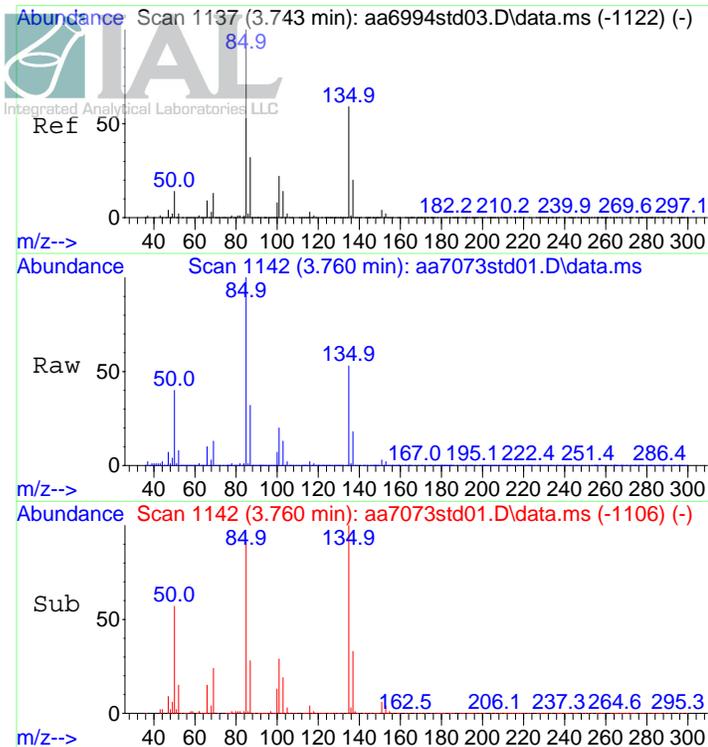
Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
52) Dibromochloromethane	13.325	129	6497747	46.17	ppbV	100
53) 1,2-Dibromoethane	13.640	107	6368628	45.54	ppbV	99
54) Tetrachloroethene	14.264	166	4164333	43.29	ppbV	98
56) Chlorobenzene	15.171	112	8496608	43.06	ppbV	98
57) Ethylbenzene	15.708	91	15102551	41.76	ppbV	97
58) Xylenes (m&p)	15.975	91	22878040	81.58	ppbV	97
59) Bromoform	16.039	173	5118623	43.16	ppbV	98
60) Styrene	16.479	104	8673250	43.41	ppbV	99
61) Xylene (o)	16.637	91	12096421	40.88	ppbV	97
62) 1,1,2,2-Tetrachloroethane	16.618	83	9635267	40.65	ppbV	99
63) n-Nonane	17.010	43	11843056	39.77	ppbV	92
65) Cumene	17.528	105	14984429	40.62	ppbV	100
66) 2-Chlorotoluene	18.261	91	12791930	41.89	ppbV	99
67) n-Propyl benzene	18.338	91	19967781	39.44	ppbV	98
68) 4-Ethyltoluene	18.585	105	14654612	40.57	ppbV	98
69) 1,3,5-Trimethylbenzene	18.724	105	13337284	40.54	ppbV	98
70) 1,2,4-Trimethylbenzene	19.415	105	13717500	41.33	ppbV	100
71) Benzyl chloride	19.634	91	11131592	44.97	ppbV	98
72) 1,3-Dichlorobenzene	19.653	146	7471637	39.75	ppbV	100
73) 1,4-Dichlorobenzene	19.778	146	7924734	40.86	ppbV	100
74) 1,2-Dichlorobenzene	20.386	146	7865403	41.19	ppbV	99
75) 1,2,4-Trichlorobenzene	22.762	180	6225065	42.16	ppbV	99
76) Naphthalene	22.875	127	2274212	45.09	ppbV	100
77) 1,3-Hexachlorobutadiene	23.276	225	3704125	36.00	ppbV	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\05-18-18\
 Data File : aa7073std01.D
 Acq On : 18 May 2018 9:46 am
 Operator : jls
 Sample : 40 ppbv Std
 Misc : CC483586
 ALS Vial : 3 Sample Multiplier: 1

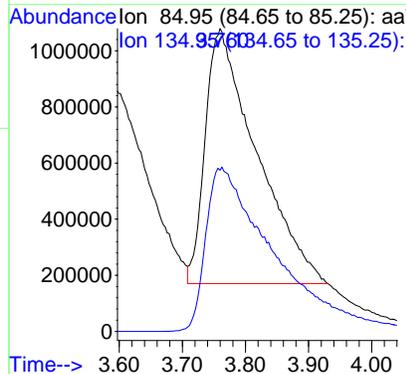
Quant Time: May 18 11:37:38 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 11:37:26 2018
 Response via : Initial Calibration



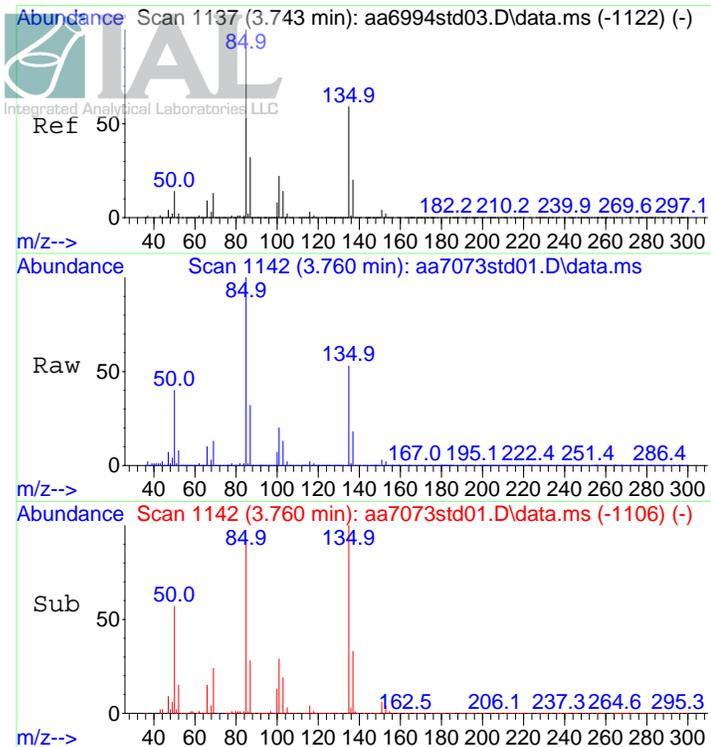


#5
 1,2-Dichlorotetrafluoroethane
 Concen: 32.24 ppbV
 RT: 3.760 min Scan# 1142
 Delta R.T. 0.016 min
 Lab File: aa7073std01.D
 Acq: 18 May 2018 9:46 am

Tgt Ion: 85 Resp: 5105732
 Ion Ratio Lower Upper
 85 100
 135 78.7 46.4 69.6#

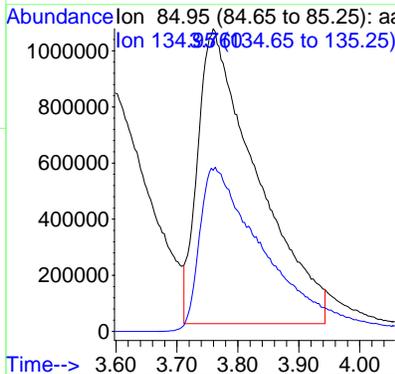


Before manual integration for 1, 2-Dichlorotetrafluoroethane.



#5
 1,2-Dichlorotetrafluoroethane
 Concen: 44.67 ppbV m
 RT: 3.760 min Scan# 1142
 Delta R.T. 0.016 min
 Lab File: aa7073std01.D
 Acq: 18 May 2018 9:46 am

Tgt Ion: 85 Resp: 7073452
 Ion Ratio Lower Upper
 85 100
 135 56.8 46.4 69.6



After manual integration for 1, 2-Dichlorotetrafluoroethane.

Data Path : C:\DATA\05-18-18\
 Data File : aa7074std02.D
 Acq On : 18 May 2018 10:19 am
 Operator : jls
 Sample : 20 ppbv Std
 Misc : CC483586
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: May 18 11:37:08 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 11:36:21 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.740	130	531970	10.00	ppbV	0.02	
38) 1,4-Difluorobenzene (IS)	9.785	114	2238046	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.112	117	2008933	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...	17.305	95	1835985	9.93	ppbV	0.00	
Target Compounds							
							Qvalue
2) Propene	3.512	41	1484720	22.54	ppbV		83
3) Dichlorodifluoromethane	3.541	85	4143365	22.39	ppbV		97
4) Chloromethane	3.756	52	551651	22.27	ppbV		89
5) 1,2-Dichlorotetrafluor...	3.753	85	3909499m	21.78	ppbV		
6) Vinyl chloride	3.869	62	1738287	22.38	ppbV		87
7) 1,3-Butadiene	3.968	54	1522794	22.06	ppbV		99
8) n-Butane	4.013	43	3189269	21.67	ppbV		95
9) Bromomethane	4.232	94	1129084	19.90	ppbV		98
10) Chloroethane	4.364	64	847208	20.46	ppbV		92
11) Ethanol	4.402	45	716152	19.46	ppbV		98
12) Vinyl bromide	4.666	106	1313405	22.65	ppbV		100
13) Acrolein	4.727	56	712750	25.08	ppbV		96
14) Acetone	4.820	58	963285	21.62	ppbV		98
15) Trichlorofluoromethane	5.007	101	4365295	22.32	ppbV		100
16) Isopropanol	5.004	45	3060927	20.99	ppbV #		90
17) n-Pentane	5.309	43	3694806	22.44	ppbV		94
18) 1,1-Dichloroethene	5.569	61	3134080	22.59	ppbV		91
19) Methylene chloride	5.685	49	2515790	22.37	ppbV		99
20) Tert-butyl alcohol	5.537	59	3254922	26.18	ppbV		100
21) Allyl chloride	5.778	76	721129	22.91	ppbV		100
22) 1,1,2-Trichloro-1,2,2-...	5.910	101	3134896	21.83	ppbV		88
23) Carbon disulfide	5.971	76	4508716	21.96	ppbV		100
24) 1,2-Dichloroethene (tr...	6.540	61	2787665	22.86	ppbV		100
25) 1,1-Dichloroethane	6.730	63	3481144	22.83	ppbV		98
26) Methyl tert-butyl ether	6.759	73	4062472	22.99	ppbV		96
27) Methyl ethyl ketone	7.071	43	4581098	23.35	ppbV		100
28) 1,2-Dichloroethene (cis)	7.553	61	2728274	23.02	ppbV		99
29) Ethyl acetate	7.743	45	760229	23.58	ppbV		100
30) n-Hexane	7.772	57	3066330	23.03	ppbV		100
31) Chloroform	7.852	83	3703561	22.65	ppbV		99
32) Tetrahydrofuran	8.235	42	2443244	25.85	ppbV		92
33) 1,2-Dichloroethane	8.621	62	3077517	22.86	ppbV		99
34) 1,1,1-Trichloroethane	8.894	97	3420329	22.54	ppbV		100
35) Benzene	9.396	78	5172999	22.93	ppbV		100
36) Carbon tetrachloride	9.563	117	3388067	22.84	ppbV		100
37) Cyclohexane	9.708	56	3087138	22.32	ppbV		89
39) 1,2-Dichloropropane	10.302	63	2315306	22.72	ppbV		100
40) Bromodichloromethane	10.524	83	4016130	22.66	ppbV		99
41) 2,2,4-Trimethylpentane	10.646	57	9467560	22.34	ppbV		97
42) Trichloroethene	10.579	130	1855876	22.32	ppbV		96
43) 1,4-Dioxane	10.540	88	973335	21.23	ppbV		97
44) Methyl methacrylate	10.794	41	3323272	23.52	ppbV		100
45) n-Heptane	10.933	43	3984310	21.68	ppbV		97
46) cis-1,3-Dichloropropene	11.588	75	3193269	23.46	ppbV		100
47) Methyl isobutyl ketone	11.611	43	5798007	22.71	ppbV		99
48) trans-1,3-Dichloropropene	12.232	75	2972169	22.00	ppbV		98
49) 1,1,2-Trichloroethane	12.441	97	2111440	22.54	ppbV		89
50) Toluene	12.794	91	5867864	22.72	ppbV		98
51) Methyl n-butyl ketone	13.109	43	5468033	23.47	ppbV		99

AL SDG #E18-06141

Data Path : C:\DATA\05-18-18\
 Data File : aa7074std02.D
 Acq On : 18 May 2018 10:19 am
 Operator : jls
 Sample : 20 ppbv Std
 Misc : CC483586
 ALS Vial : 4 Sample Multiplier: 1

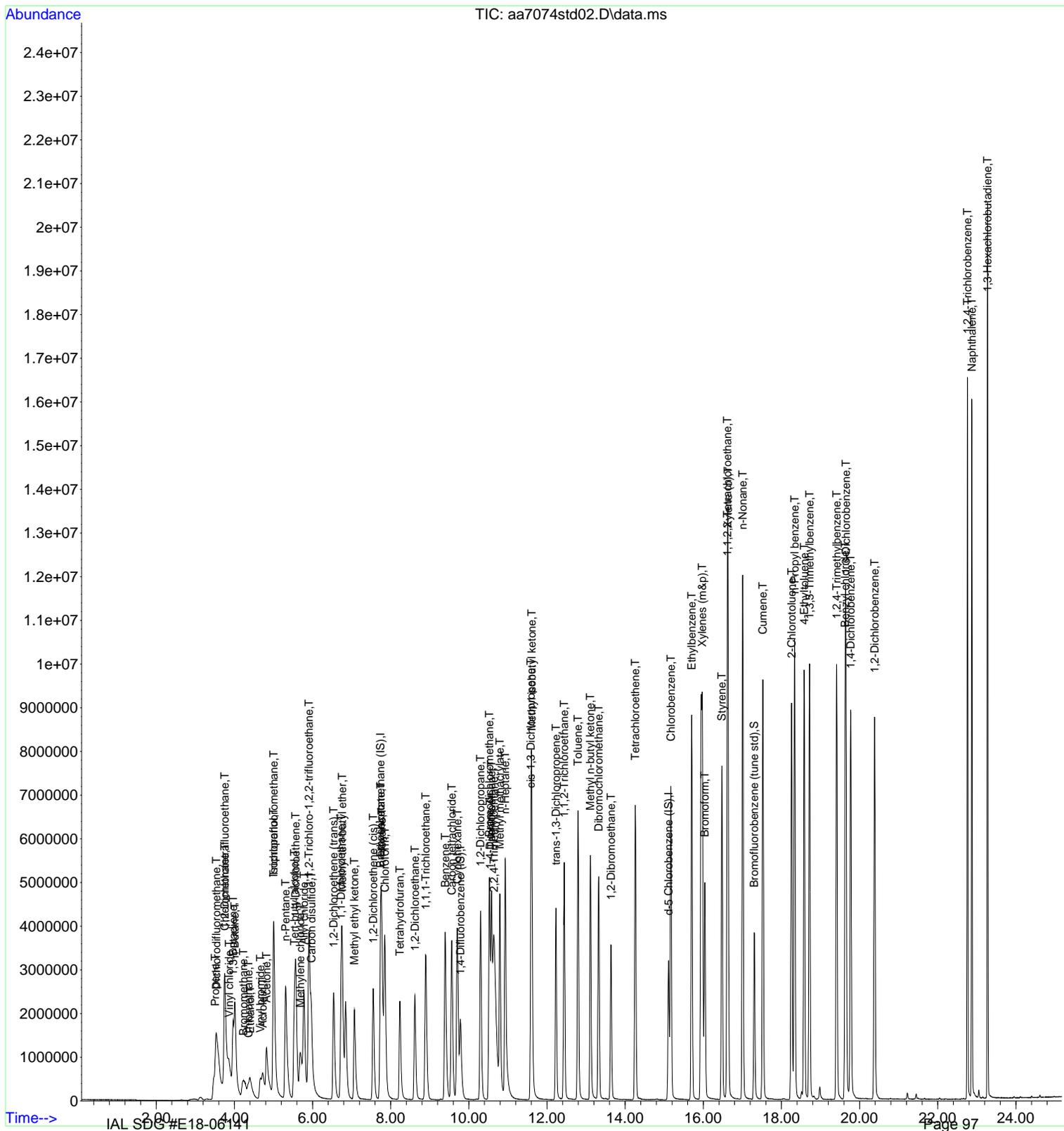
Quant Time: May 18 11:37:08 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 11:36:21 2018
 Response via : Initial Calibration

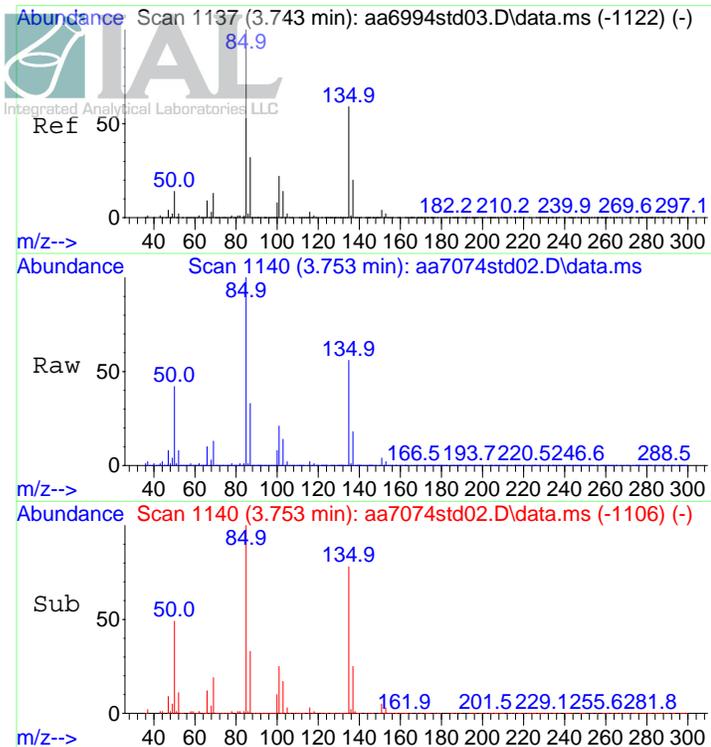
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Dibromochloromethane	13.322	129	3279041	22.84	ppbV	99
53) 1,2-Dibromoethane	13.637	107	3249507	22.71	ppbV	100
54) Tetrachloroethene	14.260	166	2203864	22.07	ppbV	98
56) Chlorobenzene	15.170	112	4355563	21.77	ppbV	99
57) Ethylbenzene	15.704	91	7936994	21.52	ppbV	96
58) Xylenes (m&p)	15.971	91	12225755	42.48	ppbV #	28
59) Bromoform	16.038	173	2626421	21.91	ppbV	97
60) Styrene	16.476	104	4444465	22.12	ppbV	98
61) Xylene (o)	16.633	91	6406380	20.96	ppbV	96
62) 1,1,2,2-Tetrachloroethane	16.614	83	5128408	20.93	ppbV	98
63) n-Nonane	17.009	43	6423191	20.81	ppbV	96
65) Cumene	17.524	105	8006340	21.06	ppbV	98
66) 2-Chlorotoluene	18.260	91	6717849	21.63	ppbV	98
67) n-Propyl benzene	18.337	91	10903452	20.74	ppbV	96
68) 4-Ethyltoluene	18.582	105	7829590	21.00	ppbV	97
69) 1,3,5-Trimethylbenzene	18.720	105	7100443	20.83	ppbV	100
70) 1,2,4-Trimethylbenzene	19.411	105	7252590	21.34	ppbV	99
71) Benzyl chloride	19.630	91	5309178	22.70	ppbV	99
72) 1,3-Dichlorobenzene	19.649	146	4062553	20.88	ppbV	99
73) 1,4-Dichlorobenzene	19.775	146	4215661	21.12	ppbV	99
74) 1,2-Dichlorobenzene	20.386	146	4163747	21.25	ppbV	100
75) 1,2,4-Trichlorobenzene	22.762	180	3486106	22.85	ppbV	98
76) Naphthalene	22.874	127	1222014	24.04	ppbV	100
77) 1,3-Hexachlorobutadiene	23.276	225	2152097	19.62	ppbV	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\05-18-18\
Data File : aa7074std02.D
Acq On : 18 May 2018 10:19 am
Operator : jls
Sample : 20 ppbv Std
Misc : CC483586
ALS Vial : 4 Sample Multiplier: 1

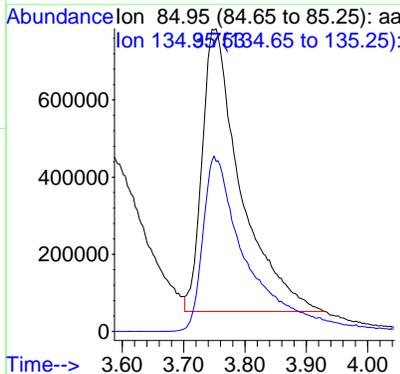
Quant Time: May 18 11:37:08 2018
Quant Method : C:\msdchem\1\METHODS\0518.M
Quant Title : TO-15 on the Agilent 7890A / 5975C
QLast Update : Fri May 18 11:36:21 2018
Response via : Initial Calibration



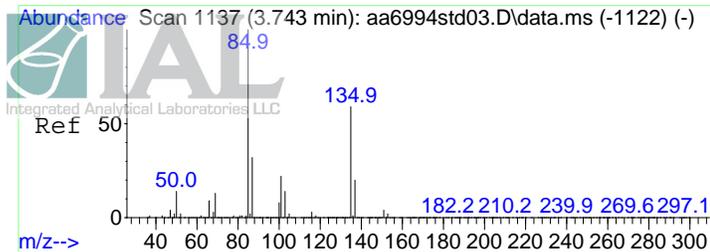


#5
 1,2-Dichlorotetrafluoroethane
 Concen: 17.72 ppbV
 RT: 3.753 min Scan# 1140
 Delta R.T. 0.010 min
 Lab File: aa7074std02.D
 Acq: 18 May 2018 10:19 am

Tgt Ion: 85 Resp: 3180717
 Ion Ratio Lower Upper
 85 100
 135 68.7 46.4 69.6

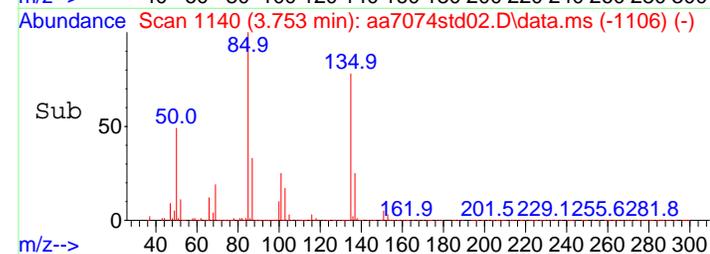
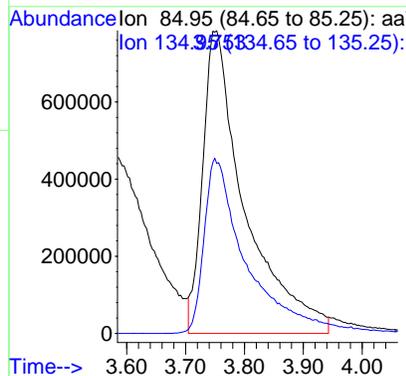
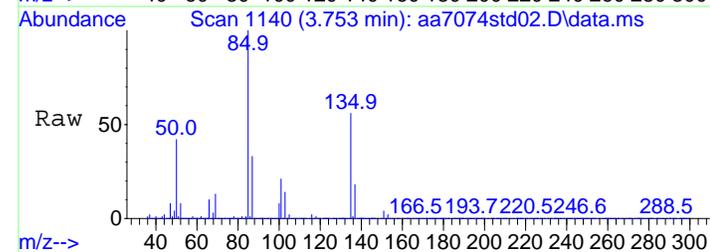


Before manual integration for 1, 2-Dichlorotetrafluoroethane.



#5
 1,2-Dichlorotetrafluoroethane
 Concen: 21.78 ppbV m
 RT: 3.753 min Scan# 1140
 Delta R.T. 0.010 min
 Lab File: aa7074std02.D
 Acq: 18 May 2018 10:19 am

Tgt Ion: 85 Resp: 3909499
 Ion Ratio Lower Upper
 85 100
 135 55.9 46.4 69.6



After manual integration for 1, 2-Dichlorotetrafluoroethane.

Data Path : C:\DATA\05-18-18\
 Data File : aa7075std03.D
 Acq On : 18 May 2018 10:53 am
 Operator : jls
 Sample : 10 ppbv Std
 Misc : CC483586
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: May 18 11:26:02 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 11:25:51 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.724	130	543623	10.00	ppbV	0.00	
38) 1,4-Difluorobenzene (IS)	9.782	114	2258745	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.109	117	1967114	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...	17.309	95	1810039	10.00	ppbV	0.00	
Target Compounds							
							Qvalue
2) Propene	3.483	41	740351	11.05	ppbV		83
3) Dichlorodifluoromethane	3.531	85	2080034	11.01	ppbV		97
4) Chloromethane	3.734	52	278411	11.00	ppbV		90
5) 1,2-Dichlorotetrafluor...	3.743	85	1834502	10.01	ppbV		95
6) Vinyl chloride	3.862	62	872965	11.01	ppbV		87
7) 1,3-Butadiene	3.965	54	775850	11.01	ppbV #		64
8) n-Butane	4.013	43	1654335	11.01	ppbV		95
9) Bromomethane	4.229	94	573956	9.95	ppbV		97
10) Chloroethane	4.357	64	423149	9.97	ppbV		93
11) Ethanol	4.396	45	353572	9.39	ppbV		96
12) Vinyl bromide	4.666	106	651915	11.11	ppbV		100
13) Acrolein	4.724	56	348524	12.01	ppbV		93
14) Acetone	4.820	58	500879	11.12	ppbV #		53
15) Trichlorofluoromethane	5.004	101	2198346	11.01	ppbV		99
16) Isopropanol	5.000	45	1490063	10.02	ppbV		98
17) n-Pentane	5.306	43	1850884	11.01	ppbV		95
18) 1,1-Dichloroethene	5.566	61	1559392	11.01	ppbV		91
19) Methylene chloride	5.682	49	1263982	11.06	ppbV #		74
20) Tert-butyl alcohol	5.531	59	1524846	12.01	ppbV		100
21) Allyl chloride	5.779	76	353885	11.07	ppbV		100
22) 1,1,2-Trichloro-1,2,2-...	5.907	101	1614484	11.01	ppbV		88
23) Carbon disulfide	5.965	76	2307921	11.01	ppbV #		85
24) 1,2-Dichloroethene (tr...	6.537	61	1370869	11.01	ppbV		100
25) 1,1-Dichloroethane	6.727	63	1714370	11.01	ppbV		98
26) Methyl tert-butyl ether	6.759	73	1986726	11.01	ppbV		96
27) Methyl ethyl ketone	7.071	43	2205827	11.01	ppbV #		89
28) 1,2-Dichloroethene (cis)	7.550	61	1332135	11.01	ppbV		100
29) Ethyl acetate	7.740	45	362368	11.01	ppbV		100
30) n-Hexane	7.766	57	1496602	11.01	ppbV #		68
31) Chloroform	7.846	83	1837950	11.01	ppbV		99
32) Tetrahydrofuran	8.235	42	1158913	12.02	ppbV		91
33) 1,2-Dichloroethane	8.618	62	1513276	11.01	ppbV		100
34) 1,1,1-Trichloroethane	8.891	97	1705454	11.01	ppbV		100
35) Benzene	9.389	78	2535720	11.01	ppbV #		91
36) Carbon tetrachloride	9.560	117	1667638	11.01	ppbV		99
37) Cyclohexane	9.704	56	1554884	11.08	ppbV		89
39) 1,2-Dichloropropane	10.299	63	1131231	11.00	ppbV		100
40) Bromodichloromethane	10.524	83	1967277	11.01	ppbV		99
41) 2,2,4-Trimethylpentane	10.637	57	4705769	10.97	ppbV		97
42) Trichloroethene	10.579	130	923228	11.00	ppbV		96
43) 1,4-Dioxane	10.544	88	462639	10.00	ppbV #		75
44) Methyl methacrylate	10.794	41	1568858	11.00	ppbV #		65
45) n-Heptane	10.929	43	2039918	11.00	ppbV #		78
46) cis-1,3-Dichloropropene	11.589	75	1511338	11.00	ppbV		100
47) Methyl isobutyl ketone	11.611	43	2834046	11.00	ppbV #		79
48) trans-1,3-Dichloropropene	12.228	75	1363556	10.00	ppbV #		63
49) 1,1,2-Trichloroethane	12.441	97	1039768	11.00	ppbV		91
50) Toluene	12.794	91	2867532	11.00	ppbV		98
51) Methyl n-butyl ketone	13.113	43	2586043	11.00	ppbV #		76

AL SDG #E18-06141

Data Path : C:\DATA\05-18-18\
 Data File : aa7075std03.D
 Acq On : 18 May 2018 10:53 am
 Operator : jls
 Sample : 10 ppbv Std
 Misc : CC483586
 ALS Vial : 5 Sample Multiplier: 1

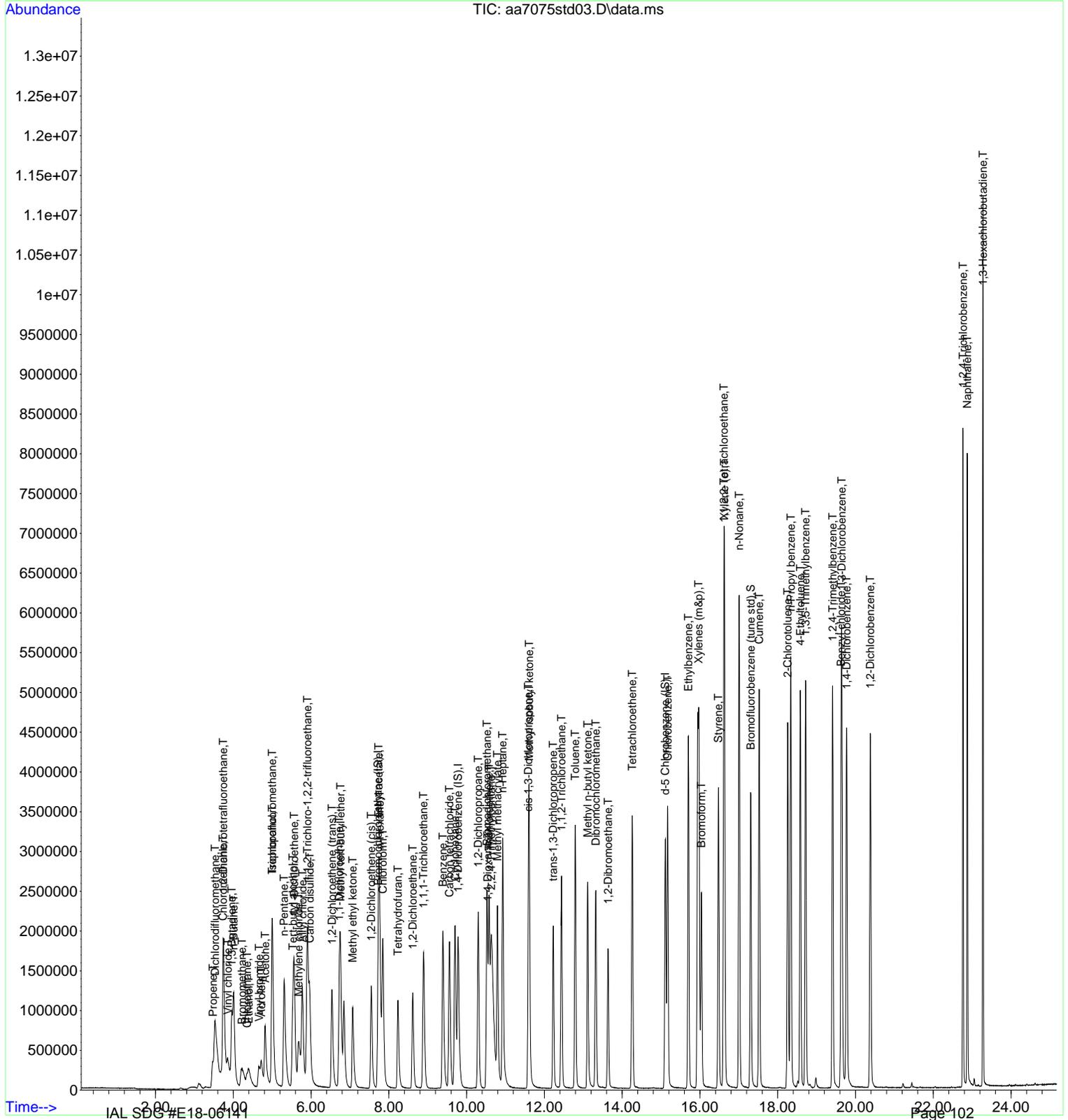
Quant Time: May 18 11:26:02 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 11:25:51 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Dibromochloromethane	13.322	129	1593950	11.00	ppbV	99
53) 1,2-Dibromoethane	13.637	107	1588612	11.00	ppbV	100
54) Tetrachloroethene	14.261	166	1108368	11.00	ppbV	98
56) Chlorobenzene	15.170	112	2154977	11.00	ppbV	100
57) Ethylbenzene	15.701	91	3972028	11.00	ppbV	96
58) Xylenes (m&p)	15.968	91	6199946	22.00	ppbV	96
59) Bromoform	16.039	173	1290866	11.00	ppbV	97
60) Styrene	16.476	104	2164639	11.00	ppbV	98
61) Xylene (o)	16.630	91	3292121	11.00	ppbV	96
62) 1,1,2,2-Tetrachloroethane	16.617	83	2639332	11.00	ppbV	99
63) n-Nonane	17.010	57	2717275	11.00	ppbV #	66
65) Cumene	17.524	105	4095384	11.00	ppbV	97
66) 2-Chlorotoluene	18.257	91	3345139	11.00	ppbV	99
67) n-Propyl benzene	18.337	91	5661498	11.00	ppbV	95
68) 4-Ethyltoluene	18.582	105	4014940	11.00	ppbV	96
69) 1,3,5-Trimethylbenzene	18.717	105	3672316	11.00	ppbV	99
70) 1,2,4-Trimethylbenzene	19.411	105	3660162	11.00	ppbV	98
71) Benzyl chloride	19.630	91	2290255	10.00	ppbV	99
72) 1,3-Dichlorobenzene	19.646	146	2095333	11.00	ppbV	99
73) 1,4-Dichlorobenzene	19.775	146	2149543	11.00	ppbV	99
74) 1,2-Dichlorobenzene	20.382	146	2110390	11.00	ppbV	100
75) 1,2,4-Trichlorobenzene	22.765	180	1792894	12.00	ppbV	99
76) Naphthalene	22.874	127	597227	12.00	ppbV	100
77) 1,3-Hexachlorobutadiene	23.279	225	1181650	11.00	ppbV	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\05-18-18\
 Data File : aa7075std03.D
 Acq On : 18 May 2018 10:53 am
 Operator : jls
 Sample : 10 ppbv Std
 Misc : CC483586
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: May 18 11:26:02 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 11:25:51 2018
 Response via : Initial Calibration



Data Path : C:\DATA\05-18-18\
 Data File : aa7076std04.D
 Acq On : 18 May 2018 12:20 pm
 Operator : jls
 Sample : 2 ppbv Std
 Misc : CC483586
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: May 18 12:46:12 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 12:06:03 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.717	130	531646	10.00	ppbV	0.00	
38) 1,4-Difluorobenzene (IS)	9.775	114	2168184	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.112	117	1841018	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...	17.308	95	1709856	10.10	ppbV	0.00	
Target Compounds							
							Qvalue
2) Propene	3.476	41	132626	1.88	ppbV		83
3) Dichlorodifluoromethane	3.531	85	354292	1.83	ppbV		96
4) Chloromethane	3.724	52	51018	2.00	ppbV		88
5) 1,2-Dichlorotetrafluor...	3.740	85	375176	1.89	ppbV		98
6) Vinyl chloride	3.865	62	145725	1.81	ppbV		84
7) 1,3-Butadiene	3.962	54	129844	1.84	ppbV		93
8) n-Butane	4.007	43	298326	1.97	ppbV		90
9) Bromomethane	4.209	94	97199	1.67	ppbV		96
10) Chloroethane	4.354	64	70288	1.63	ppbV		96
11) Ethanol	4.386	45	91666	2.15	ppbV		97
12) Vinyl bromide	4.647	106	101955	1.68	ppbV		99
13) Acrolein	4.714	56	52814	1.78	ppbV		99
14) Acetone	4.814	58	75164	1.69	ppbV		89
15) Trichlorofluoromethane	4.994	101	411260	1.98	ppbV		100
16) Isopropanol	4.997	45	217705	1.47	ppbV		98
17) n-Pentane	5.306	43	301966	1.76	ppbV		97
18) 1,1-Dichloroethene	5.563	61	246965	1.73	ppbV		91
19) Methylene chloride	5.672	49	218722	1.86	ppbV		100
20) Tert-butyl alcohol	5.531	59	243064	1.81	ppbV		100
21) Allyl chloride	5.762	76	51763	1.62	ppbV		100
22) 1,1,2-Trichloro-1,2,2-...	5.904	101	357853	2.27	ppbV		88
23) Carbon disulfide	5.955	76	380328	1.81	ppbV		96
24) 1,2-Dichloroethene (tr...	6.534	61	212690	1.68	ppbV		98
25) 1,1-Dichloroethane	6.720	63	284656	1.78	ppbV		98
26) Methyl tert-butyl ether	6.756	73	350451	1.84	ppbV		96
27) Methyl ethyl ketone	7.077	43	337043	1.64	ppbV		97
28) 1,2-Dichloroethene (cis)	7.544	61	202527	1.64	ppbV		100
29) Ethyl acetate	7.743	45	54059	1.59	ppbV		100
30) n-Hexane	7.756	57	257212	1.81	ppbV		95
31) Chloroform	7.846	83	310890	1.81	ppbV		99
32) Tetrahydrofuran	8.235	42	162029	1.65	ppbV		88
33) 1,2-Dichloroethane	8.611	62	252565	1.78	ppbV		100
34) 1,1,1-Trichloroethane	8.888	97	333228	2.03	ppbV		100
35) Benzene	9.389	78	435366	1.81	ppbV		100
36) Carbon tetrachloride	9.556	117	321683	1.98	ppbV		99
37) Cyclohexane	9.698	56	284422	1.91	ppbV		90
39) 1,2-Dichloropropane	10.296	63	210794	2.00	ppbV		99
40) Bromodichloromethane	10.521	83	342181	1.90	ppbV		100
41) 2,2,4-Trimethylpentane	10.633	57	763501	1.80	ppbV		94
42) Trichloroethene	10.576	130	161774	1.93	ppbV		97
43) 1,4-Dioxane	10.543	88	75425	1.60	ppbV		95
44) Methyl methacrylate	10.797	41	239222	1.67	ppbV		99
45) n-Heptane	10.929	43	398154	2.12	ppbV		99
46) cis-1,3-Dichloropropene	11.592	75	234782	1.69	ppbV		99
47) Methyl isobutyl ketone	11.617	43	512933	1.96	ppbV		98
48) trans-1,3-Dichloropropene	12.228	75	198542	1.42	ppbV		97
49) 1,1,2-Trichloroethane	12.440	97	202874	2.07	ppbV		89
50) Toluene	12.794	91	532944	2.00	ppbV		96
51) Methyl n-butyl ketone	13.119	43	421816	1.77	ppbV		96

AL SDG #E18-06141

Data Path : C:\DATA\05-18-18\
 Data File : aa7076std04.D
 Acq On : 18 May 2018 12:20 pm
 Operator : jls
 Sample : 2 ppbv Std
 Misc : CC483586
 ALS Vial : 6 Sample Multiplier: 1

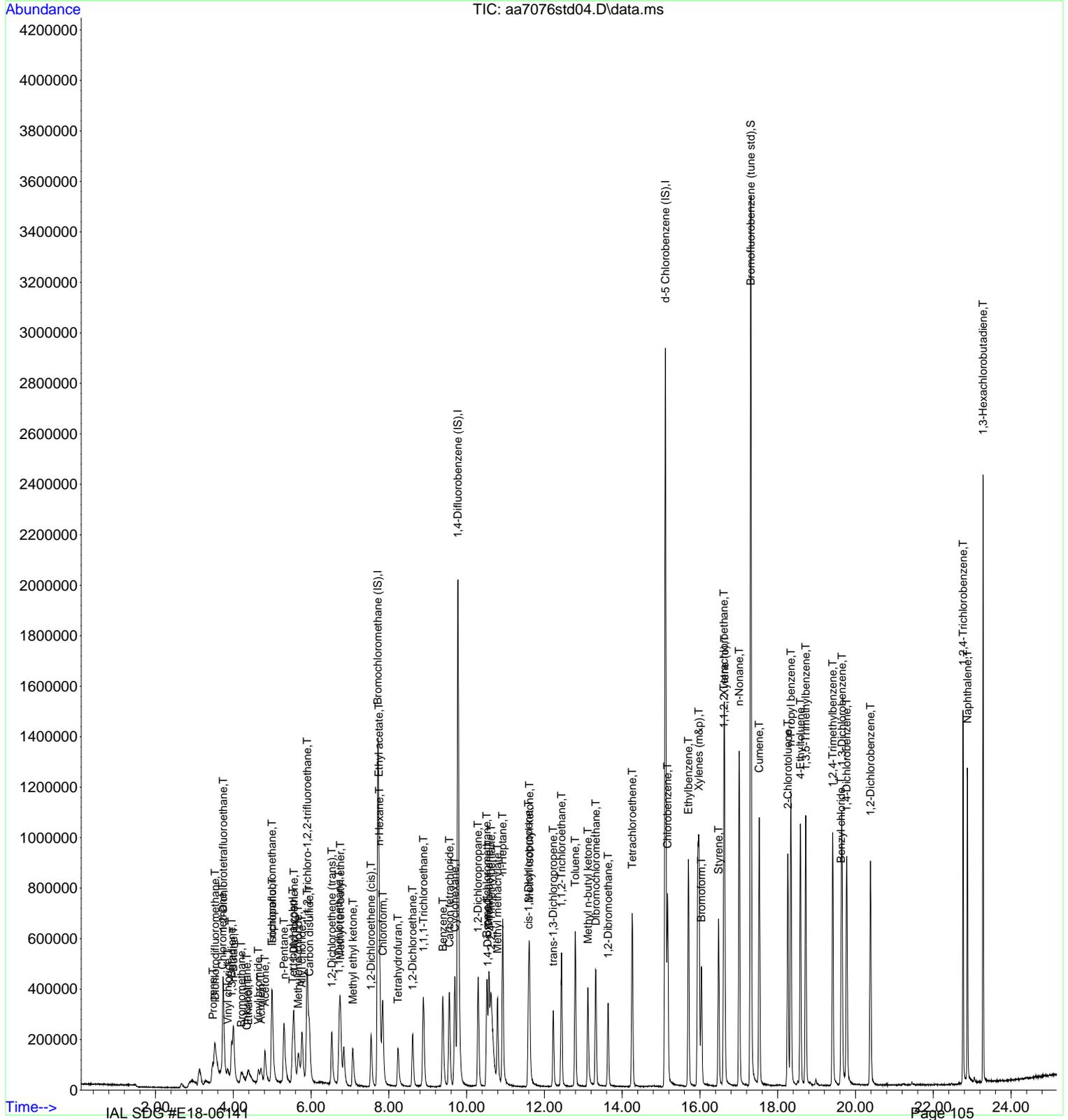
Quant Time: May 18 12:46:12 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 12:06:03 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Dibromochloromethane	13.321	129	301886	2.02	ppbV	98
53) 1,2-Dibromoethane	13.640	107	294478	1.99	ppbV	100
54) Tetrachloroethene	14.260	166	225984	2.19	ppbV	97
56) Chlorobenzene	15.167	112	438067	2.22	ppbV	99
57) Ethylbenzene	15.704	91	774341	2.17	ppbV	96
58) Xylenes (m&p)	15.974	91	1258920	4.44	ppbV	97
59) Bromoform	16.038	173	255272	2.19	ppbV	95
60) Styrene	16.479	104	375106	1.95	ppbV	99
61) Xylene (o)	16.633	91	676541	2.30	ppbV	96
62) 1,1,2,2-Tetrachloroethane	16.614	83	588991	2.46	ppbV	98
63) n-Nonane	17.009	43	683291	2.31	ppbV	98
65) Cumene	17.527	105	830635	2.27	ppbV	97
66) 2-Chlorotoluene	18.260	91	661544	2.19	ppbV	100
67) n-Propyl benzene	18.337	91	1196302	2.37	ppbV	94
68) 4-Ethyltoluene	18.585	105	792642	2.23	ppbV	96
69) 1,3,5-Trimethylbenzene	18.720	105	740731	2.27	ppbV	98
70) 1,2,4-Trimethylbenzene	19.414	105	698841	2.16	ppbV	97
71) Benzyl chloride	19.630	91	304003	1.33	ppbV	100
72) 1,3-Dichlorobenzene	19.649	146	440454	2.34	ppbV	100
73) 1,4-Dichlorobenzene	19.775	146	442950	2.30	ppbV	98
74) 1,2-Dichlorobenzene	20.385	146	428607	2.25	ppbV	99
75) 1,2,4-Trichlorobenzene	22.765	180	317666	2.25	ppbV	99
76) Naphthalene	22.877	127	91044	1.92	ppbV	100
77) 1,3-Hexachlorobutadiene	23.282	225	264100	2.55	ppbV	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\05-18-18\
 Data File : aa7076std04.D
 Acq On : 18 May 2018 12:20 pm
 Operator : jls
 Sample : 2 ppbv Std
 Misc : CC483586
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: May 18 12:46:12 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 12:06:03 2018
 Response via : Initial Calibration



Data Path : C:\DATA\05-18-18\
 Data File : aa7077std05.D
 Acq On : 18 May 2018 1:03 pm
 Operator : jls
 Sample : 0.2 ppbv Std
 Misc : CC483586
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: May 18 13:32:39 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 12:46:18 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.724	130	565151	10.00	ppbV	0.00	
38) 1,4-Difluorobenzene (IS)	9.778	114	2276433	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.109	117	1973260	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...	17.309	95	1806381	9.93	ppbV	0.00	
Target Compounds							
							Qvalue
2) Propene	3.460	41	16894	0.24	ppbV		# 71
3) Dichlorodifluoromethane	3.521	85	50420	0.26	ppbV		# 83
4) Chloromethane	3.740	52	8000m	0.31	ppbV		
5) 1,2-Dichlorotetrafluor...	3.734	85	50893	0.25	ppbV		89
6) Vinyl chloride	3.856	62	19517	0.24	ppbV		# 68
7) 1,3-Butadiene	3.959	54	15122	0.22	ppbV		83
8) n-Butane	4.010	43	44711	0.30	ppbV		97
9) Bromomethane	4.219	94	13685m	0.24	ppbV		
10) Chloroethane	4.344	64	11953m	0.28	ppbV		
11) Ethanol	4.393	45	9615m	0.22	ppbV		
12) Vinyl bromide	4.676	106	11612	0.19	ppbV		94
13) Acrolein	4.724	56	6466	0.22	ppbV		# 15
14) Acetone	4.814	58	11425	0.26	ppbV		81
15) Trichlorofluoromethane	4.994	101	58754	0.28	ppbV		98
16) Isopropanol	4.994	45	29583	0.20	ppbV		# 81
17) n-Pentane	5.303	43	40200	0.24	ppbV		96
18) 1,1-Dichloroethene	5.557	61	34473	0.24	ppbV		90
19) Methylene chloride	5.679	49	41552	0.35	ppbV		98
20) Tert-butyl alcohol	5.531	59	30621	0.23	ppbV		100
21) Allyl chloride	5.762	76	7462	0.24	ppbV		100
22) 1,1,2-Trichloro-1,2,2-...	5.901	101	47494	0.30	ppbV		91
23) Carbon disulfide	5.959	76	51179	0.25	ppbV		# 75
24) 1,2-Dichloroethene (tr...	6.528	61	30920	0.24	ppbV		98
25) 1,1-Dichloroethane	6.717	63	43161	0.27	ppbV		99
26) Methyl tert-butyl ether	6.762	73	40748	0.22	ppbV		98
27) Methyl ethyl ketone	7.084	43	46402	0.23	ppbV		99
28) 1,2-Dichloroethene (cis)	7.550	61	28581	0.23	ppbV		97
29) Ethyl acetate	7.746	45	7027	0.21	ppbV		100
30) n-Hexane	7.756	57	34619	0.24	ppbV		95
31) Chloroform	7.849	83	45270	0.26	ppbV		100
32) Tetrahydrofuran	8.241	42	23222	0.24	ppbV		# 84
33) 1,2-Dichloroethane	8.614	62	39749	0.28	ppbV		97
34) 1,1,1-Trichloroethane	8.881	97	43844	0.27	ppbV		99
35) Benzene	9.392	78	61436	0.26	ppbV		98
36) Carbon tetrachloride	9.556	117	43551	0.27	ppbV		97
37) Cyclohexane	9.698	56	32827	0.22	ppbV		87
39) 1,2-Dichloropropane	10.299	63	29438	0.28	ppbV		97
40) Bromodichloromethane	10.524	83	48412	0.27	ppbV		99
41) 2,2,4-Trimethylpentane	10.624	57	81573	0.19	ppbV		93
42) Trichloroethene	10.579	130	22926	0.28	ppbV		99
43) 1,4-Dioxane	10.550	88	10037	0.22	ppbV		94
44) Methyl methacrylate	10.794	41	27164	0.19	ppbV		94
45) n-Heptane	10.929	43	43536	0.23	ppbV		97
46) cis-1,3-Dichloropropene	11.592	75	30006	0.22	ppbV		98
47) Methyl isobutyl ketone	11.627	43	52478	0.20	ppbV		96
48) trans-1,3-Dichloropropene	12.232	75	24549	0.18	ppbV		91
49) 1,1,2-Trichloroethane	12.437	97	28062	0.29	ppbV		88
50) Toluene	12.797	91	62215	0.23	ppbV		98
51) Methyl n-butyl ketone	13.119	43	42331	0.18	ppbV		93

AL SDG #E18-06141

Data Path : C:\DATA\05-18-18\
 Data File : aa7077std05.D
 Acq On : 18 May 2018 1:03 pm
 Operator : jls
 Sample : 0.2 ppbv Std
 Misc : CC483586
 ALS Vial : 7 Sample Multiplier: 1

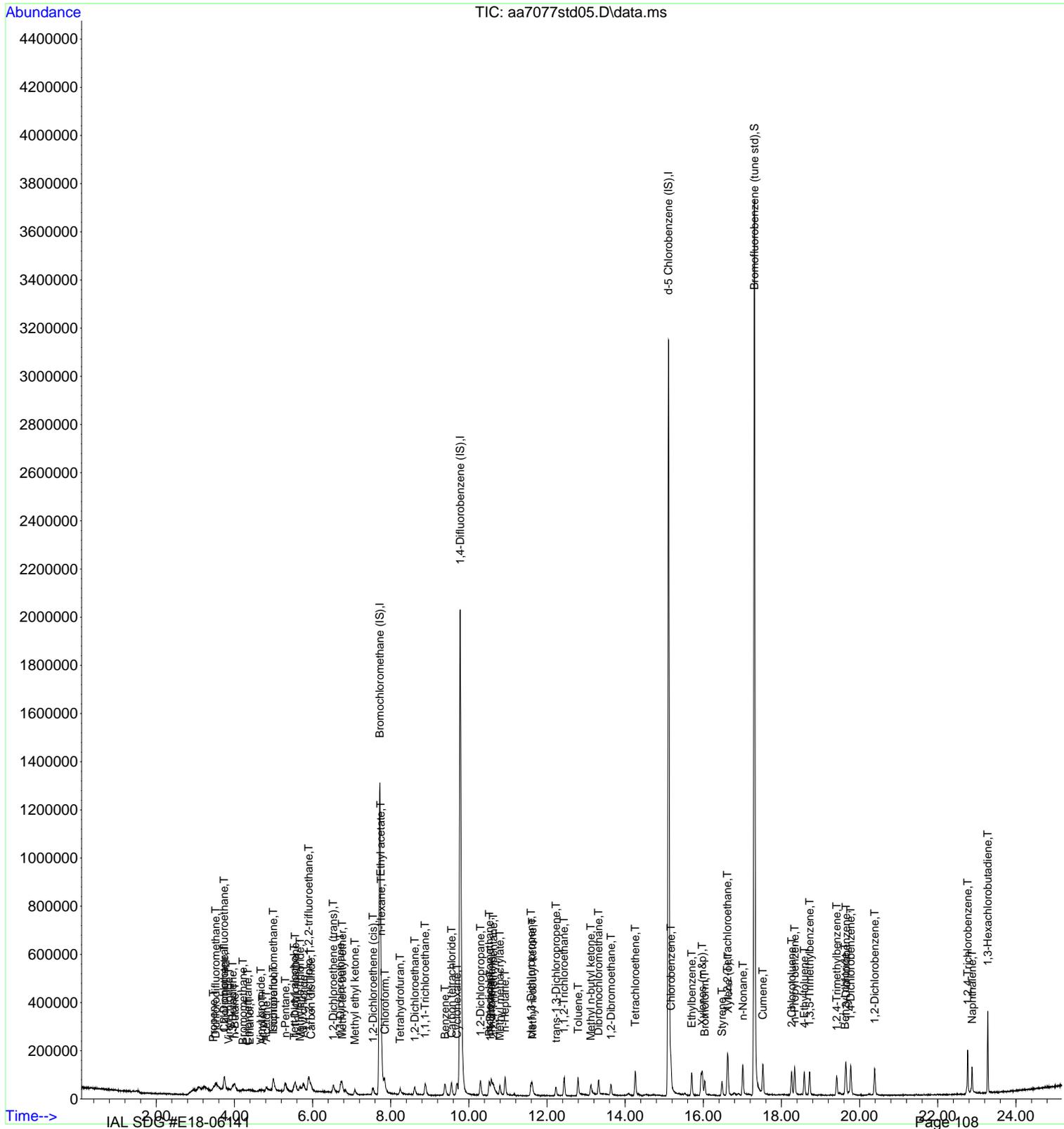
Quant Time: May 18 13:32:39 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 12:46:18 2018
 Response via : Initial Calibration

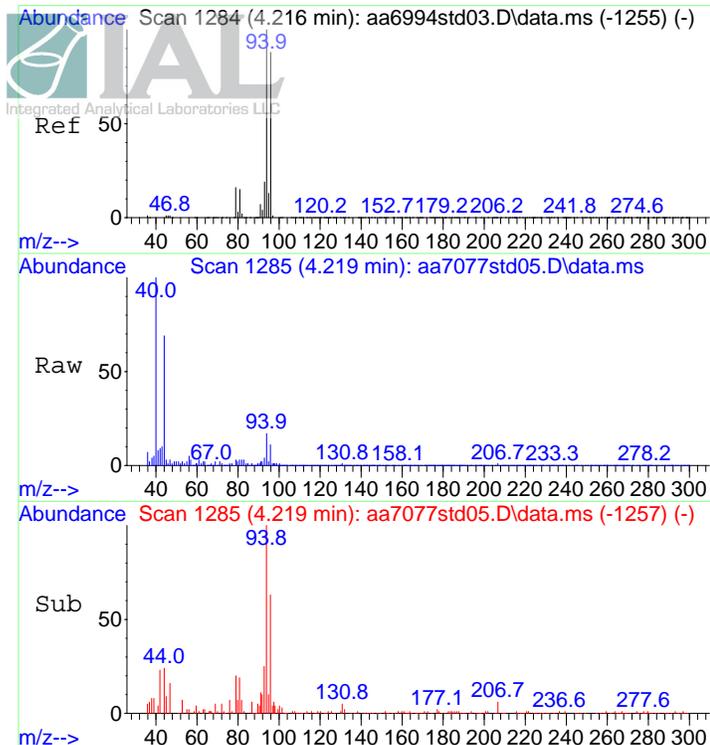
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Dibromochloromethane	13.318	129	40399	0.27	ppbV	96
53) 1,2-Dibromoethane	13.640	107	39419	0.27	ppbV	98
54) Tetrachloroethene	14.260	166	31692	0.31	ppbV	98
56) Chlorobenzene	15.170	112	62317	0.31	ppbV	96
57) Ethylbenzene	15.704	91	84252	0.24	ppbV	96
58) Xylenes (m&p)	15.971	91	126837	0.45	ppbV	97
59) Bromoform	16.039	173	30639	0.26	ppbV	97
60) Styrene	16.479	104	34517	0.18	ppbV	97
61) Xylene (o)	16.633	91	68032	0.23	ppbV	96
62) 1,1,2,2-Tetrachloroethane	16.617	83	73449	0.30	ppbV	97
63) n-Nonane	17.013	43	59250	0.20	ppbV	93
65) Cumene	17.527	105	89239	0.24	ppbV	97
66) 2-Chlorotoluene	18.264	91	71441	0.24	ppbV	99
67) n-Propyl benzene	18.341	91	110526	0.22	ppbV	96
68) 4-Ethyltoluene	18.588	105	73047	0.20	ppbV	95
69) 1,3,5-Trimethylbenzene	18.720	105	66060	0.20	ppbV	98
70) 1,2,4-Trimethylbenzene	19.415	105	59145	0.18	ppbV	99
71) Benzyl chloride	19.630	91	31049	0.13	ppbV	99
72) 1,3-Dichlorobenzene	19.649	146	57201	0.30	ppbV	100
73) 1,4-Dichlorobenzene	19.772	146	54870	0.28	ppbV	100
74) 1,2-Dichlorobenzene	20.386	146	52450	0.27	ppbV	99
75) 1,2,4-Trichlorobenzene	22.765	180	51790	0.37	ppbV	99
76) Naphthalene	22.881	127	10737	0.23	ppbV	100
77) 1,3-Hexachlorobutadiene	23.279	225	38503	0.37	ppbV	93

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\05-18-18\
 Data File : aa7077std05.D
 Acq On : 18 May 2018 1:03 pm
 Operator : jls
 Sample : 0.2 ppbv Std
 Misc : CC483586
 ALS Vial : 7 Sample Multiplier: 1

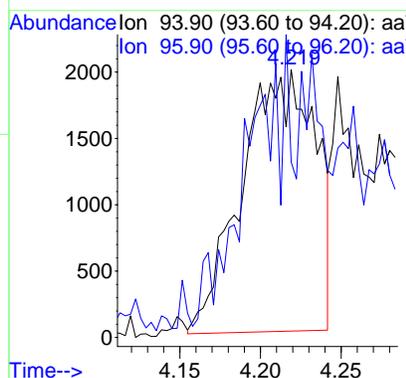
Quant Time: May 18 13:32:39 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 12:46:18 2018
 Response via : Initial Calibration



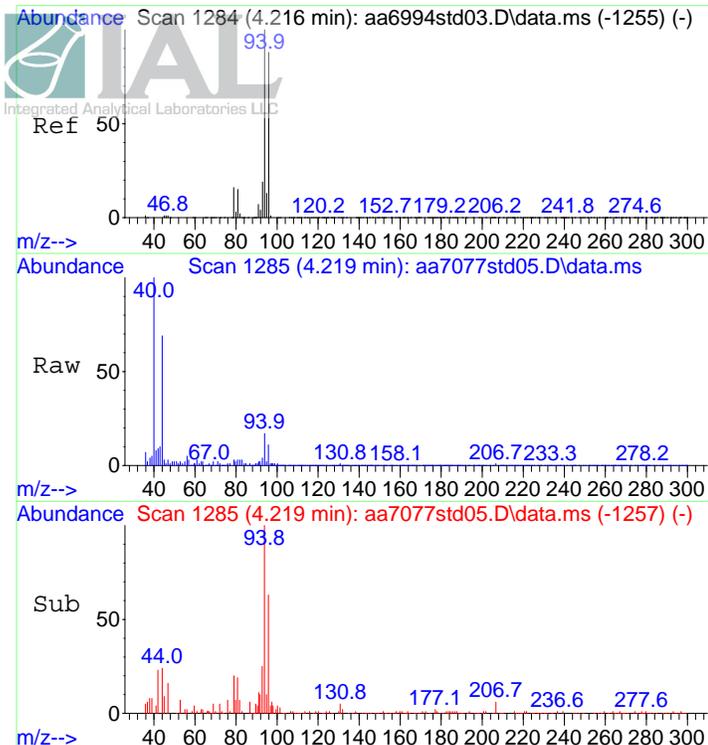


#9
 Bromomethane
 Concen: 0.11 ppbV
 RT: 4.219 min Scan# 1285
 Delta R.T. -0.010 min
 Lab File: aa7077std05.D
 Acq: 18 May 2018 1:03 pm

Tgt Ion	Resp	Lower	Upper
94	6278		
94	100		
96	60.6	75.8	113.8#

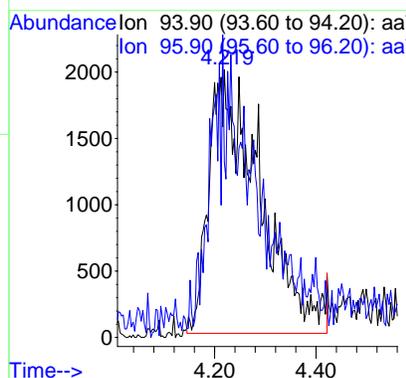


Before manual integration for Bromomethane.

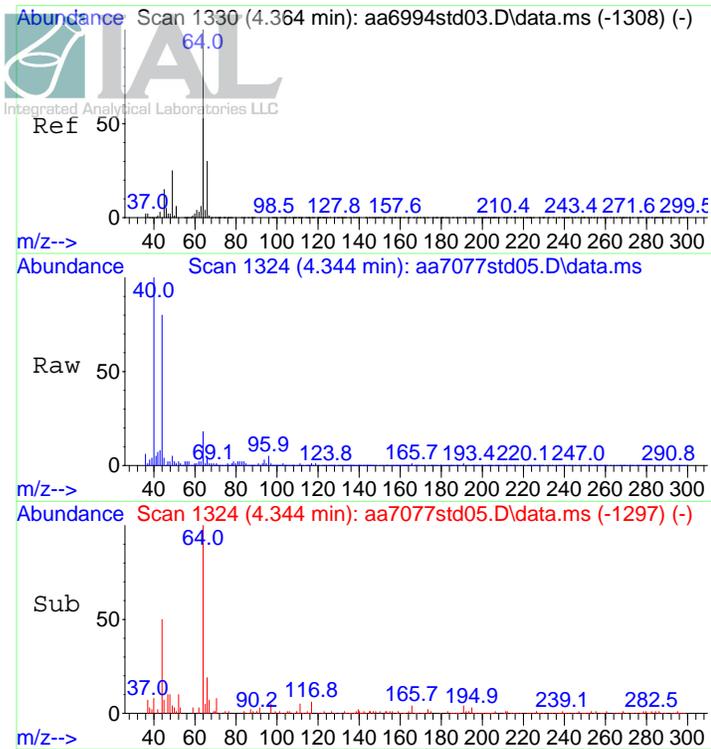


#9
 Bromomethane
 Concen: 0.24 ppbV m
 RT: 4.219 min Scan# 1285
 Delta R.T. -0.010 min
 Lab File: aa7077std05.D
 Acq: 18 May 2018 1:03 pm

Tgt Ion	Resp	Lower	Upper
94	13685		
94	100		
96	27.8	75.8	113.8#

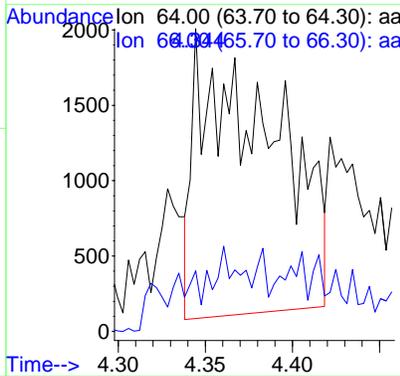


After manual integration for Bromomethane.

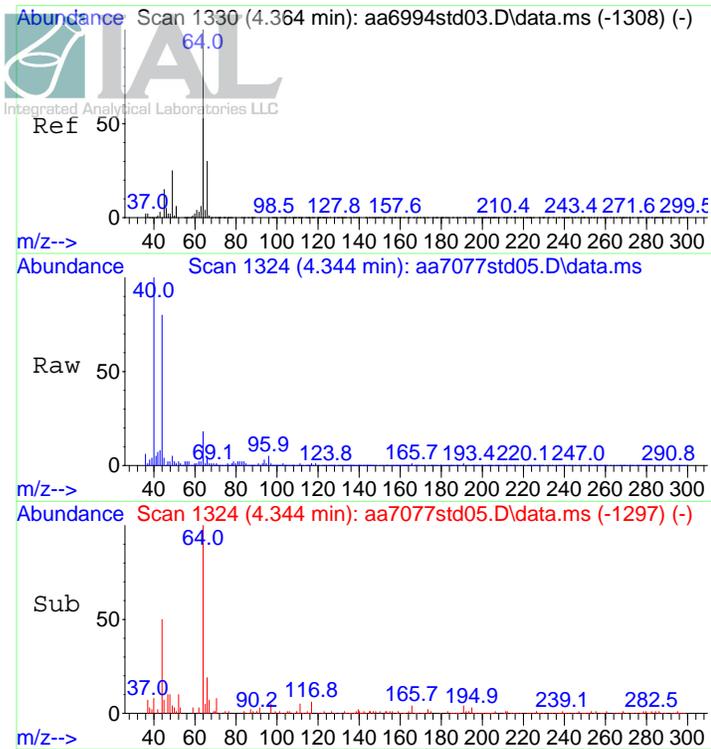


#10
 Chloroethane
 Concen: 0.13 ppbV
 RT: 4.344 min Scan# 1324
 Delta R.T. -0.013 min
 Lab File: aa7077std05.D
 Acq: 18 May 2018 1:03 pm

Tgt Ion: 64 Resp: 5717
 Ion Ratio Lower Upper
 64 100
 66 3.8 21.6 32.4#

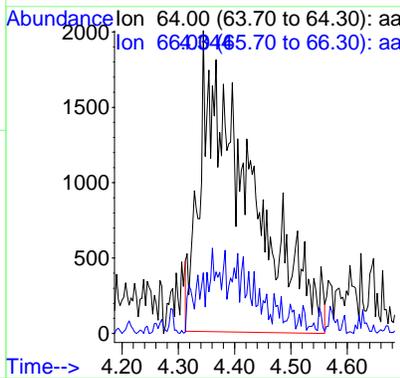


Before manual integration for Chloroethane.

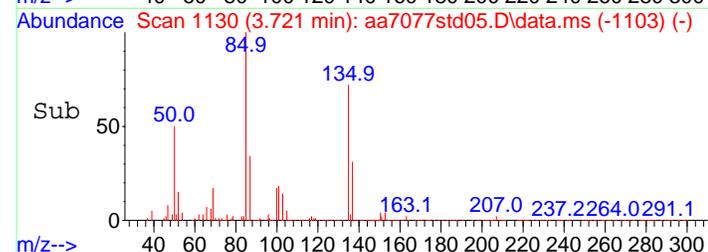
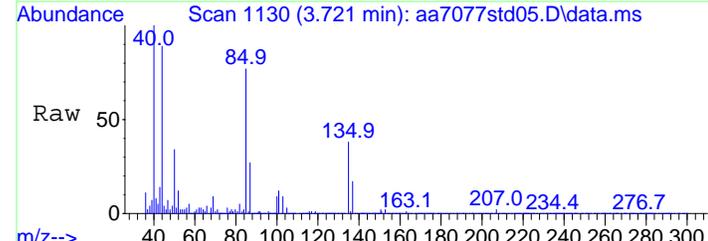
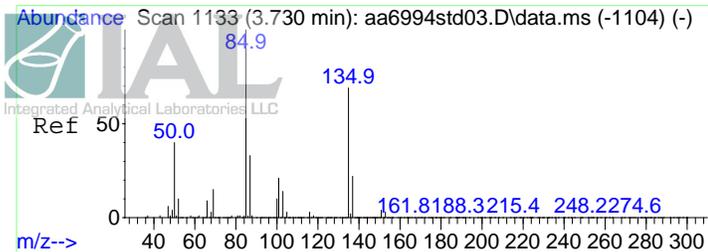


#10
 Chloroethane
 Concen: 0.28 ppbV m
 RT: 4.344 min Scan# 1324
 Delta R.T. -0.013 min
 Lab File: aa7077std05.D
 Acq: 18 May 2018 1:03 pm

Tgt Ion: 64 Resp: 11953
 Ion Ratio Lower Upper
 64 100
 66 1.8 21.6 32.4#

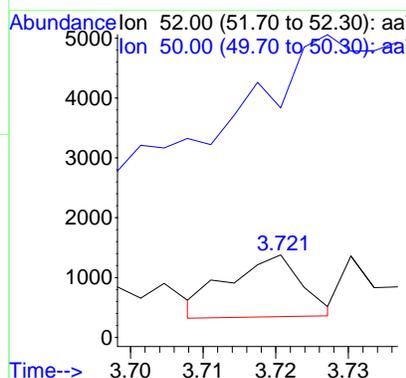


After manual integration for Chloroethane.

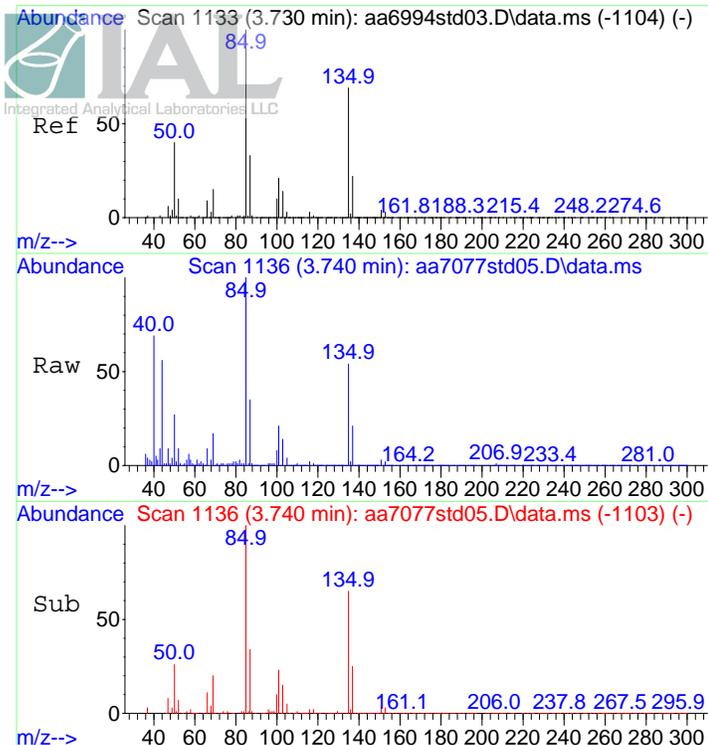


#4
 Chloromethane
 Concen: 0.03 ppbV
 RT: 3.721 min Scan# 1130
 Delta R.T. -0.013 min
 Lab File: aa7077std05.D
 Acq: 18 May 2018 1:03 pm

Tgt Ion	Resp	Lower	Upper
52	727		
52	100		
50	0.0	314.4	471.6#

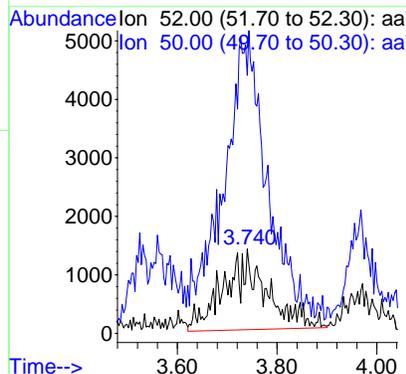


Before manual integration for Chloromethane.

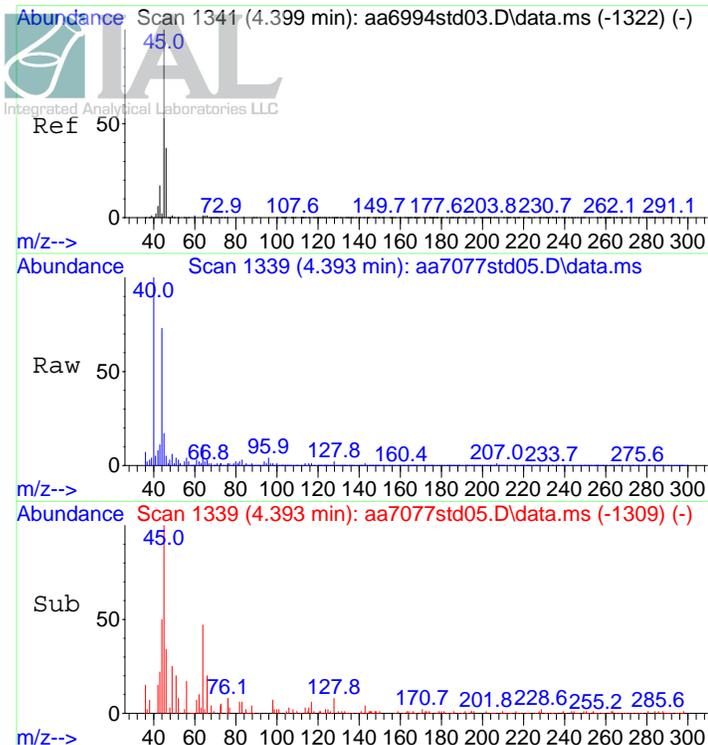


#4
 Chloromethane
 Concen: 0.31 ppbV m
 RT: 3.740 min Scan# 1136
 Delta R.T. 0.006 min
 Lab File: aa7077std05.D
 Acq: 18 May 2018 1:03 pm

Tgt Ion	Resp	Lower	Upper
52	8000		
50	100		
50	0.0	314.4	471.6#

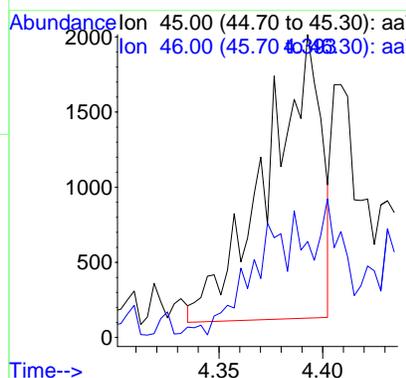


After manual integration for Chloromethane.

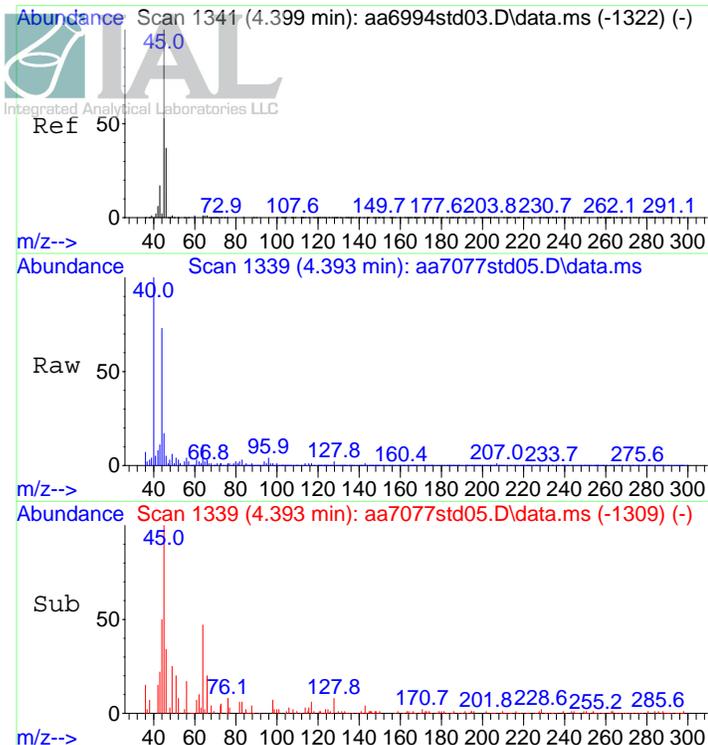


#11
 Ethanol
 Concen: 0.08 ppbV
 RT: 4.393 min Scan# 1339
 Delta R.T. -0.003 min
 Lab File: aa7077std05.D
 Acq: 18 May 2018 1:03 pm

Tgt Ion	Resp	Lower	Upper
45	100		
46	13.7	32.6	48.8#

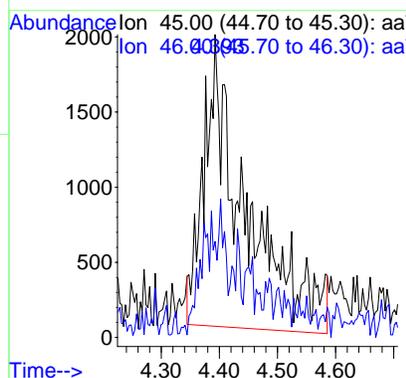


Before manual integration for Ethanol.



#11
 Ethanol
 Concen: 0.22 ppbV m
 RT: 4.393 min Scan# 1339
 Delta R.T. -0.003 min
 Lab File: aa7077std05.D
 Acq: 18 May 2018 1:03 pm

Tgt Ion	Resp	Lower	Upper
45	100		
46	4.9	32.6	48.8#



After manual integration for Ethanol.

Initial Calibration Data Summary Report

Initial Calibration Curve: 7/25/2018
 Instrument: AA

Method ID: 0725.M

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7971BFB]	07/25/2018 08:35
40 PPBV STD [AA7972STD01]	07/25/2018 09:55
20 PPBV STD [AA7973STD02]	07/25/2018 10:29
10 PPBV STD [AA7974STD03]	07/25/2018 11:02
2 PPBV STD [AA7975STD04]	07/25/2018 12:18
0.2 PPBV STD [AA7976STD05]	07/25/2018 13:42
10 PPBV ICVSS [AA7977ICVSS]	07/25/2018 15:05

RParameter	RRF 0.2ppbv	RRF 2ppbv	RRF 10ppbv	RRF 20ppbv	RRF 40ppbv	Avg ppbv	% RSD
Bromochloromethane	-----ISTD-----						
1,4-Difluorobenzene	-----ISTD-----						
d-5 Chlorobenzene	-----ISTD-----						
Acetone	0.73	0.64	0.78	0.82	0.84	0.76	11
Acrolein	0.35	0.44	0.54	0.55	0.58	0.49	19
Allyl Chloride	0.61	0.52	0.67	0.68	0.68	0.63	10
Benzene	4.4	3.9	4.5	4.6	4.8	4.4	7.4
Benzyl chloride	1.3	1.6	1.7	1.7	1.6	1.6	12
Bromodichloromethane	0.76	0.68	0.76	0.76	0.75	0.74	4.5
Bromoform	0.62	0.62	0.62	0.60	0.56	0.60	4.4
Bromomethane	0.96	0.91	1.1	1.1	1.1	1.0	7.8
1,3-Butadiene	1.1	1.0	1.2	1.2	1.2	1.2	7.5
Chlorobenzene	1.3	1.1	1.1	1.0	0.98	1.1	9.9
Chloroethane	0.69	0.63	0.77	0.80	0.84	0.75	12
Chloroform	3.0	2.5	3.0	3.1	3.2	3.0	8.4
Chloromethane	0.44	0.38	0.41	0.41	0.42	0.41	5.4
Carbon disulfide	3.7	3.4	4.1	4.2	4.3	4.0	9.2
Carbon tetrachloride	3.0	2.6	2.7	2.7	2.9	2.8	4.9
2-Chlorotoluene	1.8	1.7	1.6	1.6	1.5	1.6	8.3
Cumene	2.2	2.1	2.0	1.9	1.7	2.0	8.7
Cyclohexane	2.5	2.4	2.5	2.5	2.7	2.5	3.6
Dibromochloromethane	0.64	0.59	0.62	0.62	0.62	0.62	2.6
1,2-Dibromoethane	0.75	0.63	0.66	0.67	0.67	0.67	6.3
1,2-Dichlorobenzene	1.2	1.0	0.99	0.95	0.87	1.0	11
1,3-Dichlorobenzene	1.2	1.1	0.98	0.92	0.82	0.99	14
1,4-Dichlorobenzene	1.2	1.1	1.0	0.97	0.88	1.0	10
Dichlorodifluoromethane	3.4	2.7	3.2	3.3	3.5	3.2	9.6
1,1-Dichloroethane	2.8	2.4	2.8	2.9	3.0	2.8	7.9
1,2-Dichloroethane	2.3	1.9	2.3	2.3	2.5	2.3	8.4
1,1-Dichloroethene	2.3	2.0	2.4	2.5	2.5	2.3	8.6
1,2-Dichloroethene (trans)	2.0	1.8	2.2	2.3	2.4	2.1	10
1,2-Dichloroethene (trans)	1.9	1.8	2.1	2.2	2.3	2.1	11
1,2-Dichloropropane	0.52	0.42	0.45	0.45	0.45	0.46	8.1
1,3-Dichloropropene (cis)	0.62	0.56	0.64	0.66	0.65	0.62	6.2
1,3-Dichloropropene (trans)	0.63	0.57	0.66	0.68	0.69	0.65	7.8
1,2-Dichlorotetrafluoroethane	4.1	3.3	3.5	3.4	3.3	3.5	9.5
1,4-Dioxane	0.18	0.20	0.22	0.22	0.22	0.21	7.7
Ethanol	0.84	1.0	0.71	0.69	0.68	0.79	19
Ethylbenzene	2.2	2.1	1.9	1.9	1.7	2.0	9.7
4-Ethyltoluene	2.0	2.0	1.9	1.9	1.7	1.9	6.7
n-Heptane	0.69	0.65	0.62	0.62	0.62	0.64	4.4
1,3-Hexachlorobutadiene	0.76	0.61	0.53	0.46	0.42	0.55	24

*% RSD (Relative Standard Deviation) must be within 30%

**An exception is made for 2 compounds that must be within 40%

RRF - Relative Response Factor

Initial Calibration Data Summary Report

Initial Calibration Curve: 7/25/2018
Instrument: AA

Method ID: 0725.M

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7971BFB]	07/25/2018 08:35
40 PPBV STD [AA7972STD01]	07/25/2018 09:55
20 PPBV STD [AA7973STD02]	07/25/2018 10:29
10 PPBV STD [AA7974STD03]	07/25/2018 11:02
2 PPBV STD [AA7975STD04]	07/25/2018 12:18
0.2 PPBV STD [AA7976STD05]	07/25/2018 13:42
10 PPBV ICVSS [AA7977ICVSS]	07/25/2018 15:05

RParameter	RRF 0.2ppbv	RRF 2ppbv	RRF 10ppbv	RRF 20ppbv	RRF 40ppbv	Avg ppbv	% RSD
n-Hexane	2.5	2.2	2.4	2.5	2.6	2.4	6.3
Isopropanol	2.4	2.3	2.5	2.6	2.7	2.5	6.6
Methylene chloride	2.6	1.6	1.8	1.8	1.9	1.9	21
Methyl ethyl ketone	2.9	2.3	2.8	2.9	3.1	2.8	11
Methyl isobutyl ketone	0.91	0.85	0.83	0.83	0.82	0.85	4.6
Methyl methacrylate	0.52	0.43	0.47	0.48	0.48	0.48	6.3
Methyl n-butyl ketone	0.83	0.74	0.79	0.80	0.81	0.80	4.1
Methyl tert-butyl ether	4.1	3.8	4.0	4.1	4.2	4.0	3.3
Naphthalene	0.25	0.24	0.27	0.26	0.23	0.25	5.6
Propene	0.66	0.81	0.91	0.92	0.97	0.85	15
Styrene	1.0	1.1	1.1	1.1	1.0	1.1	2.4
Tert-butyl alcohol	3.3	2.8	3.2	3.3	3.4	3.2	7.1
1,1,2,2-Tetrachloroethane	1.5	1.3	1.2	1.2	1.1	1.2	13
Tetrachloroethene	0.55	0.46	0.44	0.43	0.42	0.46	11
Tetrahydrofuran	1.4	1.2	1.4	1.5	1.6	1.4	11
Toluene	1.4	1.2	1.2	1.2	1.2	1.3	5.2
1,2,4-Trichlorobenzene	0.95	0.77	0.75	0.70	0.62	0.76	16
1,1,1-Trichloroethane	3.2	2.8	2.9	2.9	3.0	3.0	4.9
1,1,2-Trichloroethane	0.49	0.42	0.43	0.43	0.43	0.44	6.3
Trichloroethene	0.42	0.33	0.37	0.36	0.35	0.36	9.3
Trichlorofluoromethane	3.7	3.1	3.4	3.4	3.5	3.4	6.2
1,1,2-Trichloro-1,2,2-trifluoroethane	3.2	2.8	2.7	2.6	2.7	2.8	8.1
1,2,4-Trimethylbenzene	1.7	1.8	1.8	1.7	1.6	1.7	5.7
1,3,5-Trimethylbenzene	1.9	1.8	1.8	1.7	1.5	1.7	8.2
2,2,4-Trimethylpentane	1.8	1.7	1.9	1.9	1.9	1.8	3.9
Vinyl bromide	0.96	1.0	1.2	1.3	1.3	1.2	14
Vinyl chloride	1.4	1.2	1.5	1.5	1.6	1.4	9.1
Xylenes (m&p)	1.7	1.6	1.6	1.4	1.3	1.5	10
Xylenes (o)	1.7	1.7	1.5	1.5	1.3	1.6	10

*% RSD (Relative Standard Deviation) must be within 30%

**An exception is made for 2 compounds that must be within 40%

RRF - Relative Response Factor

Method Path : C:\msdchem\1\METHODS\
 Method File : 0725.M
 Title : TO-15 on the Agilent 7890A / 5975C
 Last Update : Wed Jul 25 14:15:57 2018
 Response Via : Initial Calibration

Calibration Files

0.2 =aa7976std05.D 2 =aa7975std04.D 10 =aa7974std03.D 20 =aa7973std02.D 40 =aa7972std01.D

Compound	0.2	2	10	20	40	Avg	%RSD
1) I Bromochloromethane...	----- ISTD -----						
2) T Propene	0.658	0.808	0.908	0.923	0.972	0.854	14.62
3) T Dichlorodifluoro...	3.407	2.691	3.222	3.317	3.462	3.220	9.61
4) T Chloromethane	0.440	0.377	0.411	0.411	0.419	0.411	5.46
5) T 1,2-Dichlorotetr...	4.099	3.300	3.466	3.384	3.304	3.511	9.58
6) T Vinyl chloride	1.396	1.218	1.476	1.496	1.557	1.428	9.18
7) T 1,3-Butadiene	1.088	1.036	1.201	1.209	1.234	1.153	7.50
8) T n-Butane	2.520	2.013	2.174	2.181	2.231	2.224	8.31
9) T Bromomethane	0.960	0.905	1.085	1.073	1.058	1.016	7.82
10) T Chloroethane	0.689	0.625	0.772	0.801	0.842	0.746	11.76
11) T Ethanol	0.835	1.029	0.705	0.685	0.683	0.787	18.92
12) T Vinyl bromide	0.964	1.009	1.229	1.257	1.315	1.155	13.63
13) T Acrolein	0.348	0.443	0.537	0.552	0.582	0.493	19.44
14) T Acetone	0.732	0.639	0.779	0.815	0.842	0.762	10.50
15) T Trichlorofluorom...	3.732	3.142	3.400	3.411	3.519	3.441	6.22
16) T Isopropanol	2.370	2.258	2.541	2.592	2.655	2.483	6.63
17) T n-Pentane	2.781	2.129	2.470	2.523	2.629	2.506	9.66
18) T 1,1-Dichloroethene	2.280	2.017	2.424	2.454	2.529	2.341	8.64
19) T Methylene chloride	2.621	1.573	1.757	1.778	1.854	1.917	21.24
20) T Tert-butyl alcohol	3.268	2.827	3.151	3.263	3.444	3.191	7.17
21) T Allyl chloride	0.610	0.524	0.666	0.679	0.675	0.631	10.40
22) T 1,1,2-Trichloro-...	3.181	2.816	2.655	2.636	2.678	2.794	8.17
23) T Carbon disulfide	3.749	3.412	4.135	4.175	4.293	3.953	9.22
24) T 1,2-Dichloroethe...	2.027	1.800	2.211	2.263	2.364	2.133	10.43
25) T 1,1-Dichloroethane	2.792	2.391	2.810	2.865	2.961	2.764	7.91
26) T Methyl tert-butyl...	4.058	3.814	4.005	4.060	4.179	4.023	3.30
27) T Methyl ethyl ketone	2.891	2.334	2.803	2.942	3.134	2.821	10.55
28) T 1,2-Dichloroethe...	1.920	1.768	2.144	2.200	2.298	2.066	10.50
29) T Ethyl acetate	0.456	0.394	0.472	0.498	0.527	0.469	10.63
30) T n-Hexane	2.459	2.176	2.401	2.474	2.596	2.421	6.39
31) T Chloroform	3.035	2.533	3.036	3.065	3.174	2.969	8.43
32) T Tetrahydrofuran	1.411	1.174	1.445	1.506	1.605	1.428	11.20
33) T 1,2-Dichloroethane	2.326	1.949	2.281	2.338	2.463	2.271	8.48
34) T 1,1,1-Trichloroe...	3.160	2.787	2.883	2.916	3.046	2.959	4.93
35) T Benzene	4.366	3.921	4.452	4.604	4.801	4.429	7.41
36) T Carbon tetrachlo...	2.964	2.607	2.719	2.744	2.866	2.780	4.97
37) T Cyclohexane	2.519	2.395	2.517	2.531	2.652	2.523	3.61
38) I 1,4-Difluorobenzen...	----- ISTD -----						
39) T 1,2-Dichloropropane	0.518	0.419	0.445	0.450	0.448	0.456	8.10
40) T Bromodichloromet...	0.762	0.681	0.757	0.757	0.745	0.740	4.55
41) T 2,2,4-Trimethylp...	1.841	1.721	1.903	1.886	1.858	1.841	3.90
42) T Trichloroethene	0.420	0.329	0.366	0.356	0.347	0.364	9.37
43) T 1,4-Dioxane	0.184	0.196	0.216	0.222	0.217	0.207	7.79
44) T Methyl methacrylate	0.518	0.434	0.468	0.476	0.484	0.476	6.33
45) T n-Heptane	0.686	0.647	0.623	0.620	0.621	0.639	4.43
46) T cis-1,3-Dichloro...	0.619	0.558	0.640	0.655	0.648	0.624	6.28
47) T Methyl isobutyl ...	0.913	0.849	0.833	0.826	0.815	0.847	4.60
48) T trans-1,3-Dichlo...	0.625	0.567	0.660	0.683	0.691	0.645	7.82
49) T 1,1,2-Trichloroe...	0.488	0.422	0.427	0.430	0.426	0.438	6.31
50) T Toluene	1.372	1.209	1.235	1.235	1.223	1.255	5.29
51) T Methyl n-butyl k...	0.827	0.743	0.788	0.804	0.814	0.795	4.10
52) T Dibromochloromet...	0.637	0.593	0.618	0.624	0.615	0.618	2.60
53) T 1,2-Dibromoethane	0.745	0.632	0.656	0.667	0.665	0.673	6.36
54) T Tetrachloroethene	0.548	0.455	0.439	0.433	0.417	0.459	11.32
55) I d-5 Chlorobenzene ...	----- ISTD -----						
56) T Chlorobenzene	1.267	1.089	1.063	1.039	0.982	1.088	9.90
57) T Ethylbenzene	2.243	2.071	1.939	1.866	1.742	1.972	9.77
58) T Xylenes (m&p)	1.712	1.608	1.590	1.424	1.316	1.530	10.33
59) T Bromoform	0.624	0.616	0.621	0.601	0.558	0.604	4.46
60) T Styrene	1.043	1.063	1.094	1.081	1.030	1.062	2.49

ALSDG #E18-06141

Method Path : C:\msdchem\1\METHODS\
Method File : 0725.M

61)	T	Xylene (o)	1.742	1.679	1.546	1.462	1.340	1.554	10.45
62)	T	1,1,2,2-Tetrachl...	1.460	1.346	1.212	1.153	1.053	1.245	12.90
63)	T	n-Nonane	1.295	1.298	1.166	1.100	1.023	1.176	10.27
64)	S	Bromofluorobenze...	0.921	0.917	0.909	0.906	0.906	0.912	0.73
65)	T	Cumene	2.160	2.099	1.974	1.876	1.730	1.968	8.78
66)	T	2-Chlorotoluene	1.818	1.696	1.614	1.552	1.463	1.629	8.35
67)	T	n-Propyl benzene	2.859	2.929	2.711	2.547	2.282	2.666	9.75
68)	T	4-Ethyltoluene	1.958	2.029	1.933	1.850	1.695	1.893	6.75
69)	T	1,3,5-Trimethylb...	1.881	1.846	1.760	1.665	1.526	1.736	8.29
70)	T	1,2,4-Trimethylb...	1.722	1.833	1.764	1.692	1.567	1.715	5.75
71)	T	Benzyl chloride	1.262	1.566	1.730	1.737	1.626	1.584	12.26
72)	T	1,3-Dichlorobenzene	1.176	1.056	0.975	0.917	0.815	0.987	13.88
73)	T	1,4-Dichlorobenzene	1.155	1.076	1.013	0.970	0.884	1.020	10.12
74)	T	1,2-Dichlorobenzene	1.169	1.048	0.986	0.945	0.870	1.004	11.26
75)	T	1,2,4-Trichlorob...	0.953	0.769	0.754	0.703	0.616	0.759	16.32
76)	T	Naphthalene	0.252	0.238	0.265	0.264	0.234	0.251	5.61
77)	T	1,3-Hexachlorobu...	0.758	0.608	0.526	0.461	0.416	0.554	24.44

(#) = Out of Range

Data Path : C:\DATA\07-25-18\
 Data File : aa7972std01.D
 Acq On : 25 Jul 2018 9:55 am
 Operator : jls
 Sample : 40 ppbv Std
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Jul 25 11:49:13 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 11:48:47 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.702	130	459428	10.00	ppbV	0.01	
38) 1,4-Difluorobenzene (IS)	9.743	114	2108198	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.058	117	1916771	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...	17.251	95	1737147	9.98	ppbV	0.00	
Target Compounds							
							Qvalue
2) Propene	3.477	41	1965032	46.71	ppbV	#	79
3) Dichlorodifluoromethane	3.531	85	6999220	46.59	ppbV		97
4) Chloromethane	3.734	52	846080	44.82	ppbV	#	64
5) 1,2-Dichlorotetrafluor...	3.737	85	6071502	38.59	ppbV		90
6) Vinyl chloride	3.859	62	3148204	46.13	ppbV		87
7) 1,3-Butadiene	3.956	54	2494533	45.07	ppbV	#	76
8) n-Butane	4.001	43	4509713	45.08	ppbV		98
9) Bromomethane	4.203	94	1924371	38.82	ppbV		100
10) Chloroethane	4.342	64	1546620	42.80	ppbV		92
11) Ethanol	4.377	45	1180335	36.98	ppbV		97
12) Vinyl bromide	4.644	106	2658643	46.55	ppbV		97
13) Acrolein	4.708	56	1283660	51.33	ppbV		98
14) Acetone	4.808	58	1702732	46.49	ppbV		71
15) Trichlorofluoromethane	4.985	101	7112735	45.46	ppbV		99
16) Isopropanol	4.985	45	4879981	41.38	ppbV		93
17) n-Pentane	5.287	43	5313474	46.33	ppbV		92
18) 1,1-Dichloroethene	5.544	61	5112666	45.63	ppbV		99
19) Methylene chloride	5.660	49	3746902	46.14	ppbV	#	78
20) Tert-butyl alcohol	5.509	59	7595366	51.55	ppbV		100
21) Allyl chloride	5.750	76	1363505	44.15	ppbV		100
22) 1,1,2-Trichloro-1,2,2-...	5.882	101	5414219	44.54	ppbV		88
23) Carbon disulfide	5.946	76	8677369	45.46	ppbV	#	90
24) 1,2-Dichloroethene (tr...	6.512	61	4778182	46.49	ppbV		90
25) 1,1-Dichloroethane	6.699	63	5986100	45.92	ppbV		99
26) Methyl tert-butyl ether	6.728	73	8447001	45.60	ppbV		100
27) Methyl ethyl ketone	7.039	43	6334927	48.00	ppbV	#	83
28) 1,2-Dichloroethene (cis)	7.518	61	4645622	46.56	ppbV		89
29) Ethyl acetate	7.708	45	1065569	47.82	ppbV		100
30) n-Hexane	7.737	57	5248464	46.87	ppbV	#	80
31) Chloroform	7.814	83	6415960	45.78	ppbV		99
32) Tetrahydrofuran	8.197	42	3538735	52.19	ppbV		95
33) 1,2-Dichloroethane	8.580	62	4979239	46.93	ppbV		99
34) 1,1,1-Trichloroethane	8.856	97	6158415	46.23	ppbV		100
35) Benzene	9.354	78	9704796	46.65	ppbV	#	89
36) Carbon tetrachloride	9.518	117	5792951	46.17	ppbV		100
37) Cyclohexane	9.663	56	5361381	46.24	ppbV		98
39) 1,2-Dichloropropane	10.258	63	4158869	44.11	ppbV		99
40) Bromodichloromethane	10.483	83	6906930	43.29	ppbV		98
41) 2,2,4-Trimethylpentane	10.592	57	17233646	43.16	ppbV		99
42) Trichloroethene	10.538	130	3222063	42.33	ppbV		99
43) 1,4-Dioxane	10.496	88	1833462	39.70	ppbV		82
44) Methyl methacrylate	10.753	41	4490218	45.15	ppbV	#	70
45) n-Heptane	10.888	43	5759111	43.94	ppbV	#	76
46) cis-1,3-Dichloropropene	11.544	75	6011272	44.02	ppbV		99
47) Methyl isobutyl ketone	11.567	43	7556480	43.21	ppbV	#	83
48) trans-1,3-Dichloropropene	12.181	75	5825154	41.15	ppbV	#	77
49) 1,1,2-Trichloroethane	12.393	97	3947997	43.72	ppbV		92
50) Toluene	12.747	91	11346790	43.58	ppbV		99
51) Methyl n-butyl ketone	13.062	43	7546602	44.97	ppbV	#	80

AL SDG #E18-06141

Data Path : C:\DATA\07-25-18\
 Data File : aa7972std01.D
 Acq On : 25 Jul 2018 9:55 am
 Operator : jls
 Sample : 40 ppbv Std
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

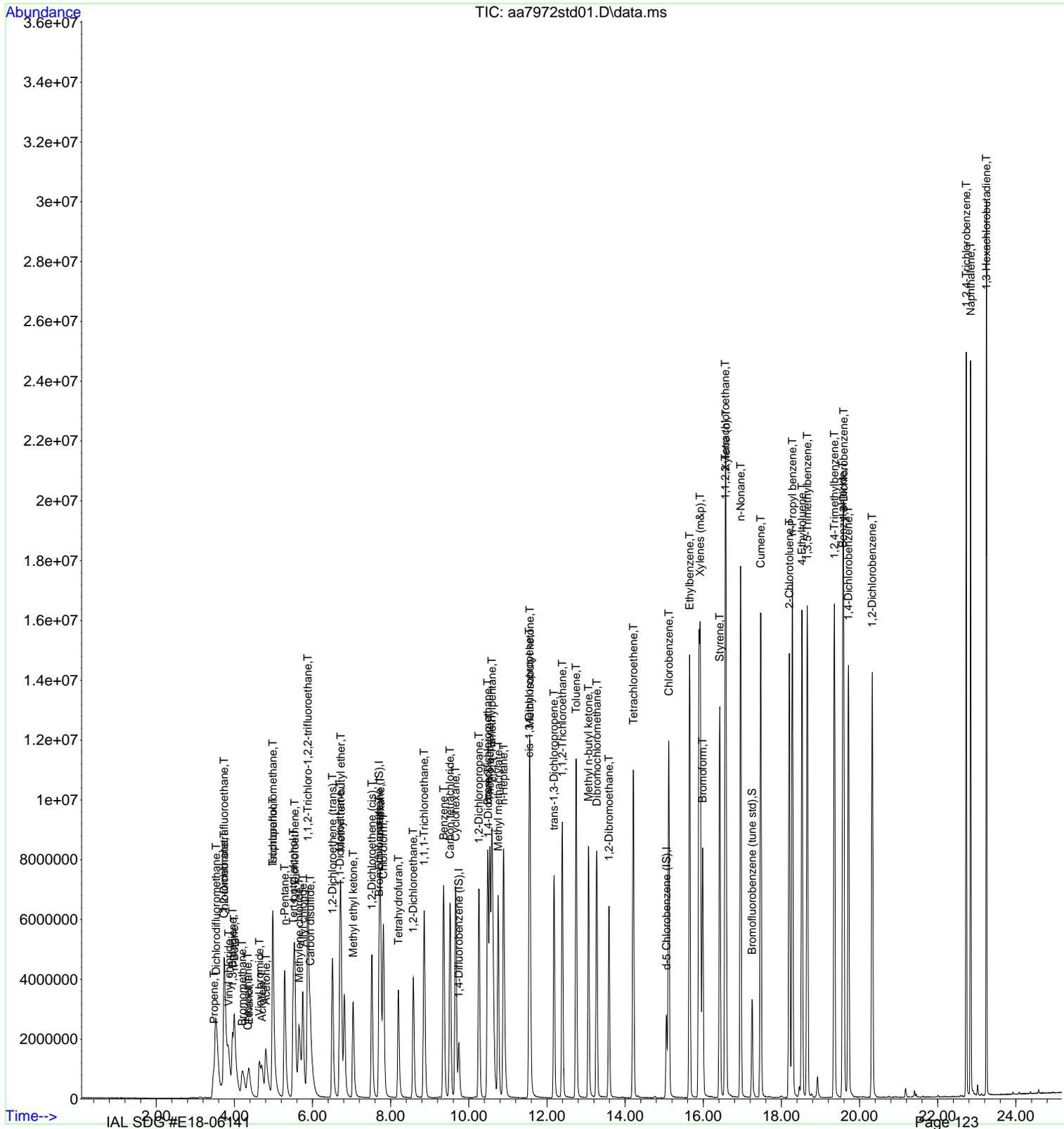
Quant Time: Jul 25 11:49:13 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 11:48:47 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Dibromochloromethane	13.274	129	5704164	43.57	ppbV	100
53) 1,2-Dibromoethane	13.586	107	6166796	44.24	ppbV	96
54) Tetrachloroethene	14.210	166	3866986	42.04	ppbV	98
56) Chlorobenzene	15.116	112	8282237	41.11	ppbV	95
57) Ethylbenzene	15.650	91	14690673	40.29	ppbV	100
58) Xylenes (m&p)	15.917	91	22193499	76.81	ppbV	99
59) Bromoform	15.984	173	4710263	40.23	ppbV	98
60) Styrene	16.422	104	8683428	41.65	ppbV	97
61) Xylene (o)	16.579	91	11299455	39.20	ppbV	100
62) 1,1,2,2-Tetrachloroethane	16.560	83	8878843	39.18	ppbV	98
63) n-Nonane	16.955	43	8627836	39.73	ppbV #	75
65) Cumene	17.470	105	14588254	39.54	ppbV	98
66) 2-Chlorotoluene	18.203	91	12340907	40.68	ppbV	97
67) n-Propyl benzene	18.280	91	19245579	38.19	ppbV	98
68) 4-Ethyltoluene	18.528	105	14294954	39.43	ppbV	98
69) 1,3,5-Trimethylbenzene	18.663	105	12872347	39.21	ppbV	99
70) 1,2,4-Trimethylbenzene	19.354	105	13212625	39.90	ppbV	100
71) Benzyl chloride	19.569	91	12469556	37.52	ppbV	97
72) 1,3-Dichlorobenzene	19.592	146	6869883	37.90	ppbV	99
73) 1,4-Dichlorobenzene	19.714	146	7454051	39.21	ppbV	98
74) 1,2-Dichlorobenzene	20.325	146	7337521	39.65	ppbV	99
75) 1,2,4-Trichlorobenzene	22.730	180	5665233	40.58	ppbV	100
76) Naphthalene	22.843	127	2155413	42.56	ppbV	100
77) 1,3-Hexachlorobutadiene	23.248	225	3506387	37.07	ppbV	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\07-25-18\
 Data File : aa7972std01.D
 Acq On : 25 Jul 2018 9:55 am
 Operator : jls
 Sample : 40 ppbv Std
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Jul 25 11:49:13 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 11:48:47 2018
 Response via : Initial Calibration



Data Path : C:\DATA\07-25-18\
 Data File : aa7973std02.D
 Acq On : 25 Jul 2018 10:29 am
 Operator : jls
 Sample : 20 ppbv Std
 Misc : CC483586
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Jul 25 11:48:30 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 11:48:15 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.695	130	486694	10.00	ppbV	0.00	
38) 1,4-Difluorobenzene (IS)	9.744	114	2128550	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.058	117	1865915	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...	17.251	95	1691328	9.97	ppbV	0.00	
Target Compounds							
							Qvalue
2) Propene	3.470	41	988321	22.35	ppbV	#	79
3) Dichlorodifluoromethane	3.522	85	3551951	22.65	ppbV		97
4) Chloromethane	3.715	52	439853	21.99	ppbV		87
5) 1,2-Dichlorotetrafluor...	3.731	85	3294031	19.53	ppbV		93
6) Vinyl chloride	3.850	62	1601376	22.30	ppbV		87
7) 1,3-Butadiene	3.950	54	1294497	22.15	ppbV	#	76
8) n-Butane	3.995	43	2335614	22.08	ppbV		98
9) Bromomethane	4.197	94	1033740	19.58	ppbV		99
10) Chloroethane	4.342	64	779909	20.76	ppbV		91
11) Ethanol	4.380	45	626678	18.27	ppbV		98
12) Vinyl bromide	4.638	106	1345719	22.49	ppbV		96
13) Acrolein	4.702	56	644716	24.68	ppbV		97
14) Acetone	4.802	58	872734	23.01	ppbV	#	68
15) Trichlorofluoromethane	4.982	101	3652093	22.07	ppbV		100
16) Isopropanol	4.975	45	2523205	20.40	ppbV		94
17) n-Pentane	5.287	43	2701076	22.47	ppbV		91
18) 1,1-Dichloroethene	5.538	61	2627511	22.28	ppbV		98
19) Methylene chloride	5.657	49	1904128	22.27	ppbV	#	78
20) Tert-butyl alcohol	5.506	59	3811059	24.85	ppbV		100
21) Allyl chloride	5.747	76	726511	22.42	ppbV		100
22) 1,1,2-Trichloro-1,2,2-...	5.879	101	2822496	21.84	ppbV		89
23) Carbon disulfide	5.933	76	4469998	22.21	ppbV	#	89
24) 1,2-Dichloroethene (tr...	6.506	61	2422888	22.51	ppbV		90
25) 1,1-Dichloroethane	6.695	63	3067293	22.43	ppbV		99
26) Methyl tert-butyl ether	6.724	73	4346736	22.30	ppbV		100
27) Methyl ethyl ketone	7.039	43	3149923	23.09	ppbV	#	83
28) 1,2-Dichloroethene (cis)	7.515	61	2355895	22.58	ppbV		89
29) Ethyl acetate	7.705	45	533489	23.24	ppbV		100
30) n-Hexane	7.737	57	2649428	22.68	ppbV	#	80
31) Chloroform	7.811	83	3281622	22.21	ppbV		99
32) Tetrahydrofuran	8.194	42	1759578	25.02	ppbV		94
33) 1,2-Dichloroethane	8.576	62	2503054	22.54	ppbV		99
34) 1,1,1-Trichloroethane	8.856	97	3122551	22.25	ppbV		99
35) Benzene	9.351	78	4929159	22.75	ppbV	#	89
36) Carbon tetrachloride	9.518	117	2937632	22.20	ppbV		100
37) Cyclohexane	9.663	56	2710024	22.12	ppbV		97
39) 1,2-Dichloropropane	10.258	63	2107013	22.26	ppbV		99
40) Bromodichloromethane	10.480	83	3543384	22.00	ppbV		98
41) 2,2,4-Trimethylpentane	10.592	57	8830251	21.80	ppbV		98
42) Trichloroethene	10.535	130	1667714	21.40	ppbV		100
43) 1,4-Dioxane	10.496	88	946441	20.60	ppbV	#	80
44) Methyl methacrylate	10.750	41	2228132	22.39	ppbV	#	70
45) n-Heptane	10.888	43	2903562	21.89	ppbV	#	76
46) cis-1,3-Dichloropropene	11.544	75	3067165	22.50	ppbV		99
47) Methyl isobutyl ketone	11.567	43	3867016	21.80	ppbV	#	85
48) trans-1,3-Dichloropropene	12.181	75	2906551	20.68	ppbV	#	76
49) 1,1,2-Trichloroethane	12.393	97	2013416	22.17	ppbV		93
50) Toluene	12.743	91	5782558	22.00	ppbV		98
51) Methyl n-butyl ketone	13.062	43	3763374	22.43	ppbV	#	81

AL SDG #E18-06141

Data Path : C:\DATA\07-25-18\
 Data File : aa7973std02.D
 Acq On : 25 Jul 2018 10:29 am
 Operator : jls
 Sample : 20 ppbv Std
 Misc : CC483586
 ALS Vial : 3 Sample Multiplier: 1

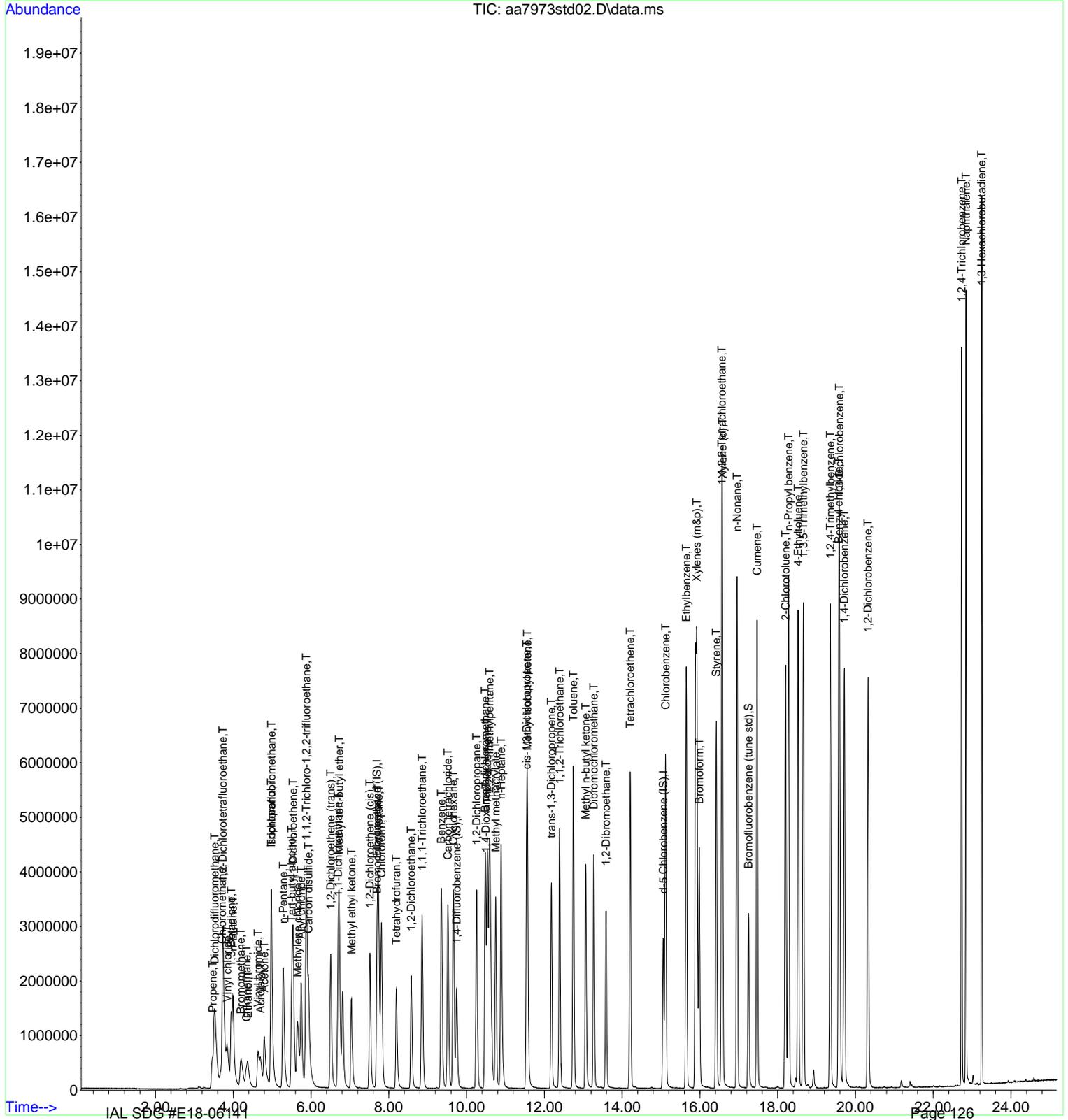
Quant Time: Jul 25 11:48:30 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 11:48:15 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
52) Dibromochloromethane	13.271	129	2923392	22.24	ppbV	99
53) 1,2-Dibromoethane	13.586	107	3121800	22.37	ppbV	96
54) Tetrachloroethene	14.206	166	2029587	21.71	ppbV	97
56) Chlorobenzene	15.116	112	4266769	21.52	ppbV	94
57) Ethylbenzene	15.650	91	7658674	21.17	ppbV	99
58) Xylenes (m&p)	15.917	91	11694300	39.41	ppbV	100
59) Bromoform	15.984	173	2467470	21.31	ppbV	98
60) Styrene	16.422	104	4438818	21.74	ppbV	96
61) Xylene (o)	16.579	91	6000746	20.80	ppbV	99
62) 1,1,2,2-Tetrachloroethane	16.560	83	4734111	20.94	ppbV	98
63) n-Nonane	16.955	43	4515097	20.75	ppbV #	77
65) Cumene	17.470	105	7699058	20.90	ppbV	97
66) 2-Chlorotoluene	18.203	91	6371298	21.16	ppbV	97
67) n-Propyl benzene	18.280	91	10455124	20.67	ppbV	96
68) 4-Ethyltoluene	18.528	105	7595419	21.06	ppbV	96
69) 1,3,5-Trimethylbenzene	18.663	105	6835220	20.81	ppbV	99
70) 1,2,4-Trimethylbenzene	19.354	105	6944267	21.10	ppbV	99
71) Benzyl chloride	19.570	91	6483719	20.08	ppbV	97
72) 1,3-Dichlorobenzene	19.589	146	3762437	20.69	ppbV	98
73) 1,4-Dichlorobenzene	19.714	146	3983041	21.07	ppbV	98
74) 1,2-Dichlorobenzene	20.325	146	3878212	21.08	ppbV	100
75) 1,2,4-Trichlorobenzene	22.730	180	3148298	22.38	ppbV	99
76) Naphthalene	22.843	127	1181278	23.92	ppbV	100
77) 1,3-Hexachlorobutadiene	23.248	225	1892848	19.29	ppbV	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\07-25-18\
 Data File : aa7973std02.D
 Acq On : 25 Jul 2018 10:29 am
 Operator : jls
 Sample : 20 ppbv Std
 Misc : CC483586
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Jul 25 11:48:30 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 11:48:15 2018
 Response via : Initial Calibration



Data Path : C:\DATA\07-25-18\
 Data File : aa7974std03.D
 Acq On : 25 Jul 2018 11:02 am
 Operator : jls
 Sample : 10 ppbv Std
 Misc : CC483586
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jul 25 11:47:45 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 11:47:37 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.689	130	485778	10.00	ppbV	0.00	
38) 1,4-Difluorobenzene (IS)	9.740	114	2080600	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.058	117	1771224	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...)	17.248	95	1610179	10.00	ppbV	0.00	
Target Compounds							
							Qvalue
2) Propene	3.457	41	485401	11.00	ppbV		# 80
3) Dichlorodifluoromethane	3.519	85	1721657	11.00	ppbV		97
4) Chloromethane	3.692	52	219571	11.00	ppbV	#	1
5) 1,2-Dichlorotetrafluor...	3.728	85	1683545	10.00	ppbV		95
6) Vinyl chloride	3.846	62	788471	11.00	ppbV		87
7) 1,3-Butadiene	3.943	54	641540	11.00	ppbV	#	77
8) n-Butane	3.991	43	1161584	11.00	ppbV		97
9) Bromomethane	4.194	94	521785	9.90	ppbV		99
10) Chloroethane	4.335	64	375017	10.00	ppbV		93
11) Ethanol	4.377	45	321790	9.40	ppbV		100
12) Vinyl bromide	4.637	106	656891	11.00	ppbV		97
13) Acrolein	4.695	56	312853	12.00	ppbV		96
14) Acetone	4.795	58	416502	11.00	ppbV		72
15) Trichlorofluoromethane	4.978	101	1816579	11.00	ppbV		100
16) Isopropanol	4.975	45	1234391	10.00	ppbV		93
17) n-Pentane	5.284	43	1319811	11.00	ppbV		92
18) 1,1-Dichloroethene	5.535	61	1295082	11.00	ppbV		98
19) Methylene chloride	5.650	49	938843	11.00	ppbV	#	78
20) Tert-butyl alcohol	5.502	59	1837100	12.00	ppbV		100
21) Allyl chloride	5.744	76	355787	11.00	ppbV		100
22) 1,1,2-Trichloro-1,2,2-...	5.879	101	1418961	11.00	ppbV		89
23) Carbon disulfide	5.933	76	2209326	11.00	ppbV	#	91
24) 1,2-Dichloroethene (tr...	6.506	61	1181605	11.00	ppbV		89
25) 1,1-Dichloroethane	6.692	63	1501637	11.00	ppbV		99
26) Methyl tert-butyl ether	6.724	73	2140023	11.00	ppbV		100
27) Methyl ethyl ketone	7.039	43	1497882	11.00	ppbV	#	83
28) 1,2-Dichloroethene (cis)	7.515	61	1145414	11.00	ppbV		90
29) Ethyl acetate	7.705	45	252052	11.00	ppbV		100
30) n-Hexane	7.734	57	1282758	11.00	ppbV	#	81
31) Chloroform	7.808	83	1622409	11.00	ppbV		98
32) Tetrahydrofuran	8.197	42	842429	12.00	ppbV		95
33) 1,2-Dichloroethane	8.576	62	1219030	11.00	ppbV		99
34) 1,1,1-Trichloroethane	8.853	97	1540529	11.00	ppbV		99
35) Benzene	9.348	78	2378901	11.00	ppbV	#	89
36) Carbon tetrachloride	9.518	117	1452782	11.00	ppbV		100
37) Cyclohexane	9.660	56	1344949	11.00	ppbV		97
39) 1,2-Dichloropropane	10.258	63	1017552	11.00	ppbV		99
40) Bromodichloromethane	10.480	83	1731985	11.00	ppbV		98
41) 2,2,4-Trimethylpentane	10.595	57	4354544	11.00	ppbV		98
42) Trichloroethene	10.538	130	837753	11.00	ppbV		99
43) 1,4-Dioxane	10.499	88	449087	10.00	ppbV		82
44) Methyl methacrylate	10.750	41	1070120	11.00	ppbV	#	71
45) n-Heptane	10.888	43	1426371	11.00	ppbV	#	76
46) cis-1,3-Dichloropropene	11.547	75	1465634	11.00	ppbV		99
47) Methyl isobutyl ketone	11.570	43	1906864	11.00	ppbV	#	86
48) trans-1,3-Dichloropropene	12.184	75	1373612	10.00	ppbV	#	77
49) 1,1,2-Trichloroethane	12.393	97	976409	11.00	ppbV		93
50) Toluene	12.743	91	2826398	11.00	ppbV		97
51) Methyl n-butyl ketone	13.065	43	1803996	11.00	ppbV	#	83

AL SDG #E18-06141

Data Path : C:\DATA\07-25-18\
 Data File : aa7974std03.D
 Acq On : 25 Jul 2018 11:02 am
 Operator : jls
 Sample : 10 ppbv Std
 Misc : CC483586
 ALS Vial : 4 Sample Multiplier: 1

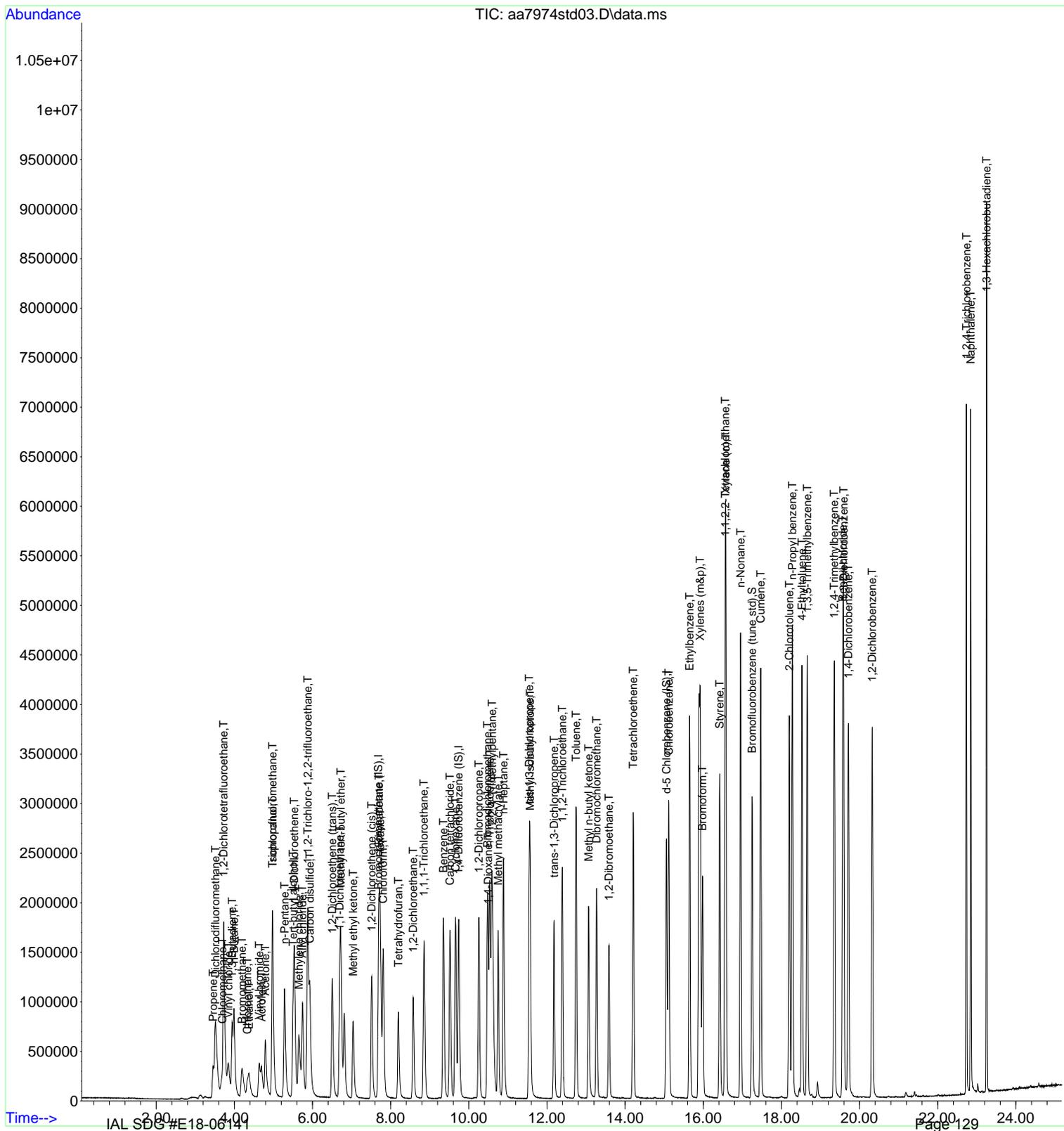
Quant Time: Jul 25 11:47:45 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 11:47:37 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Dibromochloromethane	13.271	129	1413643	11.00	ppbV	99
53) 1,2-Dibromoethane	13.586	107	1500610	11.00	ppbV	96
54) Tetrachloroethene	14.206	166	1005355	11.00	ppbV	98
56) Chlorobenzene	15.116	112	2070449	11.00	ppbV	93
57) Ethylbenzene	15.650	91	3778299	11.00	ppbV	98
58) Xylenes (m&p)	15.917	91	6197500	22.00	ppbV	96
59) Bromoform	15.981	173	1209271	11.00	ppbV	97
60) Styrene	16.422	104	2131498	11.00	ppbV	96
61) Xylene (o)	16.576	91	3011718	11.00	ppbV	100
62) 1,1,2,2-Tetrachloroethane	16.560	83	2360623	11.00	ppbV	98
63) n-Nonane	16.955	43	2272106	11.00	ppbV #	79
65) Cumene	17.470	105	3846808	11.00	ppbV	97
66) 2-Chlorotoluene	18.203	91	3143676	11.00	ppbV	97
67) n-Propyl benzene	18.280	91	5281946	11.00	ppbV	95
68) 4-Ethyltoluene	18.528	105	3765828	11.00	ppbV	96
69) 1,3,5-Trimethylbenzene	18.659	105	3429338	11.00	ppbV	100
70) 1,2,4-Trimethylbenzene	19.351	105	3436295	11.00	ppbV	99
71) Benzyl chloride	19.569	91	3064617	10.00	ppbV	97
72) 1,3-Dichlorobenzene	19.592	146	1898980	11.00	ppbV	98
73) 1,4-Dichlorobenzene	19.714	146	1974346	11.00	ppbV	98
74) 1,2-Dichlorobenzene	20.325	146	1921004	11.00	ppbV	99
75) 1,2,4-Trichlorobenzene	22.733	180	1602141	12.00	ppbV	99
76) Naphthalene	22.843	127	562619	12.00	ppbV	100
77) 1,3-Hexachlorobutadiene	23.251	225	1024427	11.00	ppbV	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\07-25-18\
 Data File : aa7974std03.D
 Acq On : 25 Jul 2018 11:02 am
 Operator : jls
 Sample : 10 ppbv Std
 Misc : CC483586
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jul 25 11:47:45 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 11:47:37 2018
 Response via : Initial Calibration



Data Path : C:\DATA\07-25-18\
 Data File : aa7975std04.D
 Acq On : 25 Jul 2018 12:18 pm
 Operator : jls
 Sample : 2 ppbv Std
 Misc : CC483586
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Jul 25 12:45:06 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 12:03:50 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.679	130	442159	10.00	ppbV	0.00	
38) 1,4-Difluorobenzene (IS)	9.737	114	1836044	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.055	117	1531011	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...	17.251	95	1404561	10.07	ppbV	0.00	
Target Compounds							
							Qvalue
2) Propene	3.448	41	78631	1.98	ppbV	#	80
3) Dichlorodifluoromethane	3.519	85	261761	1.88	ppbV		97
4) Chloromethane	3.683	52	36696	2.10	ppbV	#	1
5) 1,2-Dichlorotetrafluor...	3.721	85	291795	1.99	ppbV		97
6) Vinyl chloride	3.840	62	118454	1.87	ppbV		89
7) 1,3-Butadiene	3.937	54	100761	1.96	ppbV		84
8) n-Butane	3.988	43	195789	2.09	ppbV		94
9) Bromomethane	4.187	94	79223	1.76	ppbV		98
10) Chloroethane	4.329	64	55246	1.64	ppbV	#	88
11) Ethanol	4.371	45	85520	2.37	ppbV		99
12) Vinyl bromide	4.625	106	98171	1.85	ppbV		96
13) Acrolein	4.692	56	47056	2.05	ppbV		97
14) Acetone	4.789	58	62165	1.84	ppbV		75
15) Trichlorofluoromethane	4.975	101	305604	2.06	ppbV		99
16) Isopropanol	4.969	45	199685	1.81	ppbV		92
17) n-Pentane	5.281	43	207084	1.94	ppbV		94
18) 1,1-Dichloroethene	5.525	61	196205	1.91	ppbV		100
19) Methylene chloride	5.647	49	153043	2.01	ppbV	#	83
20) Tert-butyl alcohol	5.502	59	299995	2.15	ppbV		100
21) Allyl chloride	5.737	76	50997	1.83	ppbV		100
22) 1,1,2-Trichloro-1,2,2-...	5.872	101	273970	2.33	ppbV		89
23) Carbon disulfide	5.927	76	331928	1.89	ppbV	#	91
24) 1,2-Dichloroethene (tr...	6.499	61	175124	1.85	ppbV		90
25) 1,1-Dichloroethane	6.689	63	232574	1.92	ppbV		99
26) Methyl tert-butyl ether	6.727	73	371042	2.11	ppbV		99
27) Methyl ethyl ketone	7.043	43	227065	1.85	ppbV	#	86
28) 1,2-Dichloroethene (cis)	7.512	61	171952	1.87	ppbV		91
29) Ethyl acetate	7.711	45	38370	1.85	ppbV		100
30) n-Hexane	7.727	57	211648	2.01	ppbV		84
31) Chloroform	7.805	83	246384	1.90	ppbV		99
32) Tetrahydrofuran	8.200	42	124601	1.99	ppbV		91
33) 1,2-Dichloroethane	8.573	62	189562	1.91	ppbV		100
34) 1,1,1-Trichloroethane	8.846	97	271151	2.14	ppbV		100
35) Benzene	9.345	78	381445	1.96	ppbV		91
36) Carbon tetrachloride	9.515	117	253566	2.12	ppbV		99
37) Cyclohexane	9.653	56	232948	2.09	ppbV		98
39) 1,2-Dichloropropane	10.251	63	169395	2.11	ppbV		100
40) Bromodichloromethane	10.473	83	275138	2.08	ppbV		97
41) 2,2,4-Trimethylpentane	10.586	57	694979	2.10	ppbV		98
42) Trichloroethene	10.531	130	132953	2.08	ppbV		99
43) 1,4-Dioxane	10.505	88	72147	1.85	ppbV		83
44) Methyl methacrylate	10.750	41	175407	2.08	ppbV	#	75
45) n-Heptane	10.885	43	261435	2.30	ppbV	#	79
46) cis-1,3-Dichloropropene	11.538	75	225544	1.98	ppbV		98
47) Methyl isobutyl ketone	11.570	43	342759	2.29	ppbV		89
48) trans-1,3-Dichloropropene	12.181	75	208336	1.77	ppbV	#	81
49) 1,1,2-Trichloroethane	12.390	97	170605	2.22	ppbV		94
50) Toluene	12.743	91	488169	2.20	ppbV		98
51) Methyl n-butyl ketone	13.071	43	299937	2.10	ppbV	#	85

AL SDG #E18-06141

Data Path : C:\DATA\07-25-18\
 Data File : aa7975std04.D
 Acq On : 25 Jul 2018 12:18 pm
 Operator : jls
 Sample : 2 ppbv Std
 Misc : CC483586
 ALS Vial : 5 Sample Multiplier: 1

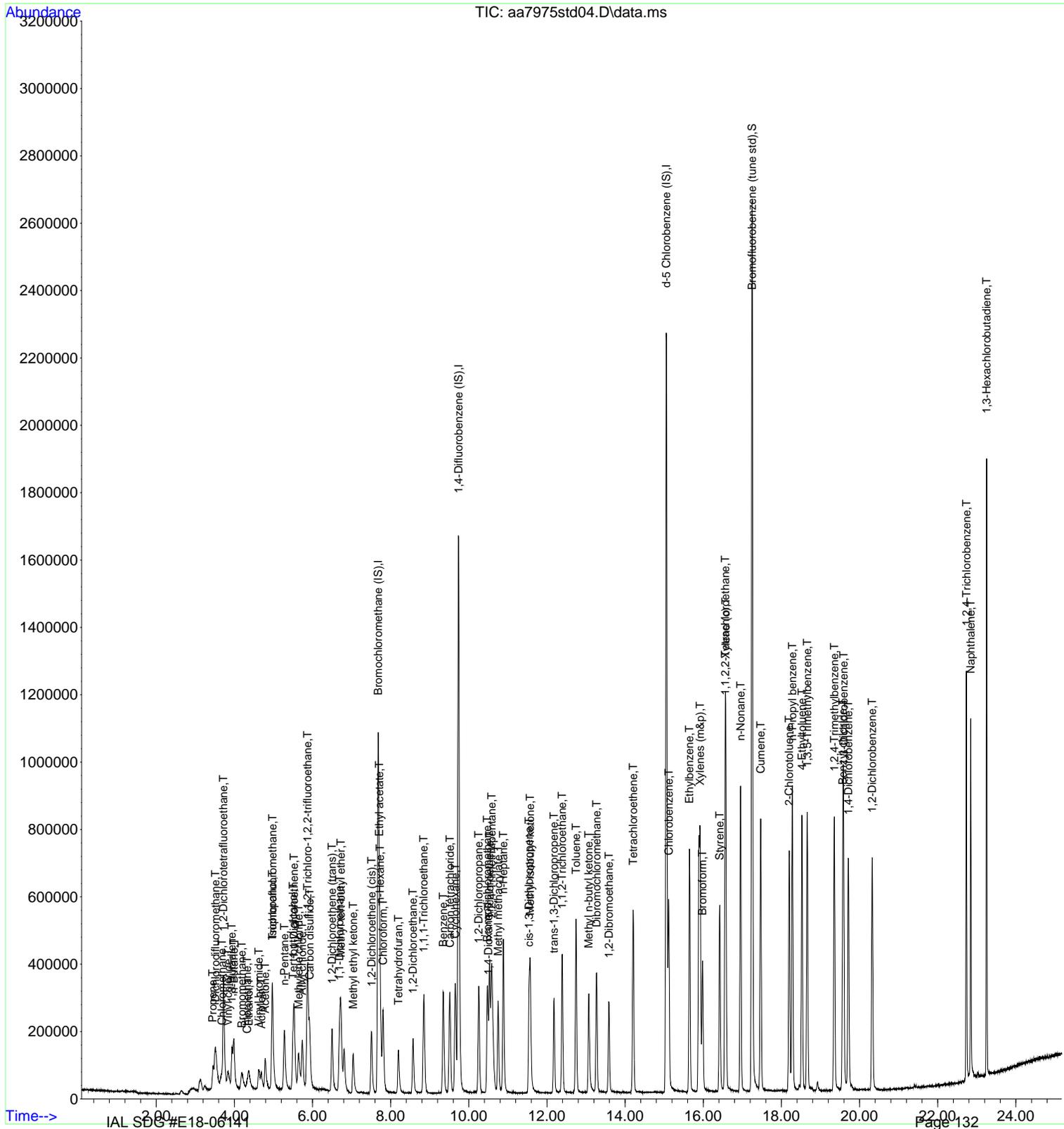
Quant Time: Jul 25 12:45:06 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 12:03:50 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Dibromochloromethane	13.271	129	239686	2.17	ppbV	99
53) 1,2-Dibromoethane	13.586	107	255131	2.15	ppbV	96
54) Tetrachloroethene	14.206	166	183734	2.33	ppbV	99
56) Chlorobenzene	15.116	112	366932	2.31	ppbV	93
57) Ethylbenzene	15.647	91	697526	2.42	ppbV	98
58) Xylenes (m&p)	15.920	91	1083055	4.74	ppbV	99
59) Bromoform	15.981	173	207538	2.29	ppbV	98
60) Styrene	16.422	104	357882	2.21	ppbV	97
61) Xylene (o)	16.573	91	565479	2.47	ppbV	100
62) 1,1,2,2-Tetrachloroethane	16.560	83	453501	2.50	ppbV	98
63) n-Nonane	16.955	43	437271	2.53	ppbV #	82
65) Cumene	17.470	105	707068	2.43	ppbV	97
66) 2-Chlorotoluene	18.200	91	571345	2.38	ppbV	98
67) n-Propyl benzene	18.280	91	986527	2.49	ppbV	95
68) 4-Ethyltoluene	18.524	105	683257	2.40	ppbV	97
69) 1,3,5-Trimethylbenzene	18.663	105	621617	2.42	ppbV	98
70) 1,2,4-Trimethylbenzene	19.354	105	617452	2.38	ppbV	98
71) Benzyl chloride	19.569	91	479594	1.89	ppbV	97
72) 1,3-Dichlorobenzene	19.592	146	355664	2.49	ppbV	98
73) 1,4-Dichlorobenzene	19.714	146	362548	2.43	ppbV	96
74) 1,2-Dichlorobenzene	20.322	146	353034	2.42	ppbV	98
75) 1,2,4-Trichlorobenzene	22.733	180	282410	2.61	ppbV	98
76) Naphthalene	22.846	127	87604	2.29	ppbV	100
77) 1,3-Hexachlorobutadiene	23.251	225	204886	2.68	ppbV	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\07-25-18\
 Data File : aa7975std04.D
 Acq On : 25 Jul 2018 12:18 pm
 Operator : jls
 Sample : 2 ppbv Std
 Misc : CC483586
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Jul 25 12:45:06 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 12:03:50 2018
 Response via : Initial Calibration



Data Path : C:\DATA\07-25-18\
 Data File : aa7976std05.D
 Acq On : 25 Jul 2018 1:42 pm
 Operator : jls
 Sample : 0.2 ppbv Std
 Misc : CC483586
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Jul 25 14:07:24 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 13:25:42 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.679	130	505975	10.00	ppbV	0.00	
38) 1,4-Difluorobenzene (IS)	9.734	114	2064481	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.055	117	1704012	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...	17.248	95	1569127	10.08	ppbV	0.00	
Target Compounds							
							Qvalue
2) Propene	3.454	41	7319	0.14	ppbV	#	4
3) Dichlorodifluoromethane	3.522	85	37924	0.22	ppbV	#	86
4) Chloromethane	3.673	52	4895m	0.26	ppbV		
5) 1,2-Dichlorotetrafluor...	3.721	85	41484	0.22	ppbV		99
6) Vinyl chloride	3.840	62	15540	0.20	ppbV		90
7) 1,3-Butadiene	3.937	54	12107	0.19	ppbV		95
8) n-Butane	3.985	43	28051	0.23	ppbV		96
9) Bromomethane	4.194	94	9614	0.17	ppbV	#	59
10) Chloroethane	4.313	64	6976	0.19	ppbV	#	59
11) Ethanol	4.374	45	7947m	0.20	ppbV		
12) Vinyl bromide	4.622	106	10734	0.17	ppbV	#	42
13) Acrolein	4.689	56	4231	0.15	ppbV	#	51
14) Acetone	4.805	58	8147	0.20	ppbV		81
15) Trichlorofluoromethane	4.969	101	41547	0.22	ppbV		98
16) Isopropanol	4.975	45	23985	0.17	ppbV	#	88
17) n-Pentane	5.284	43	30958	0.23	ppbV		90
18) 1,1-Dichloroethene	5.531	61	25377	0.20	ppbV		94
19) Methylene chloride	5.644	49	29180	0.28	ppbV	#	81
20) Tert-butyl alcohol	5.509	59	39682	0.23	ppbV		100
21) Allyl chloride	5.744	76	6789	0.20	ppbV		100
22) 1,1,2-Trichloro-1,2,2-...	5.872	101	35414	0.23	ppbV		88
23) Carbon disulfide	5.917	76	41730	0.19	ppbV	#	81
24) 1,2-Dichloroethene (tr...	6.499	61	22565	0.19	ppbV	#	82
25) 1,1-Dichloroethane	6.692	63	31076	0.20	ppbV		99
26) Methyl tert-butyl ether	6.728	73	45176	0.21	ppbV		100
27) Methyl ethyl ketone	7.056	43	32178	0.21	ppbV		95
28) 1,2-Dichloroethene (cis)	7.509	61	21368	0.19	ppbV	#	84
29) Ethyl acetate	7.721	45	5075	0.20	ppbV		100
30) n-Hexane	7.731	57	27375	0.21	ppbV		88
31) Chloroform	7.808	83	33786	0.21	ppbV		99
32) Tetrahydrofuran	8.210	42	17131	0.22	ppbV	#	85
33) 1,2-Dichloroethane	8.570	62	25895	0.21	ppbV		94
34) 1,1,1-Trichloroethane	8.850	97	35176	0.22	ppbV		99
35) Benzene	9.342	78	48602	0.20	ppbV		96
36) Carbon tetrachloride	9.502	117	32991	0.22	ppbV		94
37) Cyclohexane	9.647	56	28043	0.21	ppbV		98
39) 1,2-Dichloropropane	10.252	63	23544	0.23	ppbV		98
40) Bromodichloromethane	10.477	83	34629	0.21	ppbV		98
41) 2,2,4-Trimethylpentane	10.586	57	83595	0.20	ppbV		99
42) Trichloroethene	10.531	130	19056	0.24	ppbV		92
43) 1,4-Dioxane	10.522	88	7613	0.17	ppbV	#	36
44) Methyl methacrylate	10.756	41	23522	0.22	ppbV	#	79
45) n-Heptane	10.885	43	31156	0.21	ppbV	#	77
46) cis-1,3-Dichloropropene	11.541	75	28118	0.20	ppbV		98
47) Methyl isobutyl ketone	11.589	43	41480	0.22	ppbV		89
48) trans-1,3-Dichloropropene	12.181	75	25807	0.18	ppbV		86
49) 1,1,2-Trichloroethane	12.390	97	22149	0.23	ppbV		92
50) Toluene	12.740	91	62307	0.22	ppbV		96
51) Methyl n-butyl ketone	13.084	43	37567	0.21	ppbV		96

AL SDG #E18-06141

Data Path : C:\DATA\07-25-18\
 Data File : aa7976std05.D
 Acq On : 25 Jul 2018 1:42 pm
 Operator : jls
 Sample : 0.2 ppbv Std
 Misc : CC483586
 ALS Vial : 6 Sample Multiplier: 1

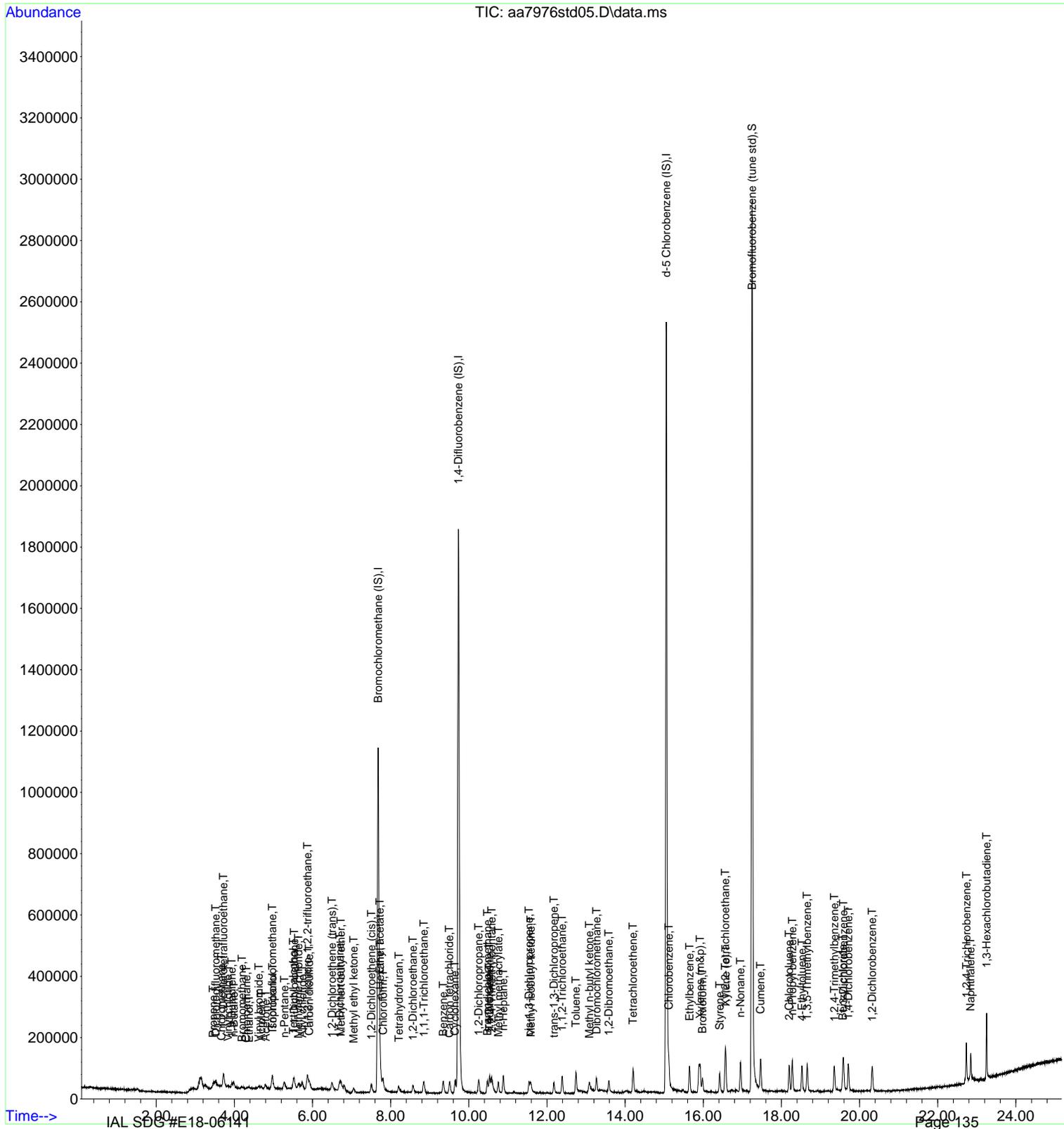
Quant Time: Jul 25 14:07:24 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 13:25:42 2018
 Response via : Initial Calibration

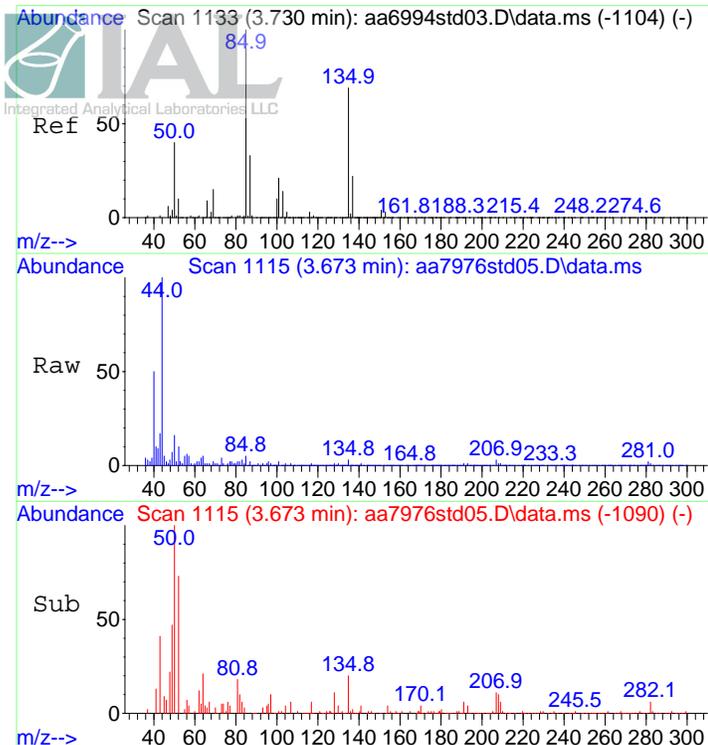
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Dibromochloromethane	13.267	129	28953	0.21	ppbV	100
53) 1,2-Dibromoethane	13.583	107	33845	0.23	ppbV	95
54) Tetrachloroethene	14.203	166	24894	0.24	ppbV	96
56) Chlorobenzene	15.116	112	47512	0.24	ppbV	95
57) Ethylbenzene	15.650	91	84088	0.23	ppbV	98
58) Xylenes (m&p)	15.917	91	128386	0.44	ppbV #	28
59) Bromoform	15.981	173	23394	0.21	ppbV	98
60) Styrene	16.422	104	39109	0.20	ppbV	99
61) Xylene (o)	16.573	91	65305	0.23	ppbV	100
62) 1,1,2,2-Tetrachloroethane	16.560	83	54745	0.24	ppbV	97
63) n-Nonane	16.952	43	48551	0.22	ppbV	90
65) Cumene	17.467	105	80967	0.22	ppbV	98
66) 2-Chlorotoluene	18.200	91	68160	0.23	ppbV	99
67) n-Propyl benzene	18.280	91	107193	0.22	ppbV	97
68) 4-Ethyltoluene	18.525	105	73402	0.21	ppbV	97
69) 1,3,5-Trimethylbenzene	18.656	105	70531	0.22	ppbV	98
70) 1,2,4-Trimethylbenzene	19.354	105	64555	0.20	ppbV	98
71) Benzyl chloride	19.569	91	42999	0.15	ppbV	96
72) 1,3-Dichlorobenzene	19.589	146	44070	0.24	ppbV	100
73) 1,4-Dichlorobenzene	19.711	146	43316	0.22	ppbV	94
74) 1,2-Dichlorobenzene	20.322	146	43839	0.24	ppbV	96
75) 1,2,4-Trichlorobenzene	22.733	180	38969	0.26	ppbV	95
76) Naphthalene	22.839	127	10301	0.23	ppbV	100
77) 1,3-Hexachlorobutadiene	23.251	225	28434	0.28	ppbV	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\07-25-18\
 Data File : aa7976std05.D
 Acq On : 25 Jul 2018 1:42 pm
 Operator : jls
 Sample : 0.2 ppbv Std
 Misc : CC483586
 ALS Vial : 6 Sample Multiplier: 1

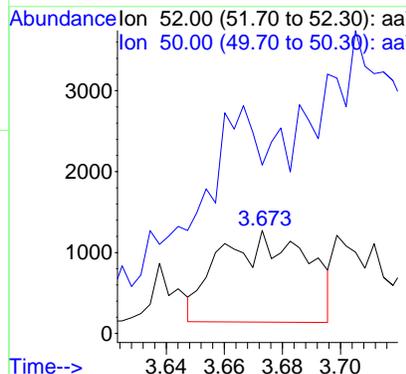
Quant Time: Jul 25 14:07:24 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 13:25:42 2018
 Response via : Initial Calibration



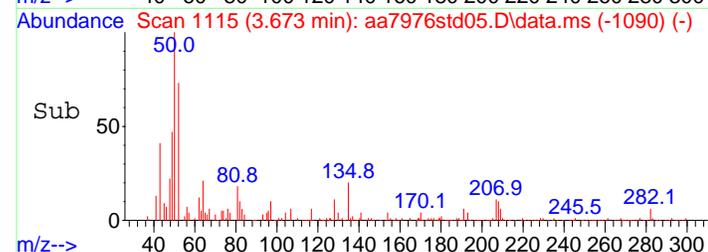
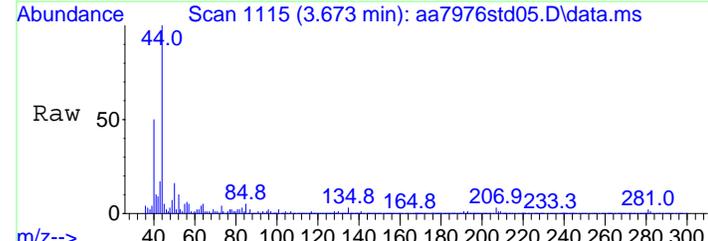
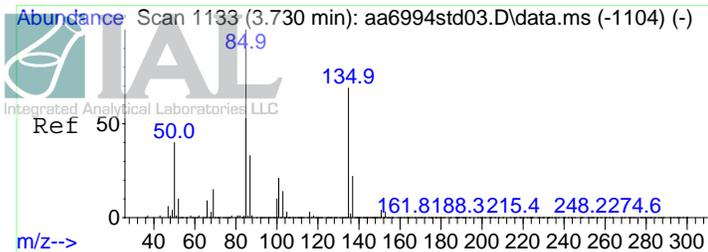


#4
 Chloromethane
 Concen: 0.12 ppbV
 RT: 3.673 min Scan# 1115
 Delta R.T. -0.019 min
 Lab File: aa7976std05.D
 Acq: 25 Jul 2018 1:42 pm

Tgt Ion	Resp	Lower	Upper
52	2327		
52	100		
50	282.1	314.4	471.6#

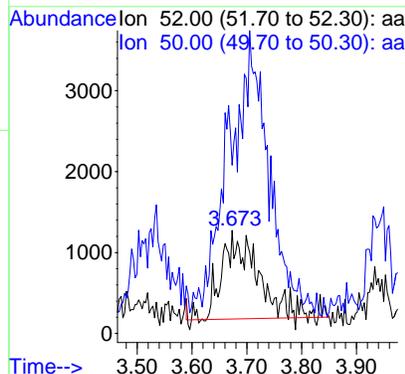


Before manual integration for Chloromethane.

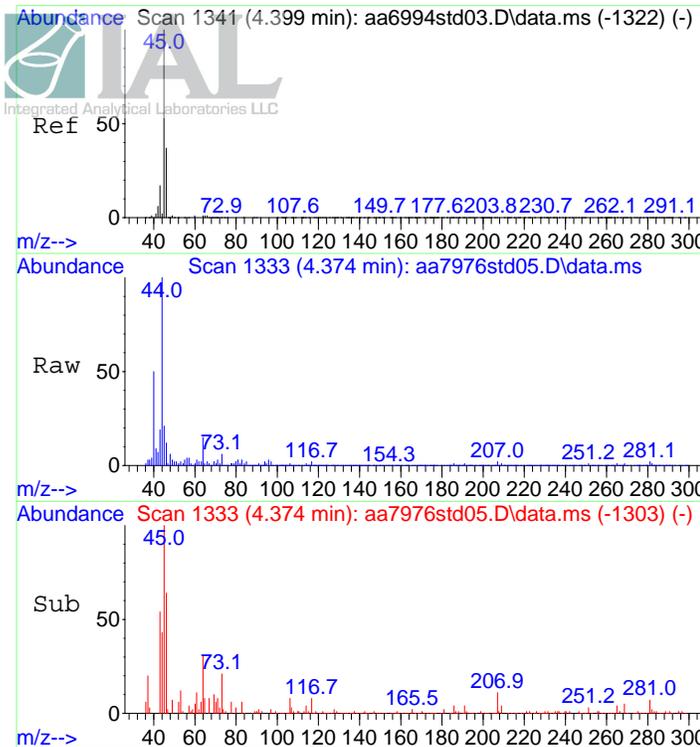


#4
 Chloromethane
 Concen: 0.26 ppbV m
 RT: 3.673 min Scan# 1115
 Delta R.T. -0.019 min
 Lab File: aa7976std05.D
 Acq: 25 Jul 2018 1:42 pm

Tgt Ion	Resp	Lower	Upper
52	4895		
52	100		
50	134.1	314.4	471.6#

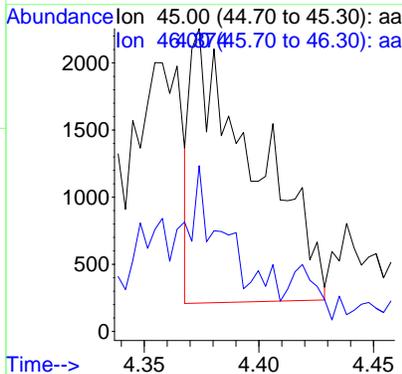


After manual integration for Chloromethane.

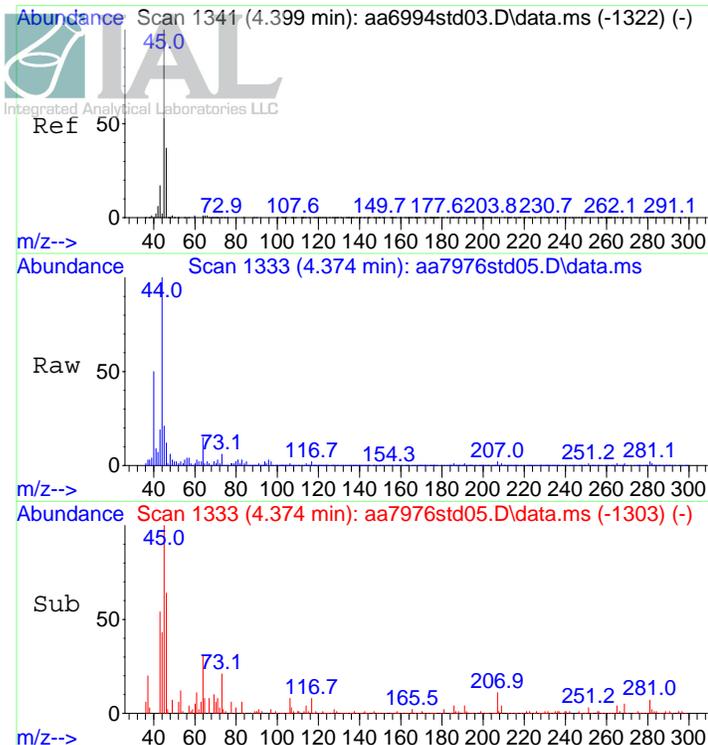


#11
Ethanol
Concen: 0.10 ppbV
RT: 4.374 min Scan# 1333
Delta R.T. -0.003 min
Lab File: aa7976std05.D
Acq: 25 Jul 2018 1:42 pm

Tgt Ion:	45	Resp:	3891
Ion Ratio	Lower	Upper	
45	100		
46	36.8	32.6	48.8

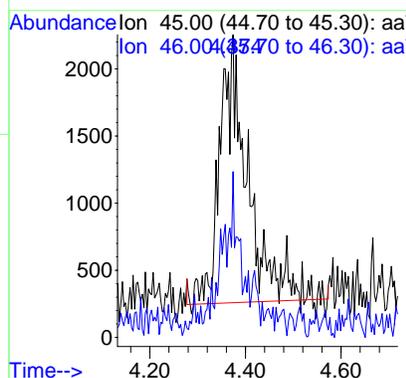


Before manual integration for Ethanol.



#11
Ethanol
Concen: 0.20 ppbV m
RT: 4.374 min Scan# 1333
Delta R.T. -0.003 min
Lab File: aa7976std05.D
Acq: 25 Jul 2018 1:42 pm

Tgt Ion:	45	Resp:	7947
Ion Ratio	Lower	Upper	
45	100		
46	18.0	32.6	48.8#



After manual integration for Ethanol.

Initial Calibration Verification Sample Standard

Lab Sample Name: 10 PPBV ICVSS
Spike Amount: 10 ppbv, except m&p-Xylenes at 20 ppbv
Amount of standard injected (ml): 50

Data File: AA7078ICVSS
Date Analyzed: 5/18/2018

Runs with this ICVSS:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7071BFB]	05/18/2018 8:17
40 PPBV STD [AA7073STD01]	05/18/2018 9:46
20 PPBV STD [AA7074STD02]	05/18/2018 10:19
10 PPBV STD [AA7075STD03]	05/18/2018 10:53
2 PPBV STD [AA7076STD04]	05/18/2018 12:20
0.2 PPBV STD [AA7077STD05]	05/18/2018 13:03
10 PPBV ICVSS [AA7078ICVSS]	05/18/2018 15:14

Compound	CAS #	Injected Amount (ppbv)	Recovered Amount (ppbv)	% Recovery	QC Limit
Acetone	67-64-1	11	11	100	
Acrolein	107-02-8	7.1	8.6	120	
Allyl Chloride	107-05-1	11	11	100	
Benzene	71-43-2	10	11	110	
Benzyl chloride	100-44-7	11	11	100	
Bromodichloromethane	75-27-4	11	11	100	
Bromoform	75-25-2	11	11	100	
Bromomethane	74-83-9	11	10	91	
1,3-Butadiene	106-99-0	11	11	100	
n-Butane	106-97-8	11	11	100	
Chlorobenzene	108-90-7	10	10.0	100	
Chloroethane	75-00-3	11	9.4	85	
Chloroform	67-66-3	11	11	100	
Chloromethane	74-87-3	10	11	110	
Carbon disulfide	75-15-0	11	11	100	
Carbon tetrachloride	56-23-5	11	10	91	
2-Chlorotoluene	95-49-8	11	11	100	
Cumene	98-82-8	11	11	100	
Cyclohexane	110-82-7	11	11	100	
Dibromochloromethane	124-48-1	11	10	91	
1,2-Dibromoethane	106-93-4	10	10	100	
1,2-Dichlorobenzene	95-50-1	10	10	100	
1,3-Dichlorobenzene	541-73-1	10	10	100	
1,4-Dichlorobenzene	106-46-7	10	10	100	
Dichlorodifluoromethane	75-71-8	11	11	100	
1,1-Dichloroethane	75-34-3	11	11	100	
1,2-Dichloroethane	107-06-2	11	11	100	
1,1-Dichloroethene	75-35-4	11	11	100	
1,2-Dichloroethene (cis)	156-59-2	10	11	110	
1,2-Dichloroethene (trans)	156-60-5	11	11	100	
1,2-Dichloropropane	78-87-5	11	10	91	
1,3-Dichloropropene (cis)	10061-01-5	11	11	100	
1,3-Dichloropropene (trans)	10061-02-6	10	10	100	
1,2-Dichlorotetrafluoroethane	76-14-2	11	9.0	82	
1,4-Dioxane	123-91-1	8.3	9.7	120	
Ethanol	64-17-5	8.7	8.0	92	
Ethyl acetate	141-78-6	11	11	100	
Ethylbenzene	100-41-4	11	11	100	
4-Ethyltoluene	622-96-8	11	11	100	
n-Heptane	142-82-5	11	11	100	

ICVSS recovery must be within 70-130% of the spiked value for all compounds.

*** Values outside of QC limits**

Initial Calibration Verification Sample Standard

Lab Sample Name: 10 PPBV ICVSS
Spike Amount: 10 ppbv, except m&p-Xylenes at 20 ppbv
Amount of standard injected (ml): 50

Data File: AA7078ICVSS
Date Analyzed: 5/18/2018

Runs with this ICVSS:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7071BFB]	05/18/2018 8:17
40 PPBV STD [AA7073STD01]	05/18/2018 9:46
20 PPBV STD [AA7074STD02]	05/18/2018 10:19
10 PPBV STD [AA7075STD03]	05/18/2018 10:53
2 PPBV STD [AA7076STD04]	05/18/2018 12:20
0.2 PPBV STD [AA7077STD05]	05/18/2018 13:03
10 PPBV ICVSS [AA7078ICVSS]	05/18/2018 15:14

Compound	CAS #	Injected Amount (ppbv)	Recovered Amount (ppbv)	% Recovery	QC Limit
1,3-Hexachlorobutadiene	87-68-3	10	10.0	100	
n-Hexane	110-54-3	10	11	110	
Isopropanol	67-63-0	8.2	8.3	100	
Methylene chloride	75-09-2	11	10	91	
Methyl ethyl ketone	78-93-3	10	11	110	
Methyl isobutyl ketone	108-10-1	11	11	100	
Methyl methacrylate	80-62-6	11	11	100	
Methyl n-butyl ketone	591-78-6	9.4	11	120	
Methyl tert-butyl ether	1634-04-4	11	11	100	
Naphthalene	91-20-3	9.3	12	130	
n-Nonane	111-84-2	11	11	100	
n-Pentane	109-66-0	11	11	100	
Propene	115-07-1	11	11	100	
n-Propyl benzene	103-65-1	11	11	100	
Styrene	100-42-5	10	12	120	
Tert-butyl alcohol	75-65-0	8.0	8.5	110	
1,1,2,2-Tetrachloroethane	79-34-5	10	10	100	
Tetrachloroethene	127-18-4	10	10	100	
Tetrahydrofuran	109-99-9	11	12	110	
Toluene	108-88-3	10	11	110	
1,2,4-Trichlorobenzene	120-82-1	8.8	9.9	110	
1,1,1-Trichloroethane	71-55-6	11	10	91	
1,1,2-Trichloroethane	79-00-5	10	10	100	
Trichloroethene	79-01-6	10	11	110	
Trichlorofluoromethane	75-69-4	11	10	91	
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	11	10.0	91	
1,2,4-Trimethylbenzene	95-63-6	10	12	120	
1,3,5-Trimethylbenzene	108-67-8	11	11	100	
2,2,4-Trimethylpentane	540-84-1	11	11	100	
Vinyl bromide	593-60-2	9.7	12	120	
Vinyl chloride	75-01-4	11	11	100	
Xylenes (m&p)	179601-23-1	22	22	100	
Xylenes (o)	95-47-6	11	11	100	

ICVSS recovery must be within 70-130% of the spiked value for all compounds.

*** Values outside of QC limits**

Data Path : C:\DATA\05-18-18\
 Data File : aa70781cvss.D
 Acq On : 18 May 2018 3:14 pm
 Operator : jls
 Sample : 10 ppbv ICVSS
 Misc : CC483422
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: May 21 10:49:09 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 13:51:08 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.724	130	522049	10.00	ppbV	0.00	
38) 1,4-Difluorobenzene (IS)	9.782	114	2190215	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.113	117	1913238	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...	17.309	95	1757413	9.98	ppbV	0.00	
Target Compounds							
							Qvalue
2) Propene	3.473	41	730162	11.06	ppbV		83
3) Dichlorodifluoromethane	3.534	85	2027364	10.97	ppbV		97
4) Chloromethane	3.730	52	276132	10.79	ppbV		97
5) 1,2-Dichlorotetrafluor...	3.746	85	1759325	9.04	ppbV		94
6) Vinyl chloride	3.859	62	839310	11.09	ppbV		87
7) 1,3-Butadiene	3.962	54	733268	11.41	ppbV		96
8) n-Butane	4.013	43	1601575	10.78	ppbV		94
9) Bromomethane	4.213	94	555490	10.03	ppbV		97
10) Chloroethane	4.357	64	403118	9.41	ppbV		93
11) Ethanol	4.399	45	331087	8.00	ppbV		97
12) Vinyl bromide	4.656	106	624617	11.56	ppbV		100
13) Acrolein	4.724	56	228811	8.59	ppbV		97
14) Acetone	4.820	58	480248	11.38	ppbV		96
15) Trichlorofluoromethane	5.000	101	2123834	10.47	ppbV		99
16) Isopropanol	5.000	45	1136925	8.34	ppbV		93
17) n-Pentane	5.306	43	1783363	11.16	ppbV		95
18) 1,1-Dichloroethene	5.566	61	1496651	11.16	ppbV		92
19) Methylene chloride	5.679	49	1228354	10.08	ppbV		99
20) Tert-butyl alcohol	5.528	59	1031467	8.45	ppbV		100
21) Allyl chloride	5.775	76	333899	11.26	ppbV		100
22) 1,1,2-Trichloro-1,2,2-...	5.907	101	1565336	9.98	ppbV		88
23) Carbon disulfide	5.965	76	2223694	11.29	ppbV		97
24) 1,2-Dichloroethene (tr...	6.534	61	1304129	10.92	ppbV		99
25) 1,1-Dichloroethane	6.727	63	1656694	10.72	ppbV		99
26) Methyl tert-butyl ether	6.756	73	1903804	10.94	ppbV		96
27) Methyl ethyl ketone	7.071	43	2144860	11.22	ppbV		99
28) 1,2-Dichloroethene (cis)	7.550	61	1271567	11.06	ppbV		99
29) Ethyl acetate	7.743	45	348813	11.27	ppbV		100
30) n-Hexane	7.766	57	1425441	10.63	ppbV		97
31) Chloroform	7.843	83	1780945	10.78	ppbV		99
32) Tetrahydrofuran	8.235	42	1110856	12.25	ppbV		90
33) 1,2-Dichloroethane	8.614	62	1465335	10.58	ppbV		100
34) 1,1,1-Trichloroethane	8.888	97	1650398	10.39	ppbV		99
35) Benzene	9.389	78	2434532	10.64	ppbV		99
36) Carbon tetrachloride	9.560	117	1618838	10.33	ppbV		99
37) Cyclohexane	9.701	56	1491999	10.98	ppbV		89
39) 1,2-Dichloropropane	10.299	63	1087670	10.21	ppbV		99
40) Bromodichloromethane	10.521	83	1904826	10.55	ppbV		99
41) 2,2,4-Trimethylpentane	10.634	57	4504117	11.45	ppbV		96
42) Trichloroethene	10.579	130	888910	10.56	ppbV		98
43) 1,4-Dioxane	10.540	88	442701	9.73	ppbV		98
44) Methyl methacrylate	10.794	41	1526724	11.49	ppbV		99
45) n-Heptane	10.929	43	1977206	10.90	ppbV		100
46) cis-1,3-Dichloropropene	11.589	75	1447617	10.95	ppbV		100
47) Methyl isobutyl ketone	11.611	43	2752777	11.20	ppbV		99
48) trans-1,3-Dichloropropene	12.228	75	1305414	10.04	ppbV		99
49) 1,1,2-Trichloroethane	12.441	97	1005340	10.12	ppbV		90
50) Toluene	12.794	91	2771243	10.72	ppbV		98
51) Methyl n-butyl ketone	13.109	43	2511183	11.44	ppbV		99

AL SDG #E18-06141

Data Path : C:\DATA\05-18-18\
 Data File : aa7078icvss.D
 Acq On : 18 May 2018 3:14 pm
 Operator : jls
 Sample : 10 ppbv ICVSS
 Misc : CC483422
 ALS Vial : 8 Sample Multiplier: 1

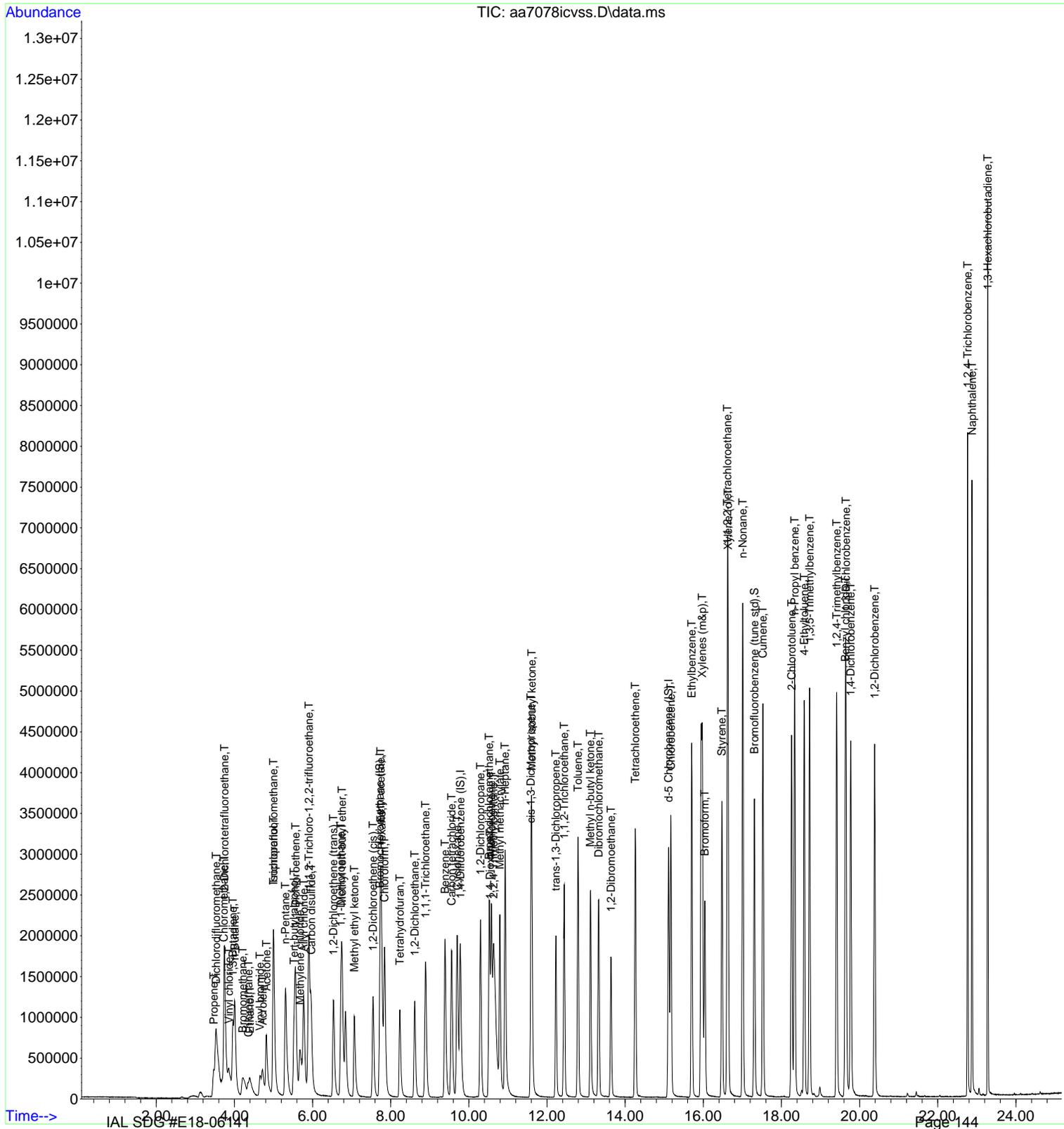
Quant Time: May 21 10:49:09 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 13:51:08 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Dibromochloromethane	13.325	129	1561436	10.38	ppbV	99
53) 1,2-Dibromoethane	13.637	107	1529144	10.34	ppbV	99
54) Tetrachloroethene	14.260	166	1076316	10.07	ppbV	97
56) Chlorobenzene	15.170	112	2087476	9.98	ppbV	99
57) Ethylbenzene	15.704	91	3845744	10.92	ppbV	96
58) Xylenes (m&p)	15.971	91	5995366	21.96	ppbV	96
59) Bromoform	16.039	173	1274264	10.70	ppbV	96
60) Styrene	16.476	104	2086221	11.56	ppbV	98
61) Xylene (o)	16.633	91	3185589	10.97	ppbV	96
62) 1,1,2,2-Tetrachloroethane	16.617	83	2560084	10.10	ppbV	99
63) n-Nonane	17.010	43	3246020	11.47	ppbV	100
65) Cumene	17.527	105	3969393	10.90	ppbV	97
66) 2-Chlorotoluene	18.260	91	3243514	10.86	ppbV	99
67) n-Propyl benzene	18.337	91	5484694	11.14	ppbV	95
68) 4-Ethyltoluene	18.585	105	3867722	11.31	ppbV	96
69) 1,3,5-Trimethylbenzene	18.720	105	3534194	11.30	ppbV	100
70) 1,2,4-Trimethylbenzene	19.415	105	3543577	11.60	ppbV	98
71) Benzyl chloride	19.633	91	2209493	10.51	ppbV	99
72) 1,3-Dichlorobenzene	19.649	146	2039322	10.36	ppbV	100
73) 1,4-Dichlorobenzene	19.775	146	2070619	10.39	ppbV	99
74) 1,2-Dichlorobenzene	20.386	146	2037844	10.48	ppbV	100
75) 1,2,4-Trichlorobenzene	22.765	180	1496366	9.91	ppbV	97
76) Naphthalene	22.878	127	558973	12.41	ppbV	100
77) 1,3-Hexachlorobutadiene	23.279	225	1144775	9.99	ppbV	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\05-18-18\
Data File : aa7078icvss.D
Acq On : 18 May 2018 3:14 pm
Operator : jls
Sample : 10 ppbv ICVSS
Misc : CC483422
ALS Vial : 8 Sample Multiplier: 1

Quant Time: May 21 10:49:09 2018
Quant Method : C:\msdchem\1\METHODS\0518.M
Quant Title : TO-15 on the Agilent 7890A / 5975C
QLast Update : Fri May 18 13:51:08 2018
Response via : Initial Calibration



Initial Calibration Verification Sample Standard

Lab Sample Name: 10 PPBV ICVSS
Spike Amount: 10 ppbv, except m&p-Xylenes at 20 ppbv
Amount of standard injected (ml): 50

Data File: AA7977ICVSS
Date Analyzed: 7/25/2018

Runs with this ICVSS:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7971BFB]	07/25/2018 8:35
40 PPBV STD [AA7972STD01]	07/25/2018 9:55
20 PPBV STD [AA7973STD02]	07/25/2018 10:29
10 PPBV STD [AA7974STD03]	07/25/2018 11:02
2 PPBV STD [AA7975STD04]	07/25/2018 12:18
0.2 PPBV STD [AA7976STD05]	07/25/2018 13:42
10 PPBV ICVSS [AA7977ICVSS]	07/25/2018 15:05

Compound	CAS #	Injected Amount (ppbv)	Recovered Amount (ppbv)	% Recovery	QC Limit
Acetone	67-64-1	11	11	100	
Acrolein	107-02-8	7.1	13	180	*
Allyl Chloride	107-05-1	11	11	100	
Benzene	71-43-2	10	10	100	
Benzyl chloride	100-44-7	11	11	100	
Bromodichloromethane	75-27-4	11	11	100	
Bromoform	75-25-2	11	12	110	
Bromomethane	74-83-9	11	10	91	
1,3-Butadiene	106-99-0	11	11	100	
n-Butane	106-97-8	11	11	100	
Chlorobenzene	108-90-7	10	11	110	
Chloroethane	75-00-3	11	10	91	
Chloroform	67-66-3	11	11	100	
Chloromethane	74-87-3	10	11	110	
Carbon disulfide	75-15-0	11	11	100	
Carbon tetrachloride	56-23-5	11	11	100	
2-Chlorotoluene	95-49-8	11	11	100	
Cumene	98-82-8	11	11	100	
Cyclohexane	110-82-7	11	11	100	
Dibromochloromethane	124-48-1	11	11	100	
1,2-Dibromoethane	106-93-4	10	10	100	
1,2-Dichlorobenzene	95-50-1	10	11	110	
1,3-Dichlorobenzene	541-73-1	10	11	110	
1,4-Dichlorobenzene	106-46-7	10	11	110	
Dichlorodifluoromethane	75-71-8	11	11	100	
1,1-Dichloroethane	75-34-3	11	11	100	
1,2-Dichloroethane	107-06-2	11	11	100	
1,1-Dichloroethene	75-35-4	11	11	100	
1,2-Dichloroethene (cis)	156-59-2	10	11	110	
1,2-Dichloroethene (trans)	156-60-5	11	11	100	
1,2-Dichloropropane	78-87-5	11	11	100	
1,3-Dichloropropene (cis)	10061-01-5	11	11	100	
1,3-Dichloropropene (trans)	10061-02-6	10	10	100	
1,2-Dichlorotetrafluoroethane	76-14-2	11	9.6	87	
1,4-Dioxane	123-91-1	8.3	9.9	120	
Ethanol	64-17-5	8.7	8.8	100	
Ethyl acetate	141-78-6	11	11	100	
Ethylbenzene	100-41-4	11	11	100	
4-Ethyltoluene	622-96-8	11	11	100	
n-Heptane	142-82-5	11	11	100	

ICVSS recovery must be within 70-130% of the spiked value for all compounds.

*** Values outside of QC limits**

Initial Calibration Verification Sample Standard

Lab Sample Name: 10 PPBV ICVSS
Spike Amount: 10 ppbv, except m&p-Xylenes at 20 ppbv
Amount of standard injected (ml): 50

Data File: AA7977ICVSS
Date Analyzed: 7/25/2018

Runs with this ICVSS:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7971BFB]	07/25/2018 8:35
40 PPBV STD [AA7972STD01]	07/25/2018 9:55
20 PPBV STD [AA7973STD02]	07/25/2018 10:29
10 PPBV STD [AA7974STD03]	07/25/2018 11:02
2 PPBV STD [AA7975STD04]	07/25/2018 12:18
0.2 PPBV STD [AA7976STD05]	07/25/2018 13:42
10 PPBV ICVSS [AA7977ICVSS]	07/25/2018 15:05

Compound	CAS #	Injected Amount (ppbv)	Recovered Amount (ppbv)	% Recovery	QC Limit
1,3-Hexachlorobutadiene	87-68-3	10	11	110	
n-Hexane	110-54-3	10	10	100	
Isopropanol	67-63-0	8.2	10	120	
Methylene chloride	75-09-2	11	10	91	
Methyl ethyl ketone	78-93-3	10	11	110	
Methyl isobutyl ketone	108-10-1	11	11	100	
Methyl methacrylate	80-62-6	11	11	100	
Methyl n-butyl ketone	591-78-6	9.4	11	120	
Methyl tert-butyl ether	1634-04-4	11	10	91	
Naphthalene	91-20-3	9.3	11	120	
n-Nonane	111-84-2	11	11	100	
n-Pentane	109-66-0	11	11	100	
Propene	115-07-1	11	12	110	
n-Propyl benzene	103-65-1	11	11	100	
Styrene	100-42-5	10	11	110	
Tert-butyl alcohol	75-65-0	8.0	9.2	120	
1,1,2,2-Tetrachloroethane	79-34-5	10	11	110	
Tetrachloroethene	127-18-4	10	10	100	
Tetrahydrofuran	109-99-9	11	12	110	
Toluene	108-88-3	10	10	100	
1,2,4-Trichlorobenzene	120-82-1	8.8	11	130	
1,1,1-Trichloroethane	71-55-6	11	11	100	
1,1,2-Trichloroethane	79-00-5	10	11	110	
Trichloroethene	79-01-6	10	11	110	
Trichlorofluoromethane	75-69-4	11	11	100	
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	11	10	91	
1,2,4-Trimethylbenzene	95-63-6	10	11	110	
1,3,5-Trimethylbenzene	108-67-8	11	11	100	
2,2,4-Trimethylpentane	540-84-1	11	11	100	
Vinyl bromide	593-60-2	9.7	11	110	
Vinyl chloride	75-01-4	11	11	100	
Xylenes (m&p)	179601-23-1	22	23	100	
Xylenes (o)	95-47-6	11	11	100	

ICVSS recovery must be within 70-130% of the spiked value for all compounds.

*** Values outside of QC limits**

Data Path : C:\DATA\07-25-18\
 Data File : aa79771cvss.D
 Acq On : 25 Jul 2018 3:05 pm
 Operator : jls
 Sample : 10 ppbv ICVSS
 Misc : CC483422
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Jul 25 15:36:45 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.686	130	434360	10.00	ppbV	0.00	
38) 1,4-Difluorobenzene (IS)	9.734	114	1814996	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.055	117	1540277	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...	17.248	95	1475009	10.50	ppbV	0.00	
Target Compounds							
							Qvalue
2) Propene	3.451	41	431460	11.63	ppbV	#	80
3) Dichlorodifluoromethane	3.519	85	1508906	10.79	ppbV		97
4) Chloromethane	3.689	52	190066	10.64	ppbV	#	1
5) 1,2-Dichlorotetrafluor...	3.727	85	1464207	9.60	ppbV		96
6) Vinyl chloride	3.846	62	664405	10.71	ppbV		86
7) 1,3-Butadiene	3.946	54	551716	11.01	ppbV		82
8) n-Butane	3.991	43	1033251	10.70	ppbV		97
9) Bromomethane	4.194	94	448735	10.17	ppbV		99
10) Chloroethane	4.338	64	328461	10.14	ppbV		93
11) Ethanol	4.371	45	299949	8.77	ppbV		98
12) Vinyl bromide	4.637	106	557257	11.11	ppbV		96
13) Acrolein	4.699	56	269452	12.60	ppbV		97
14) Acetone	4.792	58	364690	11.02	ppbV		79
15) Trichlorofluoromethane	4.972	101	1612224	10.79	ppbV		100
16) Isopropanol	4.969	45	1085784	10.07	ppbV		95
17) n-Pentane	5.277	43	1184109	10.88	ppbV		93
18) 1,1-Dichloroethene	5.535	61	1129497	11.11	ppbV		99
19) Methylene chloride	5.650	49	857233	10.30	ppbV	#	84
20) Tert-butyl alcohol	5.502	59	1281055	9.24	ppbV		100
21) Allyl chloride	5.740	76	299466	10.93	ppbV		100
22) 1,1,2-Trichloro-1,2,2-...	5.875	101	1239123	10.21	ppbV		89
23) Carbon disulfide	5.927	76	1893599	11.03	ppbV		92
24) 1,2-Dichloroethene (tr...	6.502	61	1026262	11.08	ppbV		92
25) 1,1-Dichloroethane	6.695	63	1316577	10.97	ppbV		99
26) Methyl tert-butyl ether	6.721	73	1802161	10.31	ppbV		100
27) Methyl ethyl ketone	7.036	43	1349894	11.02	ppbV	#	89
28) 1,2-Dichloroethene (cis)	7.509	61	996984	11.11	ppbV		91
29) Ethyl acetate	7.702	45	226087	11.09	ppbV		100
30) n-Hexane	7.731	57	1081737	10.29	ppbV		87
31) Chloroform	7.808	83	1402575	10.88	ppbV		98
32) Tetrahydrofuran	8.190	42	743740	11.99	ppbV		93
33) 1,2-Dichloroethane	8.573	62	1084031	10.99	ppbV		98
34) 1,1,1-Trichloroethane	8.849	97	1357207	10.56	ppbV		99
35) Benzene	9.345	78	1991345	10.35	ppbV		92
36) Carbon tetrachloride	9.515	117	1293160	10.71	ppbV		100
37) Cyclohexane	9.653	56	1160406	10.59	ppbV		100
39) 1,2-Dichloropropane	10.255	63	872412	10.54	ppbV		99
40) Bromodichloromethane	10.476	83	1514686	11.27	ppbV		98
41) 2,2,4-Trimethylpentane	10.592	57	3745203	11.21	ppbV		98
42) Trichloroethene	10.531	130	711375	10.78	ppbV		100
43) 1,4-Dioxane	10.496	88	372744	9.91	ppbV		81
44) Methyl methacrylate	10.746	41	954538	11.05	ppbV	#	78
45) n-Heptane	10.885	43	1279075	11.02	ppbV	#	82
46) cis-1,3-Dichloropropene	11.541	75	1246698	11.00	ppbV		99
47) Methyl isobutyl ketone	11.566	43	1715186	11.16	ppbV		90
48) trans-1,3-Dichloropropene	12.177	75	1174855	10.03	ppbV	#	81
49) 1,1,2-Trichloroethane	12.386	97	835731	10.50	ppbV		92
50) Toluene	12.740	91	2368634	10.40	ppbV		97
51) Methyl n-butyl ketone	13.065	43	1602395	11.10	ppbV		87

AL SDG #E18-06141

Data Path : C:\DATA\07-25-18\
 Data File : aa79771cvss.D
 Acq On : 25 Jul 2018 3:05 pm
 Operator : jls
 Sample : 10 ppbv ICVSS
 Misc : CC483422
 ALS Vial : 7 Sample Multiplier: 1

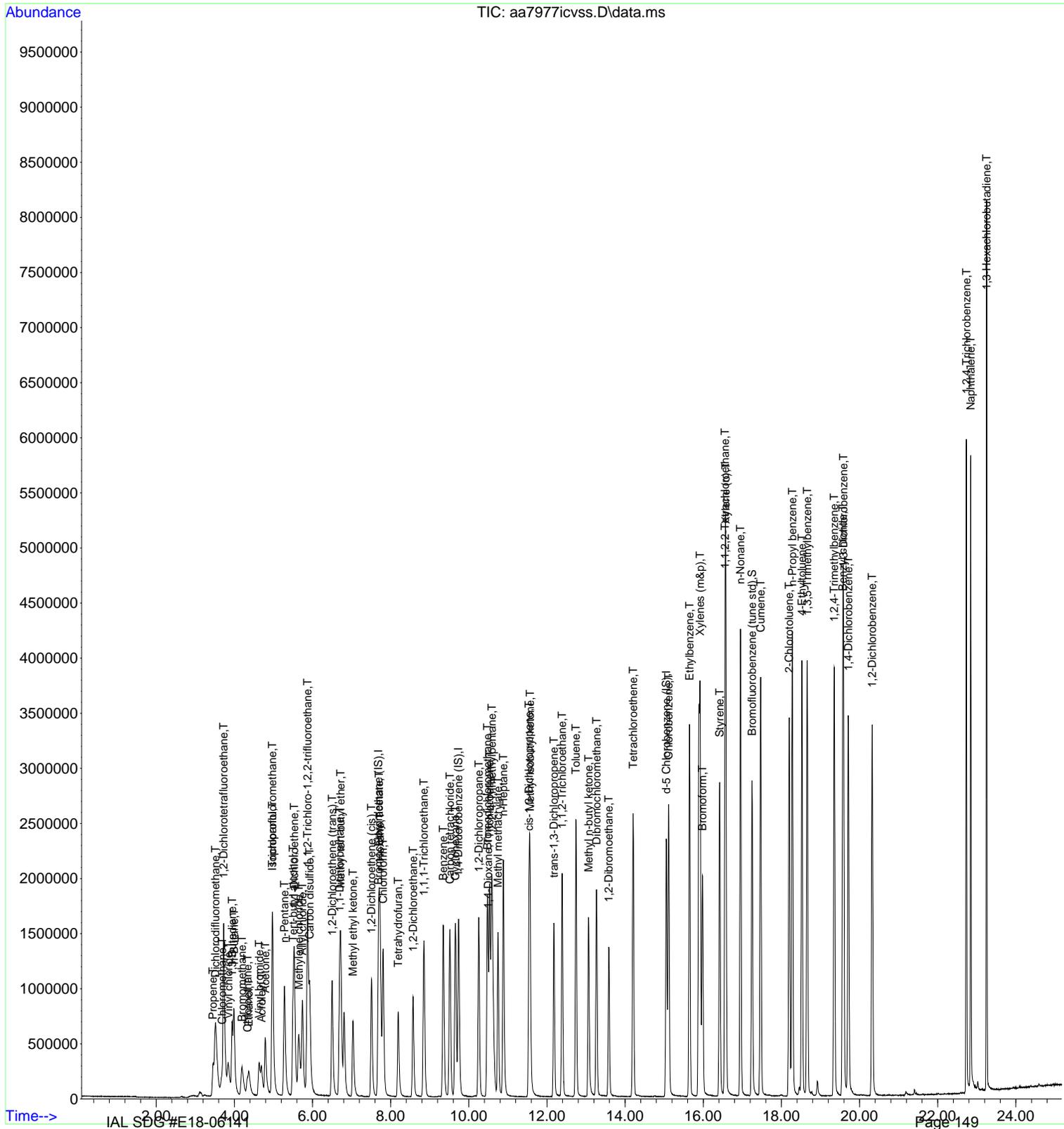
Quant Time: Jul 25 15:36:45 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Dibromochloromethane	13.267	129	1242245	11.08	ppbV	99
53) 1,2-Dibromoethane	13.582	107	1276654	10.45	ppbV	96
54) Tetrachloroethene	14.206	166	872215	10.48	ppbV	98
56) Chlorobenzene	15.113	112	1785159	10.65	ppbV	94
57) Ethylbenzene	15.650	91	3244054	10.68	ppbV	98
58) Xylenes (m&p)	15.917	91	5414325	22.97	ppbV	95
59) Bromoform	15.981	173	1100601	11.83	ppbV	97
60) Styrene	16.418	104	1821922	11.14	ppbV	97
61) Xylene (o)	16.576	91	2646162	11.06	ppbV	99
62) 1,1,2,2-Tetrachloroethane	16.557	83	2075885	10.83	ppbV	98
63) n-Nonane	16.952	43	2082795	11.49	ppbV	84
65) Cumene	17.470	105	3385196	11.17	ppbV	97
66) 2-Chlorotoluene	18.200	91	2749738	10.96	ppbV	97
67) n-Propyl benzene	18.280	91	4664264	11.36	ppbV	95
68) 4-Ethyltoluene	18.524	105	3300488	11.32	ppbV	96
69) 1,3,5-Trimethylbenzene	18.659	105	3031397	11.34	ppbV	99
70) 1,2,4-Trimethylbenzene	19.351	105	3029482	11.47	ppbV	98
71) Benzyl chloride	19.569	91	2678803	10.98	ppbV	97
72) 1,3-Dichlorobenzene	19.589	146	1694967	11.14	ppbV	98
73) 1,4-Dichlorobenzene	19.711	146	1763837	11.23	ppbV	98
74) 1,2-Dichlorobenzene	20.325	146	1706377	11.04	ppbV	99
75) 1,2,4-Trichlorobenzene	22.733	180	1288590	11.03	ppbV	99
76) Naphthalene	22.846	127	441780	11.44	ppbV	100
77) 1,3-Hexachlorobutadiene	23.251	225	917375	10.75	ppbV	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\07-25-18\
 Data File : aa7977icvss.D
 Acq On : 25 Jul 2018 3:05 pm
 Operator : jls
 Sample : 10 ppbv ICVSS
 Misc : CC483422
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Jul 25 15:36:45 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration



Continuing Calibration Data Summary Report

Initial Calibration Curve: 5/18/2018
 Instrument: AA
 Amount of standard injected (ml): 50

Date/Time of Calibration: 6/13/2018 09:18
 Sample ID: DCS
 Laboratory ID: AA7472DCVS

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7471BFB]	06/13/2018 08:30
10 PPBV DCVS [AA7472DCVS]	06/13/2018 09:18
METHOD BLANK [AA7473BLK]	06/13/2018 10:25
02 PPBV RLLCS [AA7474RLLCS]	06/13/2018 11:14
CLEAN CAN CERTIFICATION, BATCH MASTER 3059 [AA7475]	06/13/2018 12:20
10 PPBV CCCVS [AA7485CCCVS]	06/13/2018 20:47

Compound Name	Average RRF	Standard RRF	% Difference	PassFail	*
Acetone	0.81	0.88	-8.4	PASS	
Benzene	4.4	4.1	7.3	PASS	
Bromodichloromethane	0.82	0.88	-6.7	PASS	
Bromoform	0.62	0.70	-13	PASS	
Bromomethane	1.1	1.1	-0.70	PASS	
1,3-Butadiene	1.2	1.3	-4.5	PASS	
Chlorobenzene	1.1	1.1	2.6	PASS	
Chloroethane	0.82	0.81	0.90	PASS	
Chloroform	3.2	3.3	-5.0	PASS	
Chloromethane	0.49	0.52	-6.7	PASS	
Carbon disulfide	3.8	4.1	-8.4	PASS	
Carbon tetrachloride	3.0	3.0	-0.10	PASS	
Cyclohexane	2.6	2.6	0.50	PASS	
Dibromochloromethane	0.69	0.69	0.30	PASS	
1,2-Dibromoethane	0.68	0.64	4.6	PASS	
1,2-Dichlorobenzene	1.0	1.0	-2.1	PASS	
1,3-Dichlorobenzene	1.0	1.1	-3.2	PASS	
1,4-Dichlorobenzene	1.0	1.1	-1.3	PASS	
Dichlorodifluoromethane	3.5	3.7	-4.1	PASS	
1,1-Dichloroethane	3.0	3.1	-4.4	PASS	
1,2-Dichloroethane	2.7	2.7	-2.0	PASS	
1,1-Dichloroethene	2.6	2.8	-7.0	PASS	
1,2-Dichloroethene (cis)	2.2	2.2	1.9	PASS	
1,2-Dichloroethene (trans)	2.3	2.2	2.5	PASS	
1,2-Dichloropropane	0.49	0.47	3.5	PASS	
1,3-Dichloropropene (cis)	0.60	0.61	-0.50	PASS	
1,3-Dichloropropene (trans)	0.59	0.60	-0.50	PASS	
1,2-Dichlorotetrafluoroethane	3.7	3.7	1.0	PASS	
1,4-Dioxane	0.21	0.19	7.2	PASS	
Ethylbenzene	1.8	1.9	-1.8	PASS	
n-Heptane	0.83	0.93	-12	PASS	
1,3-Hexachlorobutadiene	0.60	0.63	-5.5	PASS	
n-Hexane	2.6	2.5	1.6	PASS	
Methylene chloride	2.3	2.4	-2.4	PASS	
Methyl ethyl ketone	3.7	3.9	-5.9	PASS	
Methyl isobutyl ketone	1.1	1.3	-13	PASS	
Methyl tert-butyl ether	3.3	3.2	2.7	PASS	
Styrene	0.94	0.98	-3.7	PASS	
Tert-butyl alcohol	2.3	2.5	-7.6	PASS	
1,1,2,2-Tetrachloroethane	1.3	1.4	-5.5	PASS	
Tetrachloroethene	0.49	0.43	12	PASS	
Toluene	1.2	1.1	6.2	PASS	
1,2,4-Trichlorobenzene	0.79	0.80	-0.80	PASS	
1,1,1-Trichloroethane	3.0	3.0	0.00	PASS	

*%Difference must be within +/- 30%

RRF - Relative Response Factor

Continuing Calibration Data Summary Report

Initial Calibration Curve: 5/18/2018
 Instrument: AA
 Amount of standard injected (ml): 50

Date/Time of Calibration: 6/13/2018 09:18
 Sample ID: DCS
 Laboratory ID: AA7472DCVS

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7471BFB]	06/13/2018 08:30
10 PPBV DCVS [AA7472DCVS]	06/13/2018 09:18
METHOD BLANK [AA7473BLK]	06/13/2018 10:25
02 PPBV RLLCS [AA7474RLLCS]	06/13/2018 11:14
CLEAN CAN CERTIFICATION, BATCH MASTER 3059 [AA7475]	06/13/2018 12:20
10 PPBV CCCVS [AA7485CCCVS]	06/13/2018 20:47

Compound Name	Average RRF	Standard RRF	% Difference	PassFail	*
1,1,2-Trichloroethane	0.45	0.44	4.0	PASS	
Trichloroethene	0.38	0.39	-1.0	PASS	
Trichlorofluoromethane	3.9	3.9	-1.2	PASS	
1,1,2-Trichloro-1,2,2-trifluoroethane	3.0	2.8	5.8	PASS	
1,2,4-Trimethylbenzene	1.6	1.8	-10	PASS	
1,3,5-Trimethylbenzene	1.6	1.8	-11	PASS	
2,2,4-Trimethylpentane	1.8	2.2	-22	PASS	
Vinyl bromide	1.0	1.0	-0.40	PASS	
Vinyl chloride	1.5	1.4	0.20	PASS	
Xylenes (m&p)	1.4	1.5	-7.6	PASS	
Xylenes (o)	1.5	1.6	-8.4	PASS	

*%Difference must be within +/- 30%

RRF - Relative Response Factor

Data Path : C:\DATA\06-13-18\
 Data File : aa7472dcvs.D
 Acq On : 13 Jun 2018 9:18 am
 Operator : jls
 Sample : 10 ppbv DCVS
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Jun 13 10:11:03 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 13:51:08 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 0% Max. R.T. Dev 0.40min
 Max. RRF Dev : 30% Max. Rel. Area : 500%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane (IS)	1.000	1.000	0.0	99	0.00
2 T	Propene	1.265	1.342	-6.1	108	0.00
3 T	Dichlorodifluoromethane	3.541	3.685	-4.1	105	0.00
4 T	Chloromethane	0.490	0.523	-6.7	111	-0.04
5 T	1,2-Dichlorotetrafluoroetha	3.727	3.688	1.0	109	0.00
6 T	Vinyl chloride	1.450	1.447	0.2	98	0.00
7 T	1,3-Butadiene	1.231	1.286	-4.5	98	0.00
8 T	n-Butane	2.846	3.149	-10.6	113	0.00
9 T	Bromomethane	1.061	1.068	-0.7	99	0.00
10 T	Chloroethane	0.820	0.813	0.9	104	0.00
11 T	Ethanol	0.792	0.747	5.7	107	0.00
12 T	Vinyl bromide	1.035	1.039	-0.4	95	0.00
13 T	Acrolein	0.510	0.550	-7.8	102	0.00
14 T	Acetone	0.808	0.876	-8.4	104	0.00
15 T	Trichlorofluoromethane	3.886	3.931	-1.2	106	0.00
16 T	Isopropanol	2.612	2.745	-5.1	99	0.00
17 T	n-Pentane	3.061	3.200	-4.5	103	0.00
18 T	1,1-Dichloroethene	2.569	2.750	-7.0	105	0.00
19 T	Methylene chloride	2.334	2.391	-2.4	112	0.00
20 T	Tert-butyl alcohol	2.339	2.516	-7.6	107	0.00
21 T	Allyl chloride	0.568	0.552	2.8	93	0.00
22 T	1,1,2-Trichloro-1,2,2-trifl	3.006	2.833	5.8	104	0.00
23 T	Carbon disulfide	3.772	4.088	-8.4	105	0.00
24 T	1,2-Dichloroethene (trans)	2.288	2.231	2.5	97	0.00
25 T	1,1-Dichloroethane	2.961	3.091	-4.4	107	0.00
26 T	Methyl tert-butyl ether	3.333	3.244	2.7	97	0.00
27 T	Methyl ethyl ketone	3.663	3.880	-5.9	104	0.00
28 T	1,2-Dichloroethene (cis)	2.201	2.159	1.9	96	0.00
29 T	Ethyl acetate	0.593	0.650	-9.6	106	0.00
30 T	n-Hexane	2.569	2.528	1.6	100	0.00
31 T	Chloroform	3.163	3.320	-5.0	107	0.00
32 T	Tetrahydrofuran	1.738	1.797	-3.4	100	0.00
33 T	1,2-Dichloroethane	2.652	2.705	-2.0	106	0.00
34 T	1,1,1-Trichloroethane	3.042	3.041	0.0	106	0.00
35 T	Benzene	4.381	4.059	7.3	95	0.00
36 T	Carbon tetrachloride	3.001	3.004	-0.1	107	0.00
37 T	Cyclohexane	2.604	2.591	0.5	99	0.00
38 I	1,4-Difluorobenzene (IS)	1.000	1.000	0.0	95	0.00
39 T	1,2-Dichloropropane	0.487	0.470	3.5	98	0.00
40 T	Bromodichloromethane	0.824	0.879	-6.7	106	0.00
41 T	2,2,4-Trimethylpentane	1.796	2.190	-21.9	110	0.00
42 T	Trichloroethene	0.384	0.388	-1.0	99	0.00
43 T	1,4-Dioxane	0.208	0.193	7.2	90	0.00
44 T	Methyl methacrylate	0.607	0.668	-10.0	101	0.00
45 T	n-Heptane	0.828	0.926	-11.8	107	0.00
46 T	cis-1,3-Dichloropropene	0.604	0.607	-0.5	95	0.00
47 T	Methyl isobutyl ketone	1.122	1.272	-13.4	106	0.00
48 T	trans-1,3-Dichloropropene	0.594	0.597	-0.5	94	0.00
49 T	1,1,2-Trichloroethane	0.454	0.436	4.0	99	0.00
50 T	Toluene	1.180	1.107	6.2	91	0.00
51 T	Methyl n-butyl ketone	1.002	1.108	-10.6	101	0.00
52 T	Dibromochloromethane	0.687	0.685	0.3	102	0.00
53 T	1,2-Dibromoethane	0.675	0.644	4.6	96	0.00

AL SDG #E18-06141

Data Path : C:\DATA\06-13-18\
 Data File : aa7472dcvs.D
 Acq On : 13 Jun 2018 9:18 am
 Operator : jls
 Sample : 10 ppbv DCVS
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Jun 13 10:11:03 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 13:51:08 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 0% Max. R.T. Dev 0.40min
 Max. RRF Dev : 30% Max. Rel. Area : 500%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
54 T	Tetrachloroethene	0.488	0.432	11.5	92	0.00
55 I	d-5 Chlorobenzene (IS)	1.000	1.000	0.0	90	0.00
56 T	Chlorobenzene	1.094	1.066	2.6	96	0.00
57 T	Ethylbenzene	1.841	1.874	-1.8	92	0.00
58 T	Xylenes (m&p)	1.427	1.536	-7.6	97	0.00
59 T	Bromoform	0.622	0.702	-12.9	106	0.00
60 T	Styrene	0.943	0.978	-3.7	88	0.00
61 T	Xylene (o)	1.518	1.646	-8.4	98	0.00
62 T	1,1,2,2-Tetrachloroethane	1.325	1.398	-5.5	103	0.00
63 T	n-Nonane	1.479	1.839	-24.3	108	0.00
64 S	Bromofluorobenzene (tune st	0.921	0.975	-5.9	95	0.00
65 T	Cumene	1.904	1.989	-4.5	95	0.00
66 T	2-Chlorotoluene	1.561	1.639	-5.0	96	0.00
67 T	n-Propyl benzene	2.572	2.849	-10.8	98	0.00
68 T	4-Ethyltoluene	1.788	1.938	-8.4	94	0.00
69 T	1,3,5-Trimethylbenzene	1.635	1.808	-10.6	96	0.00
70 T	1,2,4-Trimethylbenzene	1.597	1.762	-10.3	94	0.00
71 T	Benzyl chloride	1.099	1.601	-45.7#	124	0.00
72 T	1,3-Dichlorobenzene	1.029	1.062	-3.2	99	0.00
73 T	1,4-Dichlorobenzene	1.042	1.056	-1.3	96	0.00
74 T	1,2-Dichlorobenzene	1.016	1.037	-2.1	96	0.00
75 T	1,2,4-Trichlorobenzene	0.789	0.795	-0.8	94	0.00
76 T	Naphthalene	0.235	0.263	-11.9	94	0.00
77 T	1,3-Hexachlorobutadiene	0.599	0.632	-5.5	104	0.00

(#) = Out of Range

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

Data Path : C:\DATA\06-13-18\
 Data File : aa7472dcvs.D
 Acq On : 13 Jun 2018 9:18 am
 Operator : jls
 Sample : 10 ppbv DCVS
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Jun 13 10:11:03 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 13:51:08 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.724	130	539798	10.00	ppbV	0.00	
38) 1,4-Difluorobenzene (IS)	9.785	114	2150341	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.116	117	1773674	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...)	17.312	95	1728565	10.59	ppbV	0.00	
Target Compounds							
							Qvalue
2) Propene	3.473	41	796633	11.67	ppbV		84
3) Dichlorodifluoromethane	3.537	85	2187832	11.45	ppbV		98
4) Chloromethane	3.698	52	310310	11.72	ppbV #		1
5) 1,2-Dichlorotetrafluor...	3.746	85	1990852	9.90	ppbV		100
6) Vinyl chloride	3.868	62	858989	10.98	ppbV		88
7) 1,3-Butadiene	3.968	54	763581	11.49	ppbV		85
8) n-Butane	4.013	43	1869922	12.17	ppbV		90
9) Bromomethane	4.219	94	570583	9.96	ppbV		98
10) Chloroethane	4.364	64	438914	9.91	ppbV		93
11) Ethanol	4.399	45	379115	8.86	ppbV		99
12) Vinyl bromide	4.663	106	616932	11.05	ppbV		99
13) Acrolein	4.717	56	356484	12.95	ppbV		95
14) Acetone	4.817	58	520154	11.92	ppbV		81
15) Trichlorofluoromethane	5.000	101	2334372	11.13	ppbV		99
16) Isopropanol	5.000	45	1481834	10.51	ppbV #		83
17) n-Pentane	5.312	43	1900109	11.50	ppbV		99
18) 1,1-Dichloroethene	5.566	61	1633019	11.78	ppbV		88
19) Methylene chloride	5.679	49	1419869	11.27	ppbV		94
20) Tert-butyl alcohol	5.531	59	1629759	12.91	ppbV		100
21) Allyl chloride	5.778	76	327986	10.70	ppbV		100
22) 1,1,2-Trichloro-1,2,2-...	5.907	101	1681933	10.37	ppbV		85
23) Carbon disulfide	5.962	76	2427456	11.92	ppbV		96
24) 1,2-Dichloroethene (tr...	6.540	61	1324820	10.73	ppbV		95
25) 1,1-Dichloroethane	6.727	63	1835583	11.48	ppbV		99
26) Methyl tert-butyl ether	6.759	73	1926444	10.71	ppbV		94
27) Methyl ethyl ketone	7.074	43	2304085	11.65	ppbV		95
28) 1,2-Dichloroethene (cis)	7.550	61	1281740	10.79	ppbV		95
29) Ethyl acetate	7.743	45	385806	12.06	ppbV		100
30) n-Hexane	7.769	57	1501365	10.83	ppbV		87
31) Chloroform	7.846	83	1971358	11.55	ppbV		98
32) Tetrahydrofuran	8.235	42	1163814	12.41	ppbV #		85
33) 1,2-Dichloroethane	8.617	62	1605976	11.22	ppbV		100
34) 1,1,1-Trichloroethane	8.891	97	1805454	11.00	ppbV		99
35) Benzene	9.392	78	2409905	10.19	ppbV		96
36) Carbon tetrachloride	9.563	117	1783490	11.01	ppbV		100
37) Cyclohexane	9.707	56	1538354	10.95	ppbV		84
39) 1,2-Dichloropropane	10.302	63	1112181	10.63	ppbV		99
40) Bromodichloromethane	10.524	83	2079060	11.73	ppbV		99
41) 2,2,4-Trimethylpentane	10.637	57	5180610	13.41	ppbV		100
42) Trichloroethene	10.579	130	916905	11.10	ppbV		95
43) 1,4-Dioxane	10.543	88	415510	9.30	ppbV		93
44) Methyl methacrylate	10.794	41	1579870	12.11	ppbV		94
45) n-Heptane	10.932	43	2189277	12.29	ppbV		94
46) cis-1,3-Dichloropropene	11.592	75	1435942	11.06	ppbV		100
47) Methyl isobutyl ketone	11.614	43	3008902	12.47	ppbV		96
48) trans-1,3-Dichloropropene	12.231	75	1283942	10.06	ppbV		93
49) 1,1,2-Trichloroethane	12.444	97	1031155	10.57	ppbV		89
50) Toluene	12.797	91	2618304	10.32	ppbV		97
51) Methyl n-butyl ketone	13.116	43	2621325	12.16	ppbV		94

AL SDG #E18-06141

Data Path : C:\DATA\06-13-18\
 Data File : aa7472dcvs.D
 Acq On : 13 Jun 2018 9:18 am
 Operator : jls
 Sample : 10 ppbv DCVS
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

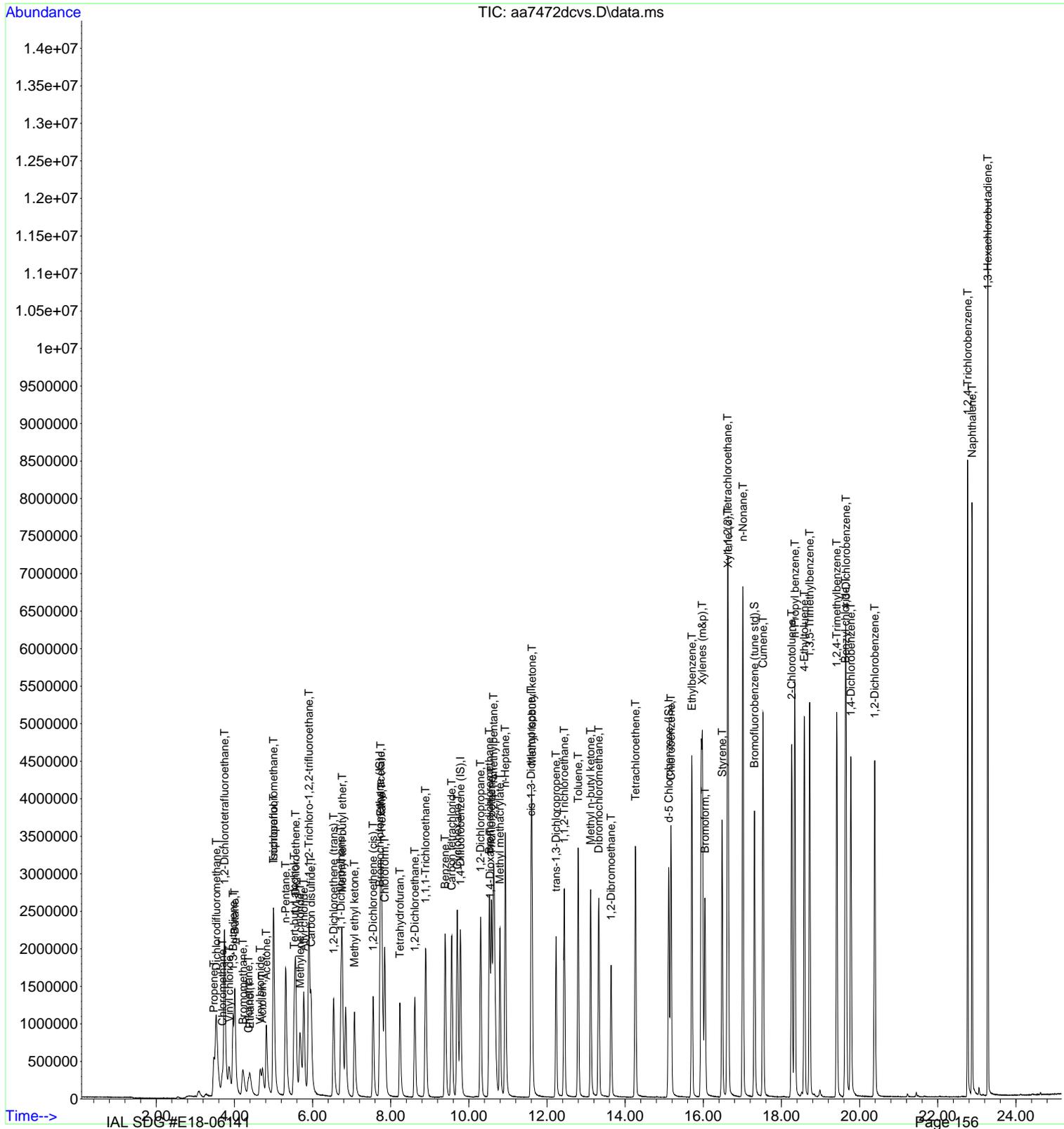
Quant Time: Jun 13 10:11:03 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 13:51:08 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Dibromochloromethane	13.325	129	1619280	10.97	ppbV	98
53) 1,2-Dibromoethane	13.640	107	1523377	10.49	ppbV	100
54) Tetrachloroethene	14.264	166	1020691	9.73	ppbV	98
56) Chlorobenzene	15.173	112	2079022	10.72	ppbV	96
57) Ethylbenzene	15.707	91	3655504	11.19	ppbV	95
58) Xylenes (m&p)	15.974	91	5994572	23.68	ppbV	95
59) Bromoform	16.042	173	1369647	12.41	ppbV	97
60) Styrene	16.482	104	1908242	11.40	ppbV	98
61) Xylene (o)	16.636	91	3212341	11.93	ppbV	95
62) 1,1,2,2-Tetrachloroethane	16.620	83	2728257	11.61	ppbV	99
63) n-Nonane	17.013	43	3588130	13.68	ppbV	96
65) Cumene	17.530	105	3881084	11.49	ppbV	97
66) 2-Chlorotoluene	18.263	91	3196887	11.55	ppbV	96
67) n-Propyl benzene	18.344	91	5557562	12.18	ppbV	94
68) 4-Ethyltoluene	18.585	105	3781802	11.93	ppbV	95
69) 1,3,5-Trimethylbenzene	18.723	105	3527242	12.16	ppbV	99
70) 1,2,4-Trimethylbenzene	19.414	105	3437135	12.13	ppbV	97
71) Benzyl chloride	19.633	91	2839391	14.57	ppbV	99
72) 1,3-Dichlorobenzene	19.652	146	2071672	11.35	ppbV	99
73) 1,4-Dichlorobenzene	19.778	146	2060344	11.15	ppbV	100
74) 1,2-Dichlorobenzene	20.385	146	2023280	11.22	ppbV	100
75) 1,2,4-Trichlorobenzene	22.765	180	1692425	12.09	ppbV	98
76) Naphthalene	22.877	127	559096	13.39	ppbV	100
77) 1,3-Hexachlorobutadiene	23.282	225	1233004	11.61	ppbV	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\06-13-18\
 Data File : aa7472dcvs.D
 Acq On : 13 Jun 2018 9:18 am
 Operator : jls
 Sample : 10 ppbv DCVS
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Jun 13 10:11:03 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 13:51:08 2018
 Response via : Initial Calibration



Continuing Calibration Data Summary Report

Initial Calibration Curve: 7/25/2018
 Instrument: AA
 Amount of standard injected (ml): 50

Date/Time of Calibration: 8/6/2018 13:07
 Sample ID: DCS
 Laboratory ID: AA8192DCVS

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA8191BFB]	08/06/2018 11:26
10 PPBV DCVS [AA8192DCVS]	08/06/2018 13:07
METHOD BLANK [AA8193BLK]	08/06/2018 13:42
02 PPBV RLLCS [AA8194RLLCS]	08/06/2018 14:23
E18-06141-07 [AA8206]	08/06/2018 22:28
E18-06141-08 [AA8207]	08/06/2018 23:02
E18-06141-09 [AA8208]	08/06/2018 23:35
10 PPBV CCCVS [AA8213CCCVS]	08/07/2018 02:22

Compound Name	Average RRF	Standard RRF	% Difference	PassFail	*
Acetone	0.76	0.87	-14	PASS	
Benzene	4.4	4.1	7.7	PASS	
Bromodichloromethane	0.74	0.86	-16	PASS	
Bromoform	0.60	0.72	-20	PASS	
Bromomethane	1.0	1.1	-4.2	PASS	
1,3-Butadiene	1.2	1.3	-16	PASS	
Chlorobenzene	1.1	1.0	4.4	PASS	
Chloroethane	0.75	0.80	-7.0	PASS	
Chloroform	3.0	3.3	-10	PASS	
Chloromethane	0.41	0.52	-26	PASS	
Carbon disulfide	4.0	4.1	-3.9	PASS	
Carbon tetrachloride	2.8	2.9	-4.0	PASS	
Cyclohexane	2.5	2.7	-5.4	PASS	
Dibromochloromethane	0.62	0.67	-7.9	PASS	
1,2-Dibromoethane	0.67	0.66	1.9	PASS	
1,2-Dichlorobenzene	1.0	1.1	-11	PASS	
1,3-Dichlorobenzene	0.99	1.1	-15	PASS	
1,4-Dichlorobenzene	1.0	1.2	-16	PASS	
Dichlorodifluoromethane	3.2	3.8	-18	PASS	
1,1-Dichloroethane	2.8	3.1	-12	PASS	
1,2-Dichloroethane	2.3	2.8	-25	PASS	
1,1-Dichloroethene	2.3	2.8	-20	PASS	
1,2-Dichloroethene (cis)	2.1	2.3	-12	PASS	
1,2-Dichloroethene (trans)	2.1	2.4	-12	PASS	
1,2-Dichloropropane	0.46	0.48	-5.5	PASS	
1,3-Dichloropropene (cis)	0.62	0.65	-4.0	PASS	
1,3-Dichloropropene (trans)	0.65	0.67	-4.2	PASS	
1,2-Dichlorotetrafluoroethane	3.5	3.6	-3.0	PASS	
1,4-Dioxane	0.21	0.20	3.4	PASS	
Ethylbenzene	2.0	1.9	3.7	PASS	
n-Heptane	0.64	0.68	-6.9	PASS	
1,3-Hexachlorobutadiene	0.55	0.64	-16	PASS	
n-Hexane	2.4	2.5	-5.2	PASS	
Methylene chloride	1.9	2.5	-29	PASS	
Methyl ethyl ketone	2.8	2.7	3.9	PASS	
Methyl isobutyl ketone	0.85	0.97	-14	PASS	
Methyl tert-butyl ether	4.0	3.5	12	PASS	
Styrene	1.1	1.0	3.9	PASS	
Tert-butyl alcohol	3.2	2.9	9.2	PASS	
1,1,2,2-Tetrachloroethane	1.2	1.4	-9.8	PASS	
Tetrachloroethene	0.46	0.41	12	PASS	
Toluene	1.3	1.2	7.5	PASS	

*%Difference must be within +/- 30%

RRF - Relative Response Factor

Continuing Calibration Data Summary Report

Initial Calibration Curve: 7/25/2018
Instrument: AA
Amount of standard injected (ml): 50

Date/Time of Calibration: 8/6/2018 13:07
Sample ID: DCS
Laboratory ID: AA8192DCVS

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA8191BFB]	08/06/2018 11:26
10 PPBV DCVS [AA8192DCVS]	08/06/2018 13:07
METHOD BLANK [AA8193BLK]	08/06/2018 13:42
02 PPBV RLLCS [AA8194RLLCS]	08/06/2018 14:23
E18-06141-07 [AA8206]	08/06/2018 22:28
E18-06141-08 [AA8207]	08/06/2018 23:02
E18-06141-09 [AA8208]	08/06/2018 23:35
10 PPBV CCCVS [AA8213CCCVS]	08/07/2018 02:22

Compound Name	Average RRF	Standard RRF	% Difference	PassFail	*
1,2,4-Trichlorobenzene	0.76	0.82	-8.2	PASS	
1,1,1-Trichloroethane	3.0	3.0	-2.9	PASS	
1,1,2-Trichloroethane	0.44	0.42	3.7	PASS	
Trichloroethene	0.36	0.36	1.6	PASS	
Trichlorofluoromethane	3.4	3.8	-11	PASS	
1,1,2-Trichloro-1,2,2-trifluoroethane	2.8	2.7	3.9	PASS	
1,2,4-Trimethylbenzene	1.7	1.9	-13	PASS	
1,3,5-Trimethylbenzene	1.7	2.0	-12	PASS	
2,2,4-Trimethylpentane	1.8	2.1	-16	PASS	
Vinyl bromide	1.2	1.1	3.1	PASS	
Vinyl chloride	1.4	1.5	-4.8	PASS	
Xylenes (m&p)	1.5	1.6	-3.6	PASS	
Xylenes (o)	1.6	1.7	-8.2	PASS	

*%Difference must be within +/- 30%

RRF - Relative Response Factor

IAL SDG #E18-06141

Data Path : C:\DATA\08-06-18\
 Data File : aa8192dcvs.D
 Acq On : 6 Aug 2018 1:07 pm
 Operator : jls
 Sample : 10 ppbv DCVS
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 06 16:03:05 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 0% Max. R.T. Dev 0.40min
 Max. RRF Dev : 30% Max. Rel. Area : 500%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane (IS)	1.000	1.000	0.0	90	0.00
2 T	Propene	0.854	0.631	26.1	63	0.00
3 T	Dichlorodifluoromethane	3.220	3.813	-18.4	107	0.00
4 T	Chloromethane	0.411	0.517	-25.8	113	-0.01
5 T	1,2-Dichlorotetrafluoroetha	3.511	3.615	-3.0	94	0.00
6 T	Vinyl chloride	1.428	1.496	-4.8	92	0.00
7 T	1,3-Butadiene	1.153	1.333	-15.6	100	0.00
8 T	n-Butane	2.224	2.248	-1.1	93	0.00
9 T	Bromomethane	1.016	1.059	-4.2	88	0.00
10 T	Chloroethane	0.746	0.798	-7.0	93	0.00
11 T	Ethanol	0.787	0.869	-10.4	111	0.00
12 T	Vinyl bromide	1.155	1.119	3.1	82	0.00
13 T	Acrolein	0.493	0.560	-13.6	94	0.00
14 T	Acetone	0.762	0.865	-13.5	100	0.00
15 T	Trichlorofluoromethane	3.441	3.831	-11.3	102	0.00
16 T	Isopropanol	2.483	3.187	-28.4	113	0.00
17 T	n-Pentane	2.506	1.764	29.6	64	0.00
18 T	1,1-Dichloroethene	2.341	2.816	-20.3	105	0.00
19 T	Methylene chloride	1.917	2.464	-28.5	127	0.00
20 T	Tert-butyl alcohol	3.191	2.898	9.2	83	0.00
21 T	Allyl chloride	0.631	0.606	4.0	82	0.00
22 T	1,1,2-Trichloro-1,2,2-trifl	2.794	2.685	3.9	91	0.00
23 T	Carbon disulfide	3.953	4.109	-3.9	90	0.00
24 T	1,2-Dichloroethene (trans)	2.133	2.393	-12.2	98	0.00
25 T	1,1-Dichloroethane	2.764	3.108	-12.4	100	0.00
26 T	Methyl tert-butyl ether	4.023	3.531	12.2	80	0.00
27 T	Methyl ethyl ketone	2.821	2.711	3.9	87	0.00
28 T	1,2-Dichloroethene (cis)	2.066	2.317	-12.1	98	0.00
29 T	Ethyl acetate	0.469	0.412	12.2	79	0.00
30 T	n-Hexane	2.421	2.548	-5.2	96	0.00
31 T	Chloroform	2.969	3.273	-10.2	97	0.00
32 T	Tetrahydrofuran	1.428	1.197	16.2	75	0.00
33 T	1,2-Dichloroethane	2.271	2.831	-24.7	112	0.00
34 T	1,1,1-Trichloroethane	2.959	3.045	-2.9	95	0.00
35 T	Benzene	4.429	4.089	7.7	83	0.00
36 T	Carbon tetrachloride	2.780	2.892	-4.0	96	0.00
37 T	Cyclohexane	2.523	2.660	-5.4	95	0.00
38 I	1,4-Difluorobenzene (IS)	1.000	1.000	0.0	85	0.00
39 T	1,2-Dichloropropane	0.456	0.481	-5.5	92	0.00
40 T	Bromodichloromethane	0.740	0.855	-15.5	96	0.00
41 T	2,2,4-Trimethylpentane	1.841	2.142	-16.3	96	0.00
42 T	Trichloroethene	0.364	0.358	1.6	83	0.00
43 T	1,4-Dioxane	0.207	0.200	3.4	79	0.00
44 T	Methyl methacrylate	0.476	0.576	-21.0	105	0.00
45 T	n-Heptane	0.639	0.683	-6.9	93	0.00
46 T	cis-1,3-Dichloropropene	0.624	0.649	-4.0	86	0.00
47 T	Methyl isobutyl ketone	0.847	0.965	-13.9	99	0.00
48 T	trans-1,3-Dichloropropene	0.645	0.672	-4.2	87	0.00
49 T	1,1,2-Trichloroethane	0.438	0.422	3.7	84	0.00
50 T	Toluene	1.255	1.161	7.5	80	0.00
51 T	Methyl n-butyl ketone	0.795	0.912	-14.7	98	0.00
52 T	Dibromochloromethane	0.618	0.667	-7.9	92	0.00
53 T	1,2-Dibromoethane	0.673	0.660	1.9	86	0.00

AL SDG #E18-06141

Data Path : C:\DATA\08-06-18\
 Data File : aa8192dcvs.D
 Acq On : 6 Aug 2018 1:07 pm
 Operator : jls
 Sample : 10 ppbv DCVS
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 06 16:03:05 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 0% Max. R.T. Dev 0.40min
 Max. RRF Dev : 30% Max. Rel. Area : 500%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
54 T	Tetrachloroethene	0.459	0.405	11.8	79	0.00
55 I	d-5 Chlorobenzene (IS)	1.000	1.000	0.0	83	0.00
56 T	Chlorobenzene	1.088	1.040	4.4	81	0.00
57 T	Ethylbenzene	1.972	1.900	3.7	82	0.00
58 T	Xylenes (m&p)	1.530	1.585	-3.6	83	0.00
59 T	Bromoform	0.604	0.724	-19.9	97	0.00
60 T	Styrene	1.062	1.021	3.9	78	0.00
61 T	Xylene (o)	1.554	1.682	-8.2	91	0.00
62 T	1,1,2,2-Tetrachloroethane	1.245	1.367	-9.8	94	0.00
63 T	n-Nonane	1.176	1.527	-29.8	109	0.00
64 S	Bromofluorobenzene (tune st	0.912	1.038	-13.8	95	0.00
65 T	Cumene	1.968	2.093	-6.4	88	0.00
66 T	2-Chlorotoluene	1.629	1.705	-4.7	88	0.00
67 T	n-Propyl benzene	2.666	3.085	-15.7	95	0.00
68 T	4-Ethyltoluene	1.893	2.090	-10.4	90	0.00
69 T	1,3,5-Trimethylbenzene	1.736	1.951	-12.4	92	0.00
70 T	1,2,4-Trimethylbenzene	1.715	1.938	-13.0	91	0.00
71 T	Benzyl chloride	1.584	1.934	-22.1	93	0.00
72 T	1,3-Dichlorobenzene	0.987	1.130	-14.5	97	0.00
73 T	1,4-Dichlorobenzene	1.020	1.178	-15.5	97	0.00
74 T	1,2-Dichlorobenzene	1.004	1.113	-10.9	94	0.00
75 T	1,2,4-Trichlorobenzene	0.759	0.821	-8.2	91	0.00
76 T	Naphthalene	0.251	0.284	-13.1	89	0.00
77 T	1,3-Hexachlorobutadiene	0.554	0.640	-15.5	101	0.00

(#) = Out of Range

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

Data Path : C:\DATA\08-06-18\
 Data File : aa8192dcvs.D
 Acq On : 6 Aug 2018 1:07 pm
 Operator : jls
 Sample : 10 ppbv DCVS
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 06 16:03:05 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.686	130	438428	10.00	ppbV	0.00	
38) 1,4-Difluorobenzene (IS)	9.734	114	1771164	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.049	117	1475100	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...)	17.242	95	1530975	11.38	ppbV	0.00	
Target Compounds							
							Qvalue
2) Propene	3.461	41	304366	8.13	ppbV		88
3) Dichlorodifluoromethane	3.519	85	1838935	13.03	ppbV		97
4) Chloromethane	3.679	52	249104	13.81	ppbV #		1
5) 1,2-Dichlorotetrafluor...	3.724	85	1584882	10.30	ppbV		99
6) Vinyl chloride	3.840	62	721657	11.52	ppbV		87
7) 1,3-Butadiene	3.943	54	643048	12.72	ppbV		92
8) n-Butane	3.988	43	1084038	11.12	ppbV		94
9) Bromomethane	4.187	94	459869	10.32	ppbV		99
10) Chloroethane	4.332	64	349689	10.69	ppbV		93
11) Ethanol	4.377	45	357948	10.37	ppbV		99
12) Vinyl bromide	4.634	106	539793	10.66	ppbV		97
13) Acrolein	4.692	56	294616	13.64	ppbV		93
14) Acetone	4.792	58	416930	12.49	ppbV		81
15) Trichlorofluoromethane	4.975	101	1847468	12.25	ppbV		100
16) Isopropanol	4.969	45	1397289	12.83	ppbV		94
17) n-Pentane	5.277	43	850703	7.74	ppbV		96
18) 1,1-Dichloroethene	5.531	61	1357985	13.23	ppbV		87
19) Methylene chloride	5.650	49	1188453	14.14	ppbV		92
20) Tert-butyl alcohol	5.502	59	1524721	10.90	ppbV		100
21) Allyl chloride	5.740	76	292135	10.57	ppbV		100
22) 1,1,2-Trichloro-1,2,2-...	5.872	101	1294943	10.57	ppbV		85
23) Carbon disulfide	5.924	76	1981881	11.44	ppbV #		90
24) 1,2-Dichloroethene (tr...	6.502	61	1154134	12.34	ppbV		94
25) 1,1-Dichloroethane	6.689	63	1499130	12.37	ppbV		99
26) Methyl tert-butyl ether	6.721	73	1703078	9.66	ppbV		94
27) Methyl ethyl ketone	7.033	43	1307380	10.57	ppbV		97
28) 1,2-Dichloroethene (cis)	7.509	61	1117186	12.33	ppbV		94
29) Ethyl acetate	7.695	45	198693	9.65	ppbV		100
30) n-Hexane	7.727	57	1228660	11.57	ppbV		89
31) Chloroform	7.805	83	1578630	12.13	ppbV		97
32) Tetrahydrofuran	8.194	42	629589	10.05	ppbV		88
33) 1,2-Dichloroethane	8.570	62	1365421	13.71	ppbV		100
34) 1,1,1-Trichloroethane	8.850	97	1468303	11.32	ppbV		99
35) Benzene	9.342	78	1972093	10.16	ppbV		95
36) Carbon tetrachloride	9.512	117	1394602	11.44	ppbV		99
37) Cyclohexane	9.657	56	1282816	11.60	ppbV		85
39) 1,2-Dichloropropane	10.255	63	936511	11.59	ppbV		100
40) Bromodichloromethane	10.470	83	1664975	12.70	ppbV		97
41) 2,2,4-Trimethylpentane	10.589	57	4173224	12.80	ppbV		100
42) Trichloroethene	10.528	130	697920	10.84	ppbV		99
43) 1,4-Dioxane	10.493	88	354021	9.64	ppbV		98
44) Methyl methacrylate	10.743	41	1122453	13.32	ppbV		91
45) n-Heptane	10.878	43	1330679	11.75	ppbV		94
46) cis-1,3-Dichloropropene	11.538	75	1263776	11.43	ppbV		98
47) Methyl isobutyl ketone	11.560	43	1880766	12.54	ppbV		96
48) trans-1,3-Dichloropropene	12.177	75	1189639	10.41	ppbV		91
49) 1,1,2-Trichloroethane	12.383	97	822438	10.59	ppbV		89
50) Toluene	12.740	91	2261627	10.18	ppbV		94
51) Methyl n-butyl ketone	13.058	43	1776612	12.62	ppbV		94

AL SDG #E18-06141

Data Path : C:\DATA\08-06-18\
 Data File : aa8192dcvs.D
 Acq On : 6 Aug 2018 1:07 pm
 Operator : jls
 Sample : 10 ppbv DCVS
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

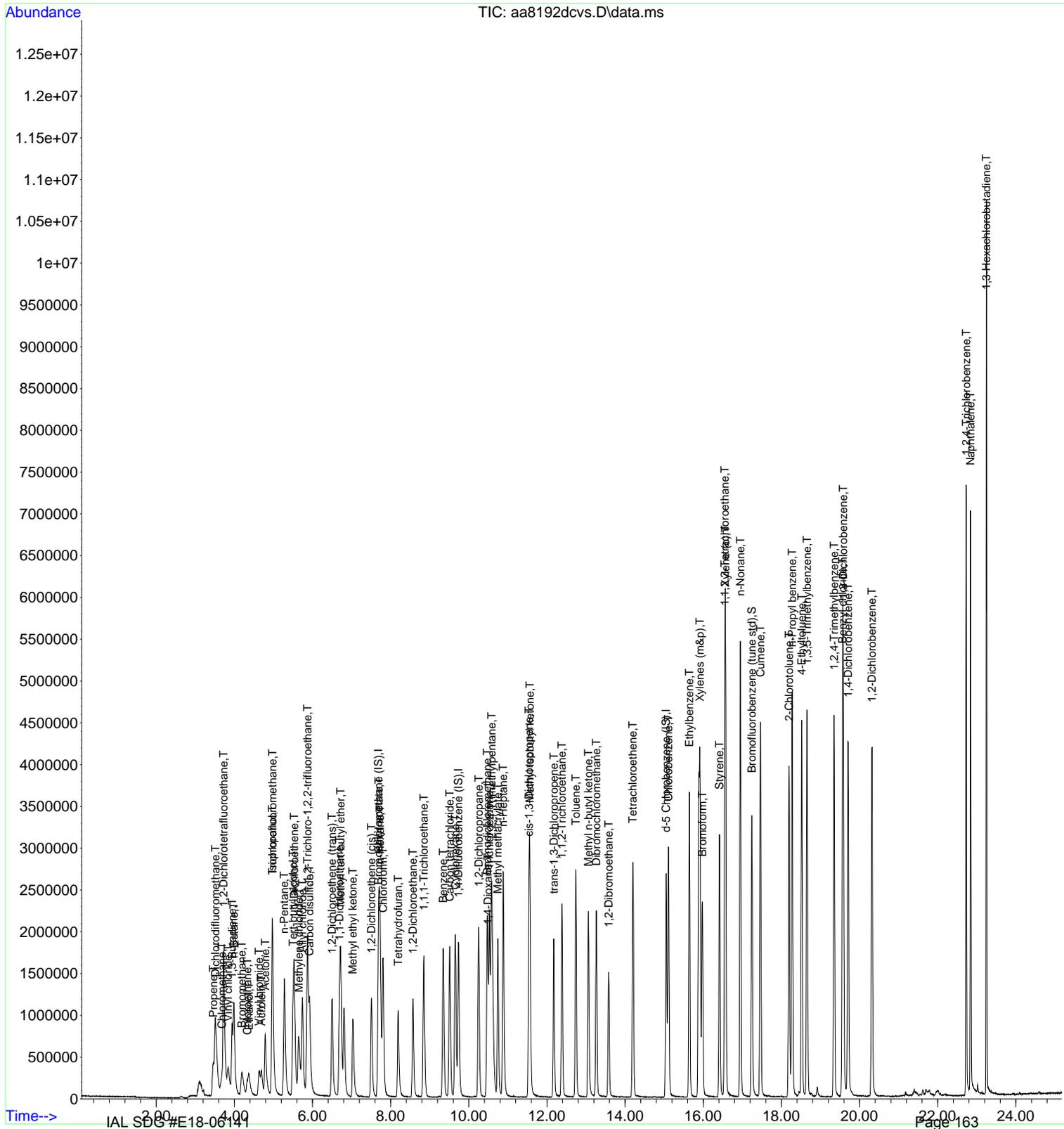
Quant Time: Aug 06 16:03:05 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Dibromochloromethane	13.264	129	1299265	11.88	ppbV	100
53) 1,2-Dibromoethane	13.579	107	1285505	10.79	ppbV	95
54) Tetrachloroethene	14.197	166	789922	9.73	ppbV	99
56) Chlorobenzene	15.110	112	1687277	10.51	ppbV	100
57) Ethylbenzene	15.644	91	3082982	10.60	ppbV	95
58) Xylenes (m&p)	15.910	91	5143044	22.79	ppbV	95
59) Bromoform	15.978	173	1173969	13.17	ppbV	97
60) Styrene	16.415	104	1657361	10.58	ppbV	98
61) Xylene (o)	16.570	91	2729707	11.91	ppbV	95
62) 1,1,2,2-Tetrachloroethane	16.553	83	2218126	12.08	ppbV	97
63) n-Nonane	16.949	43	2478302	14.28	ppbV	97
65) Cumene	17.463	105	3396814	11.70	ppbV	95
66) 2-Chlorotoluene	18.196	91	2766632	11.52	ppbV	94
67) n-Propyl benzene	18.274	91	5005472	12.73	ppbV	92
68) 4-Ethyltoluene	18.521	105	3391155	12.14	ppbV	95
69) 1,3,5-Trimethylbenzene	18.656	105	3165205	12.36	ppbV	98
70) 1,2,4-Trimethylbenzene	19.348	105	3144189	12.43	ppbV	95
71) Benzyl chloride	19.563	91	2852487	12.21	ppbV	100
72) 1,3-Dichlorobenzene	19.582	146	1834174	12.59	ppbV	98
73) 1,4-Dichlorobenzene	19.708	146	1911222	12.70	ppbV	98
74) 1,2-Dichlorobenzene	20.319	146	1805329	12.19	ppbV	99
75) 1,2,4-Trichlorobenzene	22.730	180	1453447	12.99	ppbV	99
76) Naphthalene	22.839	127	502670	13.60	ppbV	100
77) 1,3-Hexachlorobutadiene	23.248	225	1038802	12.71	ppbV	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\08-06-18\
 Data File : aa8192dcvs.D
 Acq On : 6 Aug 2018 1:07 pm
 Operator : jls
 Sample : 10 ppbv DCVS
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 06 16:03:05 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration



Continuing Calibration Data Summary Report

Initial Calibration Curve: 7/25/2018
 Instrument: AA
 Amount of standard injected (ml): 50

Date/Time of Calibration: 8/7/2018 10:17
 Sample ID: DCS
 Laboratory ID: AA8222DCVS

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA8221BFB]	08/07/2018 09:33
10 PPBV DCVS [AA8222DCVS]	08/07/2018 10:17
METHOD BLANK [AA8223BLK]	08/07/2018 11:01
02 PPBV RLLCS [AA8224RLLCS]	08/07/2018 11:47
E18-06141-08 [AA8232]	08/07/2018 16:57
E18-06141-09 [AA8233]	08/07/2018 17:31
10 PPBV CCCVS [AA8242CCCVS]	08/07/2018 22:31

Compound Name	Average RRF	Standard RRF	% Difference	PassFail	*
Acetone	0.76	0.81	-6.3	PASS	
Benzene	4.4	4.0	11	PASS	
Bromodichloromethane	0.74	0.87	-18	PASS	
Bromoform	0.60	0.72	-19	PASS	
Bromomethane	1.0	1.0	-1.9	PASS	
1,3-Butadiene	1.2	1.2	-2.9	PASS	
Chlorobenzene	1.1	1.1	0.70	PASS	
Chloroethane	0.75	0.75	-0.70	PASS	
Chloroform	3.0	3.2	-8.9	PASS	
Chloromethane	0.41	0.50	-22	PASS	
Carbon disulfide	4.0	4.1	-3.4	PASS	
Carbon tetrachloride	2.8	2.8	1.0	PASS	
Cyclohexane	2.5	2.5	1.1	PASS	
Dibromochloromethane	0.62	0.68	-10	PASS	
1,2-Dibromoethane	0.67	0.66	1.5	PASS	
1,2-Dichlorobenzene	1.0	1.1	-10	PASS	
1,3-Dichlorobenzene	0.99	1.1	-16	PASS	
1,4-Dichlorobenzene	1.0	1.2	-13	PASS	
Dichlorodifluoromethane	3.2	3.7	-15	PASS	
1,1-Dichloroethane	2.8	3.0	-8.7	PASS	
1,2-Dichloroethane	2.3	2.8	-23	PASS	
1,1-Dichloroethene	2.3	2.7	-16	PASS	
1,2-Dichloroethene (cis)	2.1	2.2	-4.2	PASS	
1,2-Dichloroethene (trans)	2.1	2.2	-3.6	PASS	
1,2-Dichloropropane	0.46	0.47	-3.1	PASS	
1,3-Dichloropropene (cis)	0.62	0.63	-1.1	PASS	
1,3-Dichloropropene (trans)	0.65	0.65	0.00	PASS	
1,2-Dichlorotetrafluoroethane	3.5	3.5	-0.10	PASS	
1,4-Dioxane	0.21	0.19	8.2	PASS	
Ethylbenzene	2.0	1.9	1.4	PASS	
n-Heptane	0.64	0.50	21	PASS	
1,3-Hexachlorobutadiene	0.55	0.65	-17	PASS	
n-Hexane	2.4	2.4	0.10	PASS	
Methylene chloride	1.9	2.4	-27	PASS	
Methyl ethyl ketone	2.8	3.2	-13	PASS	
Methyl isobutyl ketone	0.85	0.69	18	PASS	
Methyl tert-butyl ether	4.0	3.1	22	PASS	
Styrene	1.1	1.0	4.7	PASS	
Tert-butyl alcohol	3.2	2.5	22	PASS	
1,1,2,2-Tetrachloroethane	1.2	1.5	-18	PASS	
Tetrachloroethene	0.46	0.41	10	PASS	
Toluene	1.3	1.2	7.3	PASS	
1,2,4-Trichlorobenzene	0.76	0.80	-5.3	PASS	

*%Difference must be within +/- 30%

RRF - Relative Response Factor

Continuing Calibration Data Summary Report

Initial Calibration Curve: 7/25/2018
Instrument: AA
Amount of standard injected (ml): 50

Date/Time of Calibration: 8/7/2018 10:17
Sample ID: DCS
Laboratory ID: AA8222DCVS

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA8221BFB]	08/07/2018 09:33
10 PPBV DCVS [AA8222DCVS]	08/07/2018 10:17
METHOD BLANK [AA8223BLK]	08/07/2018 11:01
02 PPBV RLLCS [AA8224RLLCS]	08/07/2018 11:47
E18-06141-08 [AA8232]	08/07/2018 16:57
E18-06141-09 [AA8233]	08/07/2018 17:31
10 PPBV CCCVS [AA8242CCCVS]	08/07/2018 22:31

Compound Name	Average RRF	Standard RRF	% Difference	PassFail	*
1,1,1-Trichloroethane	3.0	2.9	1.9	PASS	
1,1,2-Trichloroethane	0.44	0.44	-0.90	PASS	
Trichloroethene	0.36	0.37	-1.9	PASS	
Trichlorofluoromethane	3.4	3.6	-5.3	PASS	
1,1,2-Trichloro-1,2,2-trifluoroethane	2.8	2.7	3.3	PASS	
1,2,4-Trimethylbenzene	1.7	1.9	-12	PASS	
1,3,5-Trimethylbenzene	1.7	2.0	-13	PASS	
2,2,4-Trimethylpentane	1.8	2.1	-15	PASS	
Vinyl bromide	1.2	1.0	12	PASS	
Vinyl chloride	1.4	1.4	4.4	PASS	
Xylenes (m&p)	1.5	1.6	-4.5	PASS	
Xylenes (o)	1.6	1.7	-11	PASS	

*%Difference must be within +/- 30%

RRF - Relative Response Factor

IAL SDG #E18-06141

Data Path : C:\DATA\08-07-18\
 Data File : aa8222dcvs.D
 Acq On : 7 Aug 2018 10:17 am
 Operator : jjw
 Sample : 10 ppbv DCVS
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 07 10:57:35 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 0% Max. R.T. Dev 0.40min
 Max. RRF Dev : 30% Max. Rel. Area : 500%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane (IS)	1.000	1.000	0.0	83	0.00
2 T	Propene	0.854	0.749	12.3	69	0.00
3 T	Dichlorodifluoromethane	3.220	3.694	-14.7	95	0.00
4 T	Chloromethane	0.411	0.502	-22.1	102	0.00
5 T	1,2-Dichlorotetrafluoroetha	3.511	3.515	-0.1	84	0.00
6 T	Vinyl chloride	1.428	1.365	4.4	77	0.00
7 T	1,3-Butadiene	1.153	1.187	-2.9	82	0.00
8 T	n-Butane	2.224	2.878	-29.4	110	0.00
9 T	Bromomethane	1.016	1.035	-1.9	79	0.00
10 T	Chloroethane	0.746	0.751	-0.7	81	0.00
11 T	Ethanol	0.787	0.727	7.6	86	0.00
12 T	Vinyl bromide	1.155	1.021	11.6	69	0.00
13 T	Acrolein	0.493	0.502	-1.8	78	0.00
14 T	Acetone	0.762	0.810	-6.3	87	0.00
15 T	Trichlorofluoromethane	3.441	3.625	-5.3	89	0.00
16 T	Isopropanol	2.483	2.772	-11.6	91	0.00
17 T	n-Pentane	2.506	3.025	-20.7	102	0.00
18 T	1,1-Dichloroethene	2.341	2.715	-16.0	93	0.00
19 T	Methylene chloride	1.917	2.427	-26.6	115	0.00
20 T	Tert-butyl alcohol	3.191	2.479	22.3	65	0.00
21 T	Allyl chloride	0.631	0.562	10.9	70	0.00
22 T	1,1,2-Trichloro-1,2,2-trifl	2.794	2.702	3.3	85	0.00
23 T	Carbon disulfide	3.953	4.086	-3.4	82	0.00
24 T	1,2-Dichloroethene (trans)	2.133	2.210	-3.6	83	0.00
25 T	1,1-Dichloroethane	2.764	3.004	-8.7	89	0.00
26 T	Methyl tert-butyl ether	4.023	3.122	22.4	65	0.00
27 T	Methyl ethyl ketone	2.821	3.198	-13.4	95	0.00
28 T	1,2-Dichloroethene (cis)	2.066	2.152	-4.2	84	0.00
29 T	Ethyl acetate	0.469	0.562	-19.8	99	0.00
30 T	n-Hexane	2.421	2.419	0.1	84	0.00
31 T	Chloroform	2.969	3.234	-8.9	89	0.00
32 T	Tetrahydrofuran	1.428	1.760	-23.2	101	0.00
33 T	1,2-Dichloroethane	2.271	2.792	-22.9	102	0.00
34 T	1,1,1-Trichloroethane	2.959	2.903	1.9	84	0.00
35 T	Benzene	4.429	3.963	10.5	74	0.00
36 T	Carbon tetrachloride	2.780	2.753	1.0	84	0.00
37 T	Cyclohexane	2.523	2.495	1.1	83	0.00
38 I	1,4-Difluorobenzene (IS)	1.000	1.000	0.0	76	0.00
39 T	1,2-Dichloropropane	0.456	0.470	-3.1	81	0.00
40 T	Bromodichloromethane	0.740	0.870	-17.6	88	0.00
41 T	2,2,4-Trimethylpentane	1.841	2.111	-14.7	85	0.00
42 T	Trichloroethene	0.364	0.371	-1.9	78	-0.01
43 T	1,4-Dioxane	0.207	0.190	8.2	67	0.00
44 T	Methyl methacrylate	0.476	0.416	12.6	68	0.00
45 T	n-Heptane	0.639	0.503	21.3	62	0.00
46 T	cis-1,3-Dichloropropene	0.624	0.631	-1.1	75	0.00
47 T	Methyl isobutyl ketone	0.847	0.694	18.1	64	0.00
48 T	trans-1,3-Dichloropropene	0.645	0.645	0.0	75	0.00
49 T	1,1,2-Trichloroethane	0.438	0.442	-0.9	79	0.00
50 T	Toluene	1.255	1.164	7.3	72	0.00
51 T	Methyl n-butyl ketone	0.795	0.704	11.4	68	0.00
52 T	Dibromochloromethane	0.618	0.680	-10.0	84	0.00
53 T	1,2-Dibromoethane	0.673	0.663	1.5	77	0.00

AL SDG #E18-06141

Data Path : C:\DATA\08-07-18\
 Data File : aa8222dcvs.D
 Acq On : 7 Aug 2018 10:17 am
 Operator : jjw
 Sample : 10 ppbv DCVS
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 07 10:57:35 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 0% Max. R.T. Dev 0.40min
 Max. RRF Dev : 30% Max. Rel. Area : 500%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
54 T	Tetrachloroethene	0.459	0.413	10.0	72	0.00
55 I	d-5 Chlorobenzene (IS)	1.000	1.000	0.0	74	0.00
56 T	Chlorobenzene	1.088	1.080	0.7	75	0.00
57 T	Ethylbenzene	1.972	1.945	1.4	74	0.00
58 T	Xylenes (m&p)	1.530	1.599	-4.5	74	0.00
59 T	Bromoform	0.604	0.717	-18.7	85	0.00
60 T	Styrene	1.062	1.012	4.7	68	0.00
61 T	Xylene (o)	1.554	1.722	-10.8	82	0.00
62 T	1,1,2,2-Tetrachloroethane	1.245	1.464	-17.6	89	0.00
63 T	n-Nonane	1.176	1.498	-27.4	95	0.00
64 S	Bromofluorobenzene (tune st	0.912	1.044	-14.5	85	0.00
65 T	Cumene	1.968	2.131	-8.3	79	0.00
66 T	2-Chlorotoluene	1.629	1.759	-8.0	80	0.00
67 T	n-Propyl benzene	2.666	3.122	-17.1	85	0.00
68 T	4-Ethyltoluene	1.893	2.105	-11.2	80	0.00
69 T	1,3,5-Trimethylbenzene	1.736	1.968	-13.4	82	0.00
70 T	1,2,4-Trimethylbenzene	1.715	1.918	-11.8	80	0.00
71 T	Benzyl chloride	1.584	1.852	-16.9	79	0.00
72 T	1,3-Dichlorobenzene	0.987	1.145	-16.0	86	0.00
73 T	1,4-Dichlorobenzene	1.020	1.151	-12.8	84	0.00
74 T	1,2-Dichlorobenzene	1.004	1.104	-10.0	82	0.00
75 T	1,2,4-Trichlorobenzene	0.759	0.799	-5.3	78	0.00
76 T	Naphthalene	0.251	0.272	-8.4	76	0.00
77 T	1,3-Hexachlorobutadiene	0.554	0.648	-17.0	91	0.00

(#) = Out of Range

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

Data Path : C:\DATA\08-07-18\
 Data File : aa8222dcvs.D
 Acq On : 7 Aug 2018 10:17 am
 Operator : jjw
 Sample : 10 ppbv DCVS
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 07 10:57:35 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.682	130	404488	10.00	ppbV	0.00	
38) 1,4-Difluorobenzene (IS)	9.734	114	1591121	10.00	ppbV	0.00	
55) d-5 Chlorobenzene (IS)	15.049	117	1304098	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...	17.245	95	1361833	11.45	ppbV	0.00	
Target Compounds							
							Qvalue
2) Propene	3.461	41	333397	9.65	ppbV		88
3) Dichlorodifluoromethane	3.522	85	1643634	12.62	ppbV		97
4) Chloromethane	3.699	52	223554	13.43	ppbV #		1
5) 1,2-Dichlorotetrafluor...	3.728	85	1421663	10.01	ppbV		97
6) Vinyl chloride	3.843	62	607411	10.51	ppbV		88
7) 1,3-Butadiene	3.946	54	528277	11.32	ppbV		82
8) n-Butane	3.988	43	1280649	14.24	ppbV #		87
9) Bromomethane	4.203	94	414307	10.08	ppbV		98
10) Chloroethane	4.332	64	303914	10.07	ppbV		91
11) Ethanol	4.374	45	276350	8.68	ppbV		97
12) Vinyl bromide	4.631	106	454469	9.73	ppbV		95
13) Acrolein	4.692	56	243851	12.24	ppbV		95
14) Acetone	4.792	58	360312	11.70	ppbV		73
15) Trichlorofluoromethane	4.975	101	1613029	11.59	ppbV		99
16) Isopropanol	4.972	45	1121332	11.16	ppbV #		87
17) n-Pentane	5.281	43	1346021	13.28	ppbV		98
18) 1,1-Dichloroethene	5.535	61	1207782	12.76	ppbV		87
19) Methylene chloride	5.644	49	1079715	13.93	ppbV		92
20) Tert-butyl alcohol	5.499	59	1203173	9.32	ppbV		100
21) Allyl chloride	5.744	76	250159	9.81	ppbV		100
22) 1,1,2-Trichloro-1,2,2-...	5.875	101	1202213	10.64	ppbV		83
23) Carbon disulfide	5.924	76	1818232	11.37	ppbV #		88
24) 1,2-Dichloroethene (tr...	6.499	61	983160	11.40	ppbV		97
25) 1,1-Dichloroethane	6.689	63	1336708	11.96	ppbV		99
26) Methyl tert-butyl ether	6.721	73	1389271	8.54	ppbV		95
27) Methyl ethyl ketone	7.033	43	1422716	12.47	ppbV		96
28) 1,2-Dichloroethene (cis)	7.509	61	957454	11.46	ppbV		97
29) Ethyl acetate	7.702	45	250199	13.18	ppbV		100
30) n-Hexane	7.724	57	1076376	10.99	ppbV		86
31) Chloroform	7.805	83	1439004	11.98	ppbV		98
32) Tetrahydrofuran	8.190	42	854272	14.79	ppbV #		83
33) 1,2-Dichloroethane	8.570	62	1242105	13.52	ppbV		99
34) 1,1,1-Trichloroethane	8.846	97	1291828	10.79	ppbV		99
35) Benzene	9.342	78	1763209	9.84	ppbV		95
36) Carbon tetrachloride	9.512	117	1225016	10.90	ppbV		100
37) Cyclohexane	9.657	56	1109929	10.88	ppbV		88
39) 1,2-Dichloropropane	10.248	63	823141	11.34	ppbV		98
40) Bromodichloromethane	10.473	83	1523040	12.93	ppbV		98
41) 2,2,4-Trimethylpentane	10.589	57	3695096	12.61	ppbV		99
42) Trichloroethene	10.525	130	649943	11.23	ppbV		97
43) 1,4-Dioxane	10.489	88	302644	9.18	ppbV		99
44) Methyl methacrylate	10.743	41	728507	9.62	ppbV		93
45) n-Heptane	10.878	43	880263	8.65	ppbV		93
46) cis-1,3-Dichloropropene	11.538	75	1104407	11.12	ppbV		99
47) Methyl isobutyl ketone	11.560	43	1215031	9.01	ppbV		93
48) trans-1,3-Dichloropropene	12.177	75	1025649	9.99	ppbV		87
49) 1,1,2-Trichloroethane	12.383	97	773565	11.09	ppbV		89
50) Toluene	12.737	91	2037850	10.21	ppbV		94
51) Methyl n-butyl ketone	13.058	43	1232417	9.74	ppbV		93

AL SDG #E18-06141

Data Path : C:\DATA\08-07-18\
 Data File : aa8222dcvs.D
 Acq On : 7 Aug 2018 10:17 am
 Operator : jjw
 Sample : 10 ppbv DCVS
 Misc : CC483586
 ALS Vial : 2 Sample Multiplier: 1

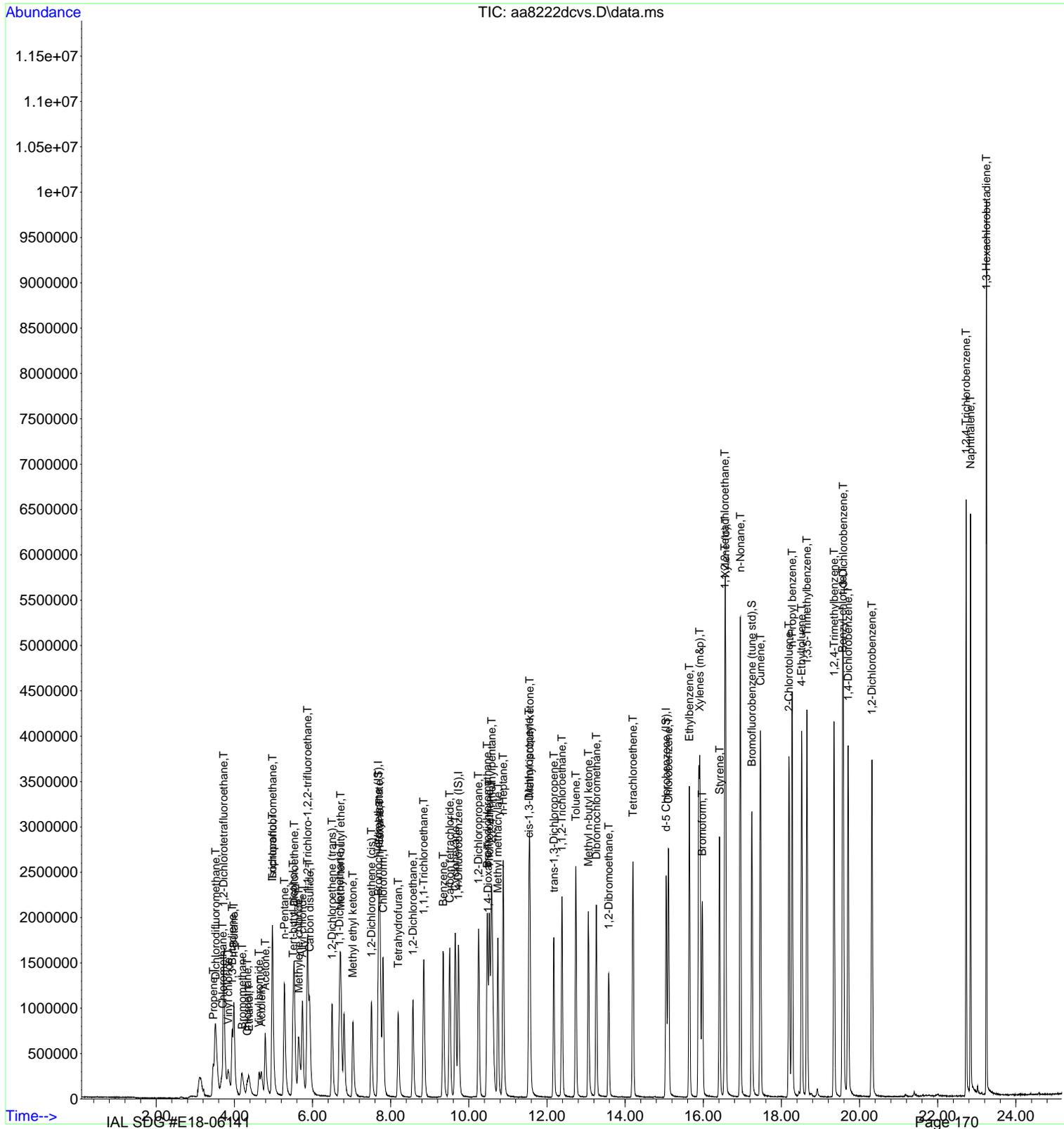
Quant Time: Aug 07 10:57:35 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Dibromochloromethane	13.264	129	1190543	12.12	ppbV	100
53) 1,2-Dibromoethane	13.579	107	1161108	10.85	ppbV	97
54) Tetrachloroethene	14.203	166	722436	9.90	ppbV	99
56) Chlorobenzene	15.110	112	1548795	10.91	ppbV	100
57) Ethylbenzene	15.644	91	2789576	10.85	ppbV	95
58) Xylenes (m&p)	15.907	91	4587138	22.99	ppbV	95
59) Bromoform	15.975	173	1029107	13.06	ppbV	97
60) Styrene	16.418	104	1451963	10.48	ppbV	99
61) Xylene (o)	16.570	91	2470278	12.19	ppbV	95
62) 1,1,2,2-Tetrachloroethane	16.553	83	2100434	12.94	ppbV	96
63) n-Nonane	16.949	43	2148325	14.00	ppbV	95
65) Cumene	17.463	105	3056880	11.91	ppbV	95
66) 2-Chlorotoluene	18.193	91	2523060	11.88	ppbV	95
67) n-Propyl benzene	18.274	91	4479204	12.89	ppbV	92
68) 4-Ethyltoluene	18.518	105	3019565	12.23	ppbV	94
69) 1,3,5-Trimethylbenzene	18.653	105	2823448	12.47	ppbV	97
70) 1,2,4-Trimethylbenzene	19.348	105	2750707	12.30	ppbV	95
71) Benzyl chloride	19.563	91	2414584	11.69	ppbV	99
72) 1,3-Dichlorobenzene	19.582	146	1641931	12.75	ppbV	98
73) 1,4-Dichlorobenzene	19.708	146	1651802	12.42	ppbV	98
74) 1,2-Dichlorobenzene	20.315	146	1583140	12.10	ppbV	98
75) 1,2,4-Trichlorobenzene	22.727	180	1250631	12.64	ppbV	100
76) Naphthalene	22.839	127	425537	13.02	ppbV	100
77) 1,3-Hexachlorobutadiene	23.248	225	929653	12.87	ppbV	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\08-07-18\
Data File : aa8222dcvs.D
Acq On : 7 Aug 2018 10:17 am
Operator : jjw
Sample : 10 ppbv DCVS
Misc : CC483586
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 07 10:57:35 2018
Quant Method : C:\msdchem\1\METHODS\0725.M
Quant Title : TO-15 on the Agilent 7890A / 5975C
QLast Update : Wed Jul 25 14:15:57 2018
Response via : Initial Calibration



Section VIII: Raw Quality Control Data Package

BFB Tune Spectra

Method Blank

Laboratory Sample Duplicate

Instrument Run Logs

Pressure Gauge Readings (initial and final)

Example Calculations

Screening Data

Clean Canister Certification

Data Path: C:\DATA\05-18-18\
Data File: AA7071BFB.D
Acq On: 5/18/2018 8:17:00AM
Operator: jls
Sample: BFB
Misc: ALM029426
ALS Vial: 1 **Multiplier:** 1
Integration File: rteint.p
Method: C:\msdchem\1\METHODS\0511.M
Last Update: Mon May 14 11:16:39 2018
Spectrum Information:

Pass/Fail	Target Mass	Rel. to Mass	Lower Limit %	Higher Limit %	Raw Abundance	% Relative Abundance
PASS	50	95	8	40	183818	27.3
PASS	75	95	30	66	378859	56.2
PASS	95	95	100	100	674347	100.0
PASS	96	95	5	9	45051	6.7
PASS	173	174	0.00	2	2912	0.8
PASS	174	95	50	100	387670	57.5
PASS	175	174	4	9	26268	6.8
PASS	176	174	93	101	373555	96.4
PASS	177	176	5	9	24192	6.5

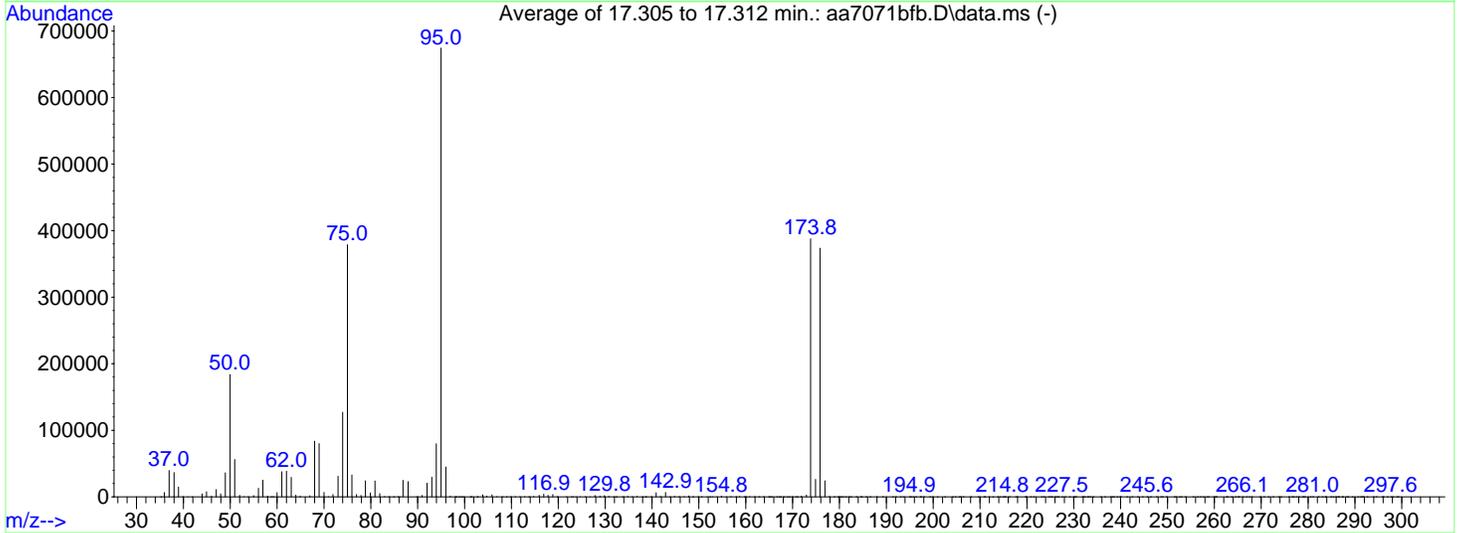
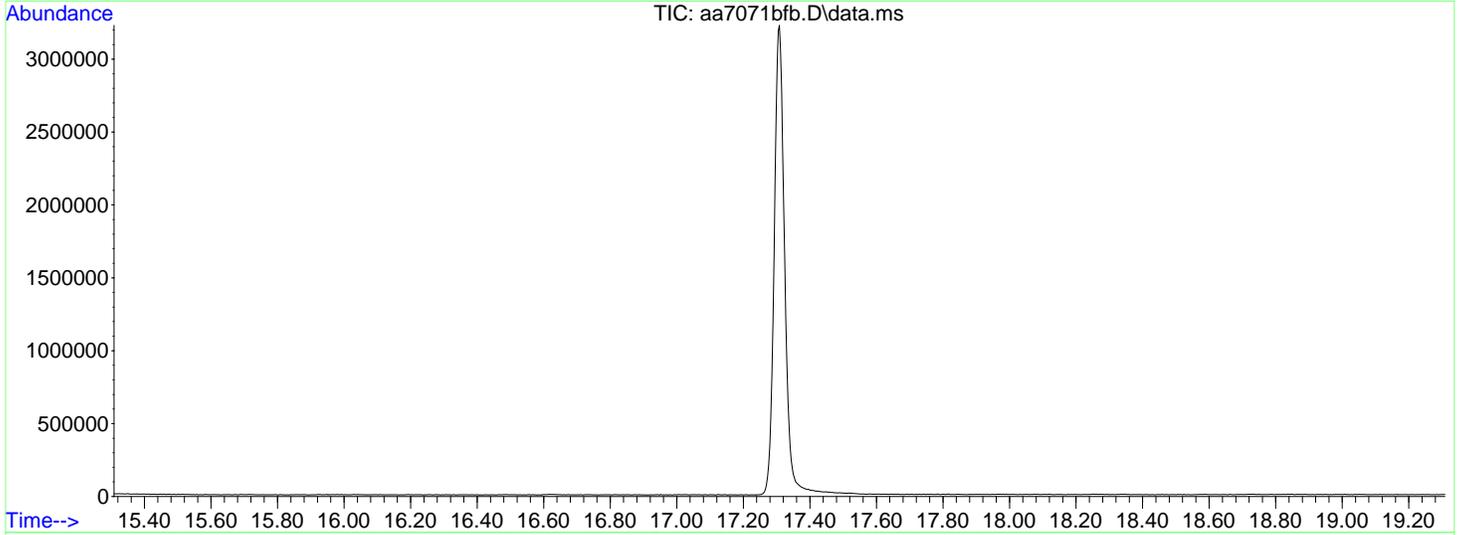
Runs with this BFB:

Lab Sample Number	Date File	Field Sample	Date/Time of Sample/Standard Analysis
BFB	AA7071BFB	NA	5/18/2018 8:17:00 AM
40 PPBV STD	AA7073STD01	NA	5/18/2018 9:46:00 AM
20 PPBV STD	AA7074STD02	NA	5/18/2018 10:19:00 AM
10 PPBV STD	AA7075STD03	NA	5/18/2018 10:53:00 AM
2 PPBV STD	AA7076STD04	NA	5/18/2018 12:20:00 PM
0.2 PPBV STD	AA7077STD05	NA	5/18/2018 1:03:00 PM
10 PPBV ICVSS	AA7078ICVSS	NA	5/18/2018 3:14:00 PM

Data Path : C:\DATA\05-18-18\
 Data File : aa7071bfb.D
 Acq On : 18 May 2018 8:17 am
 Operator : jls
 Sample : BFB
 Misc : ALM029426
 ALS Vial : 1 Sample Multiplier: 1

Integration File: rteint.p

Method : C:\msdchem\1\METHODS\0518.M
 Title : TO-15 on the Agilent 7890A / 5975C
 Last Update : Fri May 18 13:51:08 2018



AutoFind: Scans 5355, 5356, 5357; Background Corrected with Scan 5332

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	8	40	27.3	183818	PASS
75	95	30	66	56.2	378859	PASS
95	95	100	100	100.0	674347	PASS
96	95	5	9	6.7	45051	PASS
173	174	0.00	2	0.8	2912	PASS
174	95	50	100	57.5	387670	PASS
175	174	4	9	6.8	26268	PASS
176	174	93	101	96.4	373555	PASS
177	176	5	9	6.5	24192	PASS

Data Path: C:\DATA\06-13-18\
Data File: AA7471BFB.D
Acq On: 6/13/2018 8:30:00AM
Operator: jls
Sample: BFB
Misc: ALM029426
ALS Vial: 1 **Multiplier:** 1
Integration File: rteint.p
Method: C:\msdchem\1\METHODS\0518.M
Last Update: Fri May 18 13:51:08 2018
Spectrum Information:

PassFail	Target Mass	Rel. to Mass	Lower Limit %	Higher Limit %	Raw Abundance	% Relative Abundance
PASS	50	95	8	40	200102	30.6
PASS	75	95	30	66	384246	58.8
PASS	95	95	100	100	653589	100.0
PASS	96	95	5	9	42346	6.5
PASS	173	174	0.00	2	3805	1.0
PASS	174	95	50	100	363368	55.6
PASS	175	174	4	9	26616	7.3
PASS	176	174	93	101	348380	95.9
PASS	177	176	5	9	23261	6.7

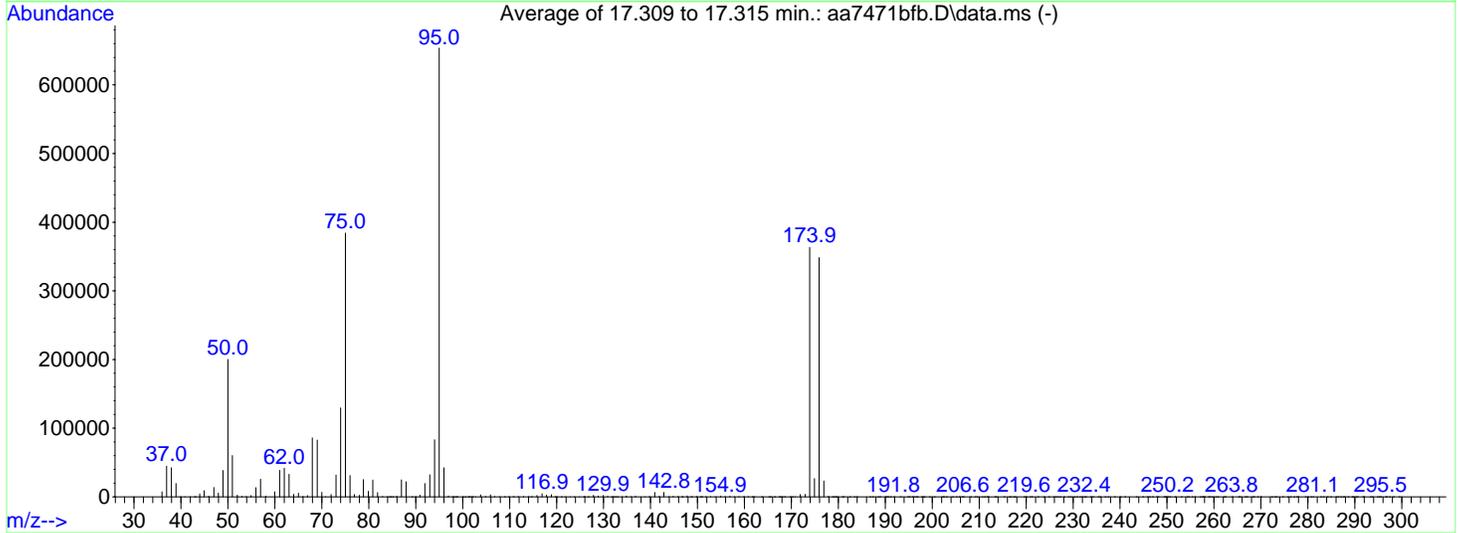
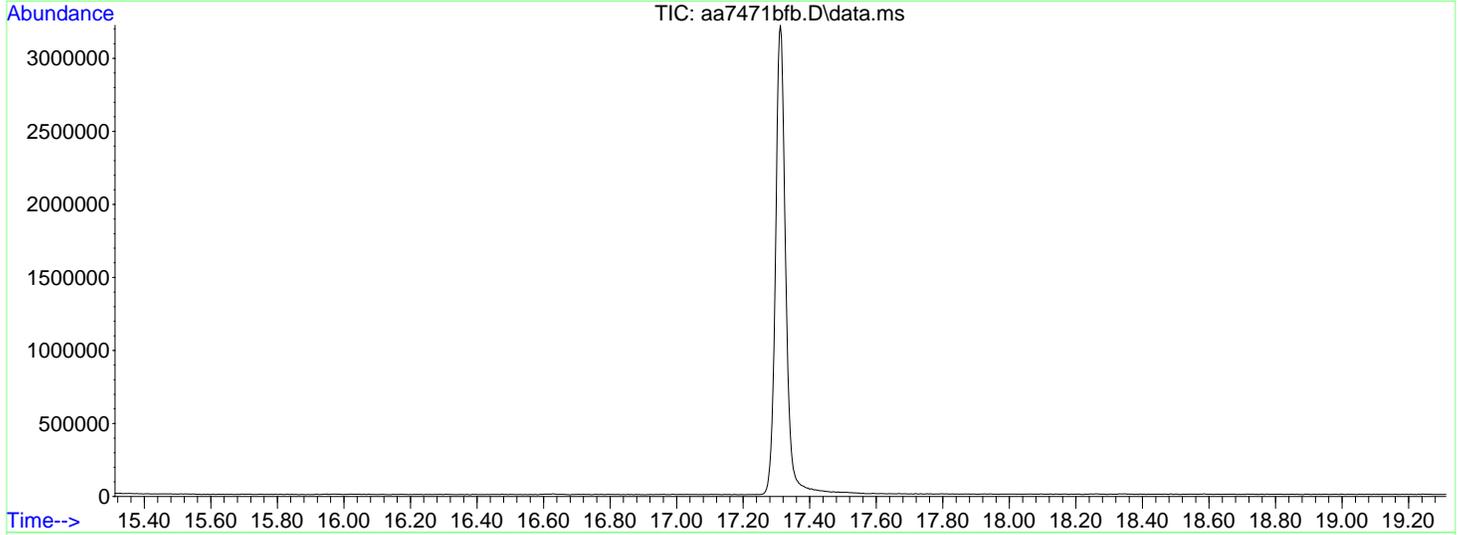
Runs with this BFB:

Lab Sample Number	Date File	Field Sample	Date/Time of Sample/Standard Analysis
BFB	AA7471BFB	NA	6/13/2018 8:30:00 AM
10 PPBV DCVS	AA7472DCVS	NA	6/13/2018 9:18:00 AM
METHOD BLANK	AA7473BLK	NA	6/13/2018 10:25:00 AM
02 PPBV RLLCS	AA7474RLLCS	NA	6/13/2018 11:14:00 AM
3059	AA7475	NA	6/13/2018 12:20:00 PM
10 PPBV CCCVS	AA7485CCCVS	NA	6/13/2018 8:47:00 PM

Data Path : C:\DATA\06-13-18\
 Data File : aa7471bfb.D
 Acq On : 13 Jun 2018 8:30 am
 Operator : jls
 Sample : BFB
 Misc : ALM029426
 ALS Vial : 1 Sample Multiplier: 1

Integration File: rteint.p

Method : C:\msdchem\1\METHODS\0518.M
 Title : TO-15 on the Agilent 7890A / 5975C
 Last Update : Fri May 18 13:51:08 2018



AutoFind: Scans 5356, 5357, 5358; Background Corrected with Scan 5334

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	8	40	30.6	200102	PASS
75	95	30	66	58.8	384246	PASS
95	95	100	100	100.0	653589	PASS
96	95	5	9	6.5	42346	PASS
173	174	0.00	2	1.0	3805	PASS
174	95	50	100	55.6	363368	PASS
175	174	4	9	7.3	26616	PASS
176	174	93	101	95.9	348380	PASS
177	176	5	9	6.7	23261	PASS

Data Path: C:\DATA\07-25-18\
Data File: AA7971BFB.D
Acq On: 7/25/2018 8:35:00AM
Operator: jls
Sample: BFB
Misc: ALM029426
ALS Vial: 1 **Multiplier:** 1
Integration File: rteint.p
Method: C:\msdchem\1\METHODS\0518.M
Last Update: Fri May 18 13:51:08 2018

Spectrum Information:

Pass/Fail	Target Mass	Rel. to Mass	Lower Limit %	Higher Limit %	Raw Abundance	% Relative Abundance
PASS	50	95	8	40	156657	20.1
PASS	75	95	30	66	381141	49.0
PASS	95	95	100	100	778176	100.0
PASS	96	95	5	9	52476	6.7
PASS	173	174	0.00	2	3085	0.7
PASS	174	95	50	100	459712	59.1
PASS	175	174	4	9	31097	6.8
PASS	176	174	93	101	438861	95.5
PASS	177	176	5	9	27982	6.4

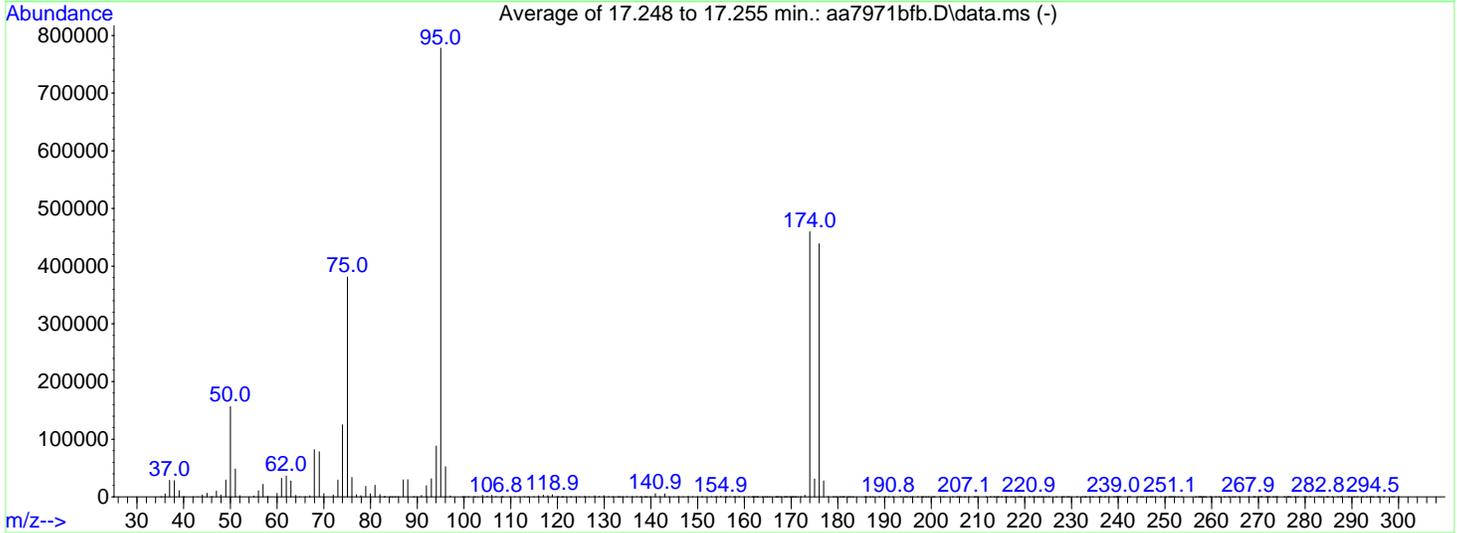
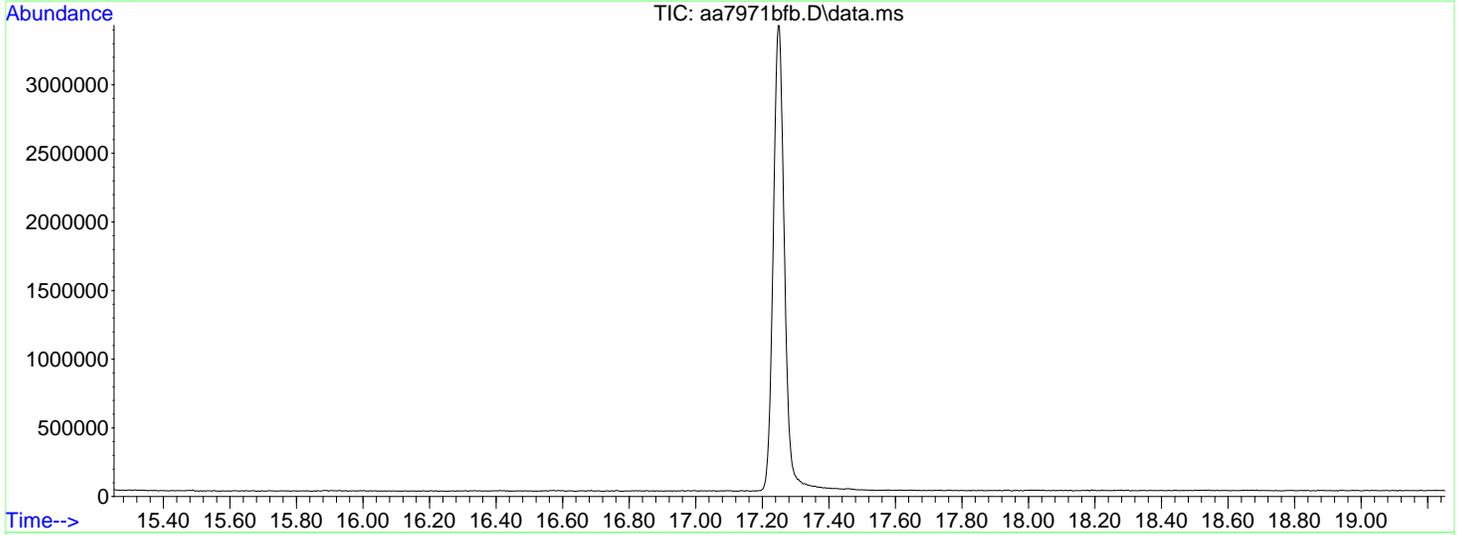
Runs with this BFB:

Lab Sample Number	Date File	Field Sample	Date/Time of Sample/Standard Analysis
BFB	AA7971BFB	NA	7/25/2018 8:35:00 AM
40 PPBV STD	AA7972STD01	NA	7/25/2018 9:55:00 AM
20 PPBV STD	AA7973STD02	NA	7/25/2018 10:29:00 AM
10 PPBV STD	AA7974STD03	NA	7/25/2018 11:02:00 AM
2 PPBV STD	AA7975STD04	NA	7/25/2018 12:18:00 PM
0.2 PPBV STD	AA7976STD05	NA	7/25/2018 1:42:00 PM
10 PPBV ICVSS	AA7977ICVSS	NA	7/25/2018 3:05:00 PM

Data Path : C:\DATA\07-25-18\
 Data File : aa7971bfb.D
 Acq On : 25 Jul 2018 8:35 am
 Operator : jls
 Sample : BFB
 Misc : ALM029426
 ALS Vial : 1 Sample Multiplier: 1

Integration File: rteint.p

Method : C:\msdchem\1\METHODS\0725.M
 Title : TO-15 on the Agilent 7890A / 5975C
 Last Update : Wed Jul 25 14:15:57 2018



AutoFind: Scans 5337, 5338, 5339; Background Corrected with Scan 5316

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	8	40	20.1	156657	PASS
75	95	30	66	49.0	381141	PASS
95	95	100	100	100.0	778176	PASS
96	95	5	9	6.7	52476	PASS
173	174	0.00	2	0.7	3085	PASS
174	95	50	100	59.1	459712	PASS
175	174	4	9	6.8	31097	PASS
176	174	93	101	95.5	438861	PASS
177	176	5	9	6.4	27982	PASS

Data Path: C:\DATA\08-06-18\
Data File: AA8191BFB.D
Acq On: 8/6/2018 11:26:00AM
Operator: jls
Sample: BFB
Misc: ALM029426
ALS Vial: 1 **Multiplier:** 1
Integration File: rteint.p
Method: C:\msdchem\1\METHODS\0725.M
Last Update: Wed Jul 25 14:15:57 2018

Spectrum Information:

PassFail	Target Mass	Rel. to Mass	Lower Limit %	Higher Limit %	Raw Abundance	% Relative Abundance
PASS	50	95	8	40	177003	32.6
PASS	75	95	30	66	334482	61.6
PASS	95	95	100	100	542971	100.0
PASS	96	95	5	9	36432	6.7
PASS	173	174	0.00	2	2635	0.9
PASS	174	95	50	100	286464	52.8
PASS	175	174	4	9	21901	7.6
PASS	176	174	93	101	278015	97.1
PASS	177	176	5	9	17336	6.2

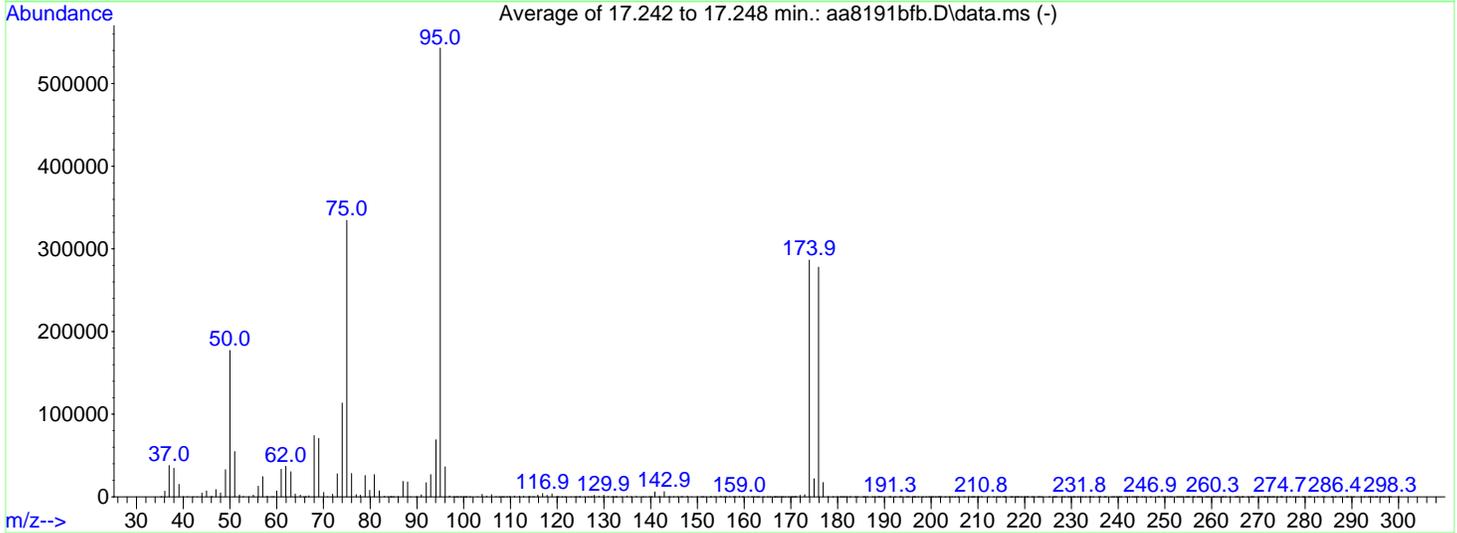
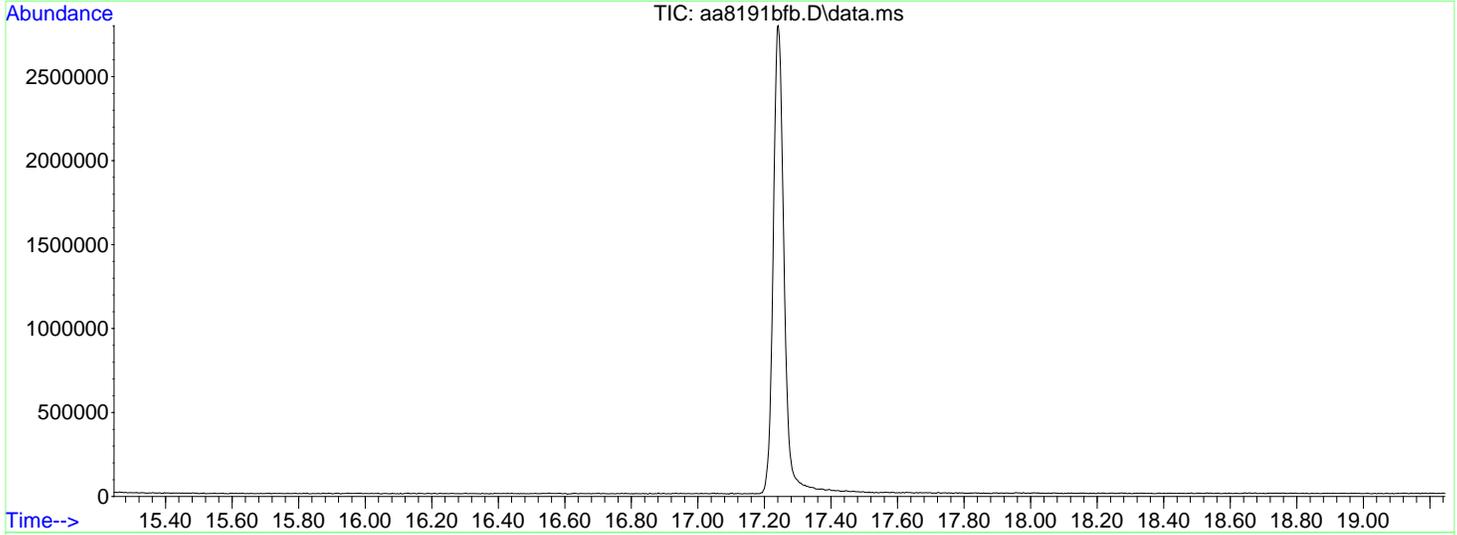
Runs with this BFB:

Lab Sample Number	Date File	Field Sample	Date/Time of Sample/Standard Analysis
BFB	AA8191BFB	NA	8/6/2018 11:26:00 AM
10 PPBV DCVS	AA8192DCVS	NA	8/6/2018 1:07:00 PM
METHOD BLANK	AA8193BLK	NA	8/6/2018 1:42:00 PM
02 PPBV RLLCS	AA8194RLLCS	NA	8/6/2018 2:23:00 PM
E18-06141-07	AA8206	SS-101	8/6/2018 10:28:00 PM
E18-06141-08	AA8207	SS-102	8/6/2018 11:02:00 PM
E18-06141-09	AA8208	SS-103	8/6/2018 11:35:00 PM
10 PPBV CCCVS	AA8213CCCVS	NA	8/7/2018 2:22:00 AM

Data Path : C:\DATA\08-06-18\
 Data File : aa8191bfb.D
 Acq On : 6 Aug 2018 11:26 am
 Operator : jls
 Sample : BFB
 Misc : ALM029426
 ALS Vial : 1 Sample Multiplier: 1

Integration File: rteint.p

Method : C:\msdchem\1\METHODS\0725.M
 Title : TO-15 on the Agilent 7890A / 5975C
 Last Update : Wed Jul 25 14:15:57 2018



AutoFind: Scans 5335, 5336, 5337; Background Corrected with Scan 5309

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	8	40	32.6	177003	PASS
75	95	30	66	61.6	334482	PASS
95	95	100	100	100.0	542971	PASS
96	95	5	9	6.7	36432	PASS
173	174	0.00	2	0.9	2635	PASS
174	95	50	100	52.8	286464	PASS
175	174	4	9	7.6	21901	PASS
176	174	93	101	97.1	278015	PASS
177	176	5	9	6.2	17336	PASS

Data Path: C:\DATA\08-07-18\
Data File: AA8221BFB.D
Acq On: 8/7/2018 9:33:00AM
Operator: jls
Sample: BFB
Misc: ALM029426
ALS Vial: 1 **Multiplier:** 1
Integration File: rteint.p
Method: C:\msdchem\1\METHODS\0725.M
Last Update: Wed Jul 25 14:15:57 2018

Spectrum Information:

Pass/Fail	Target Mass	Rel. to Mass	Lower Limit %	Higher Limit %	Raw Abundance	% Relative Abundance
PASS	50	95	8	40	155635	33.4
PASS	75	95	30	66	289323	62.1
PASS	95	95	100	100	466217	100.0
PASS	96	95	5	9	29664	6.4
PASS	173	174	0.00	2	2337	1.0
PASS	174	95	50	100	240853	51.7
PASS	175	174	4	9	17856	7.4
PASS	176	174	93	101	230955	95.9
PASS	177	176	5	9	15117	6.5

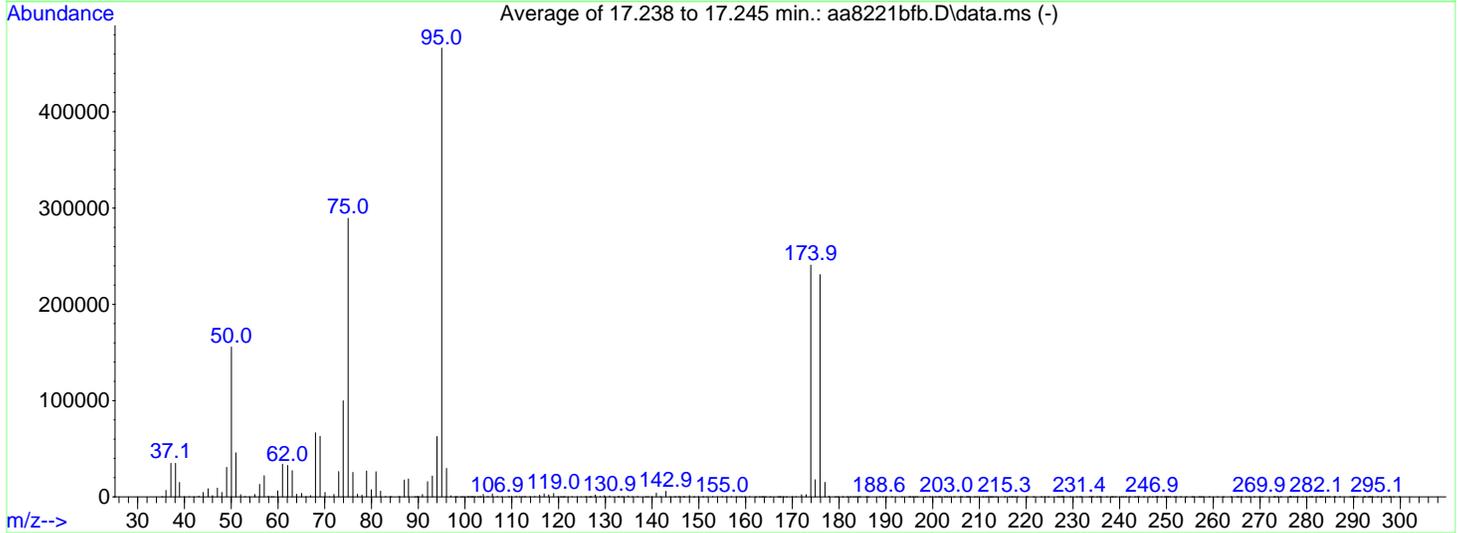
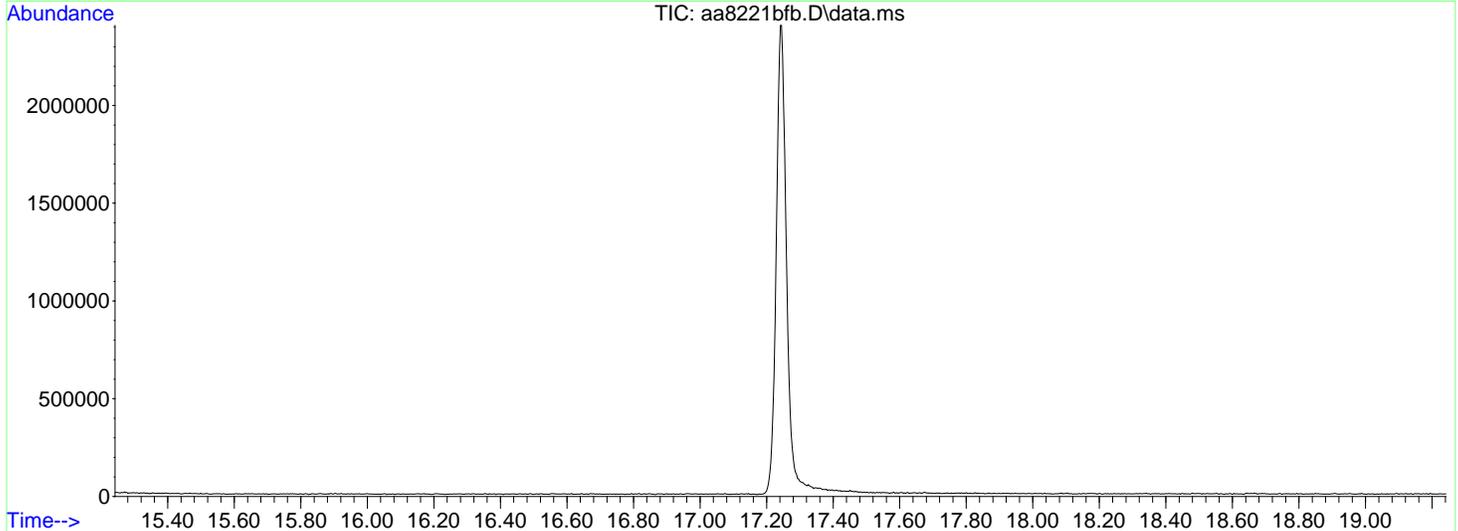
Runs with this BFB:

Lab Sample Number	Date File	Field Sample	Date/Time of Sample/Standard Analysis
BFB	AA8221BFB	NA	8/7/2018 9:33:00 AM
10 PPBV DCVS	AA8222DCVS	NA	8/7/2018 10:17:00 AM
METHOD BLANK	AA8223BLK	NA	8/7/2018 11:01:00 AM
02 PPBV RLLCS	AA8224RLLCS	NA	8/7/2018 11:47:00 AM
E18-06141-08	AA8232	SS-102	8/7/2018 4:57:00 PM
E18-06141-09	AA8233	SS-103	8/7/2018 5:31:00 PM
10 PPBV CCCVS	AA8242CCCVS	NA	8/7/2018 10:31:00 PM

Data Path : C:\DATA\08-07-18\
 Data File : aa8221bfb.D
 Acq On : 7 Aug 2018 9:33 am
 Operator : jjw
 Sample : BFB
 Misc : ALM029426
 ALS Vial : 1 Sample Multiplier: 1

Integration File: rteint.p

Method : C:\msdchem\1\METHODS\0725.M
 Title : TO-15 on the Agilent 7890A / 5975C
 Last Update : Wed Jul 25 14:15:57 2018



AutoFind: Scans 5334, 5335, 5336; Background Corrected with Scan 5312

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	8	40	33.4	155635	PASS
75	95	30	66	62.1	289323	PASS
95	95	100	100	100.0	466217	PASS
96	95	5	9	6.4	29664	PASS
173	174	0.00	2	1.0	2337	PASS
174	95	50	100	51.7	240853	PASS
175	174	4	9	7.4	17856	PASS
176	174	93	101	95.9	230955	PASS
177	176	5	9	6.5	15117	PASS

Method Blank Report

Lab Sample Name: METHOD BLANK
Field Sample Name: METHOD BLANK
Matrix: Air
Dilution Factor: 1

Data File: AA7473BLK
Date Analyzed: 6/13/2018
Sample Volume: 500ml
GC/MS Column: RTX-1, 0.32 mmID

Runs with this Method Blank:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7471BFB]	06/13/2018 8:30
10 PPBV DCVS [AA7472DCVS]	06/13/2018 9:18
METHOD BLANK [AA7473BLK]	06/13/2018 10:25
02 PPBV RLLCS [AA7474RLLCS]	06/13/2018 11:14
3059 [AA7475]	06/13/2018 12:20
10 PPBV CCCVS [AA7485CCCVS]	06/13/2018 20:47

Compound	CAS #	Reporting Limit (ppbv)	Concentration (ppbv)
Acetone	67-64-1	0.20	ND
Benzene	71-43-2	0.15	ND
Bromodichloromethane	75-27-4	0.20	ND
Bromoform	75-25-2	0.20	ND
Bromomethane	74-83-9	0.20	ND
1,3-Butadiene	106-99-0	0.20	ND
Chlorobenzene	108-90-7	0.20	ND
Chloroethane	75-00-3	0.20	ND
Chloroform	67-66-3	0.20	ND
Chloromethane	74-87-3	0.20	ND
Carbon disulfide	75-15-0	0.20	ND
Carbon tetrachloride	56-23-5	0.20	ND
Cyclohexane	110-82-7	0.12	ND
Dibromochloromethane	124-48-1	0.20	ND
1,2-Dibromoethane	106-93-4	0.20	ND
1,2-Dichlorobenzene	95-50-1	0.20	ND
1,3-Dichlorobenzene	541-73-1	0.20	ND
1,4-Dichlorobenzene	106-46-7	0.20	ND
Dichlorodifluoromethane	75-71-8	0.20	ND
1,1-Dichloroethane	75-34-3	0.20	ND
1,2-Dichloroethane	107-06-2	0.20	ND
1,1-Dichloroethene	75-35-4	0.20	ND
1,2-Dichloroethene (cis)	156-59-2	0.19	ND
1,2-Dichloroethene (trans)	156-60-5	0.20	ND
1,2-Dichloropropane	78-87-5	0.20	ND
1,3-Dichloropropene (cis)	10061-01-5	0.19	ND
1,3-Dichloropropene (trans)	10061-02-6	0.18	ND
1,2-Dichlorotetrafluoroethane	76-14-2	0.20	ND
1,4-Dioxane	123-91-1	0.17	ND
Ethylbenzene	100-41-4	0.20	ND
n-Heptane	142-82-5	0.20	ND
1,3-Hexachlorobutadiene	87-68-3	0.20	ND
n-Hexane	110-54-3	0.20	ND
Methylene chloride	75-09-2	0.20	ND
Methyl ethyl ketone	78-93-3	0.13	ND
Methyl isobutyl ketone	108-10-1	0.20	ND

Method Blank must be less than the Practical Quantitation Limit (PQL).

Method Blank Report

Lab Sample Name: METHOD BLANK
Field Sample Name: METHOD BLANK
Matrix: Air
Dilution Factor: 1

Data File: AA7473BLK
Date Analyzed: 6/13/2018
Sample Volume: 500ml
GC/MS Column: RTX-1, 0.32 mmID

Runs with this Method Blank:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7471BFB]	06/13/2018 8:30
10 PPBV DCVS [AA7472DCVS]	06/13/2018 9:18
METHOD BLANK [AA7473BLK]	06/13/2018 10:25
02 PPBV RLLCS [AA7474RLLCS]	06/13/2018 11:14
3059 [AA7475]	06/13/2018 12:20
10 PPBV CCCVS [AA7485CCCVS]	06/13/2018 20:47

Compound	CAS #	Reporting Limit (ppbv)	Concentration (ppbv)
Methyl tert-butyl ether	1634-04-4	0.11	ND
Styrene	100-42-5	0.14	ND
Tert-butyl alcohol	75-65-0	0.18	ND
1,1,2,2-Tetrachloroethane	79-34-5	0.17	ND
Tetrachloroethene	127-18-4	0.20	ND
Toluene	108-88-3	0.17	ND
1,2,4-Trichlorobenzene	120-82-1	0.20	ND
1,1,1-Trichloroethane	71-55-6	0.20	ND
1,1,2-Trichloroethane	79-00-5	0.20	ND
Trichloroethene	79-01-6	0.20	ND
Trichlorofluoromethane	75-69-4	0.20	ND
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	0.20	ND
1,2,4-Trimethylbenzene	95-63-6	0.12	ND
1,3,5-Trimethylbenzene	108-67-8	0.17	ND
2,2,4-Trimethylpentane	540-84-1	0.20	ND
Vinyl bromide	593-60-2	0.20	ND
Vinyl chloride	75-01-4	0.20	ND
Xylenes (m&p)	179601-23-1	0.20	ND
Xylenes (o)	95-47-6	0.20	ND

Method Blank must be less than the Practical Quantitation Limit (PQL).

Data Path : C:\DATA\06-13-18\
 Data File : aa7473blk.D
 Acq On : 13 Jun 2018 10:25 am
 Operator : jls
 Sample : Method Blank
 Misc : 1127, 500cc
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Jun 14 13:54:42 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 13:51:08 2018
 Response via : Initial Calibration

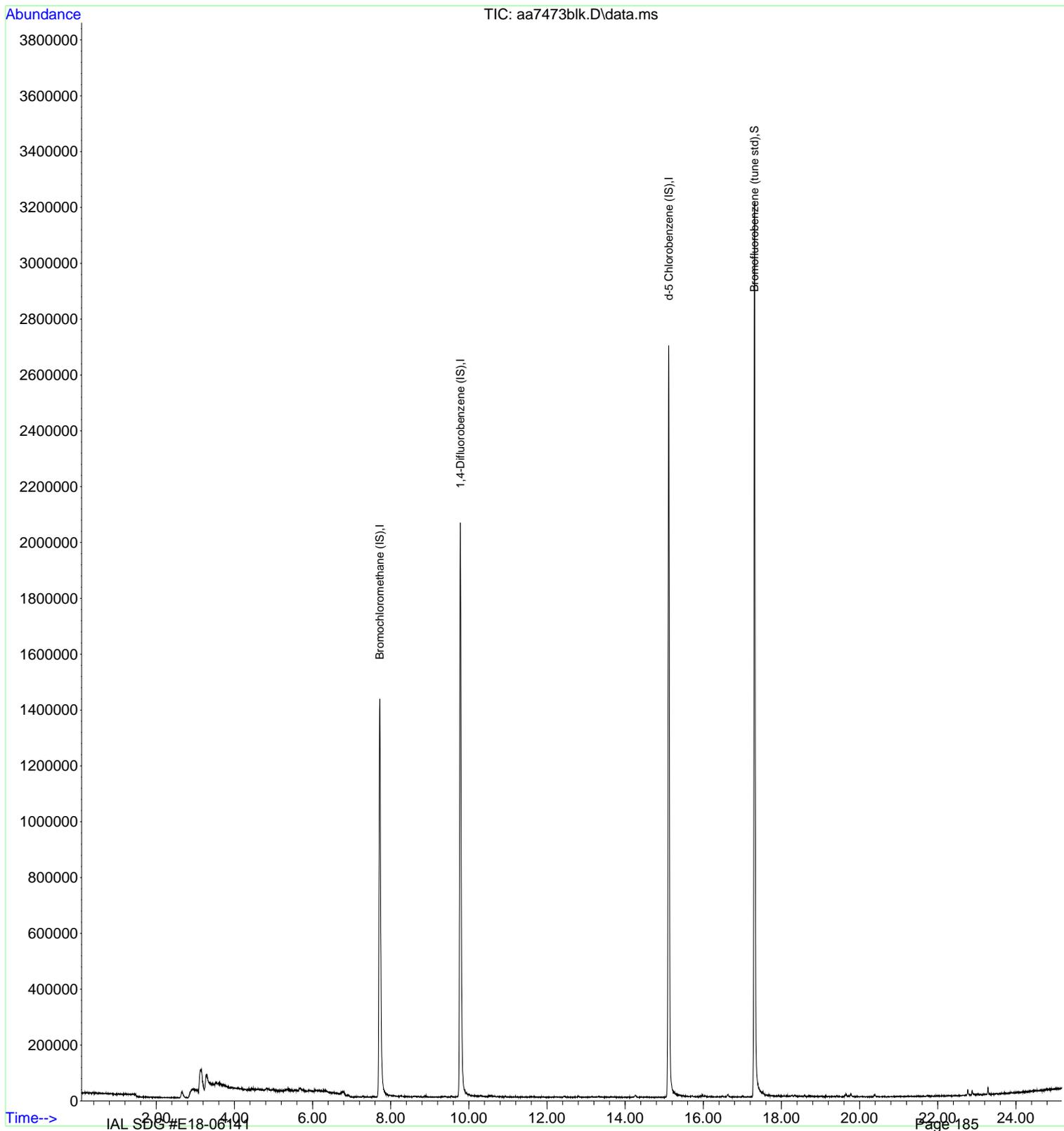
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Bromochloromethane (IS)	7.724	130	510475	10.00	ppbV	0.00
38) 1,4-Difluorobenzene (IS)	9.785	114	1863647	10.00	ppbV #	0.00
55) d-5 Chlorobenzene (IS)	15.116	117	1558649	10.00	ppbV	0.00
System Monitoring Compounds						
64) Bromofluorobenzene (tu...	17.312	95	1385718	9.66	ppbV	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\06-13-18\
 Data File : aa7473blk.D
 Acq On : 13 Jun 2018 10:25 am
 Operator : jls
 Sample : Method Blank
 Misc : 1127, 500cc
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Jun 14 13:54:42 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 13:51:08 2018
 Response via : Initial Calibration



Method Blank Report

Lab Sample Name: METHOD BLANK
Field Sample Name: METHOD BLANK
Matrix: Air
Dilution Factor: 1

Data File: AA8193BLK
Date Analyzed: 8/6/2018
Sample Volume: 500ml
GC/MS Column: RTX-1, 0.32 mmID

Runs with this Method Blank:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA8191BFB]	08/06/2018 11:26
10 PPBV DCVS [AA8192DCVS]	08/06/2018 13:07
METHOD BLANK [AA8193BLK]	08/06/2018 13:42
02 PPBV RLLCS [AA8194RLLCS]	08/06/2018 14:23
E18-06141-07 [AA8206]	08/06/2018 22:28
E18-06141-08 [AA8207]	08/06/2018 23:02
E18-06141-09 [AA8208]	08/06/2018 23:35
10 PPBV CCCVS [AA8213CCCVS]	08/07/2018 2:22

Compound	CAS #	Reporting Limit (ppbv)	Concentration (ppbv)
Acetone	67-64-1	0.20	ND
Benzene	71-43-2	0.15	ND
Bromodichloromethane	75-27-4	0.20	ND
Bromoform	75-25-2	0.20	ND
Bromomethane	74-83-9	0.20	ND
1,3-Butadiene	106-99-0	0.20	ND
Chlorobenzene	108-90-7	0.20	ND
Chloroethane	75-00-3	0.20	ND
Chloroform	67-66-3	0.20	ND
Chloromethane	74-87-3	0.20	ND
Carbon disulfide	75-15-0	0.20	ND
Carbon tetrachloride	56-23-5	0.20	ND
Cyclohexane	110-82-7	0.12	ND
Dibromochloromethane	124-48-1	0.20	ND
1,2-Dibromoethane	106-93-4	0.20	ND
1,2-Dichlorobenzene	95-50-1	0.20	ND
1,3-Dichlorobenzene	541-73-1	0.20	ND
1,4-Dichlorobenzene	106-46-7	0.20	ND
Dichlorodifluoromethane	75-71-8	0.20	ND
1,1-Dichloroethane	75-34-3	0.20	ND
1,2-Dichloroethane	107-06-2	0.20	ND
1,1-Dichloroethene	75-35-4	0.20	ND
1,2-Dichloroethene (cis)	156-59-2	0.19	ND
1,2-Dichloroethene (trans)	156-60-5	0.20	ND
1,2-Dichloropropane	78-87-5	0.20	ND
1,3-Dichloropropene (cis)	10061-01-5	0.19	ND
1,3-Dichloropropene (trans)	10061-02-6	0.18	ND
1,2-Dichlorotetrafluoroethane	76-14-2	0.20	ND
1,4-Dioxane	123-91-1	0.17	ND
Ethylbenzene	100-41-4	0.20	ND
n-Heptane	142-82-5	0.20	ND
1,3-Hexachlorobutadiene	87-68-3	0.20	ND
n-Hexane	110-54-3	0.20	ND
Methylene chloride	75-09-2	0.20	ND

Method Blank must be less than the Practical Quantitation Limit (PQL).

Method Blank Report

Lab Sample Name: METHOD BLANK
Field Sample Name: METHOD BLANK
Matrix: Air
Dilution Factor: 1

Data File: AA8193BLK
Date Analyzed: 8/6/2018
Sample Volume: 500ml
GC/MS Column: RTX-1, 0.32 mmID

Runs with this Method Blank:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA8191BFB]	08/06/2018 11:26
10 PPBV DCVS [AA8192DCVS]	08/06/2018 13:07
METHOD BLANK [AA8193BLK]	08/06/2018 13:42
02 PPBV RLLCS [AA8194RLLCS]	08/06/2018 14:23
E18-06141-07 [AA8206]	08/06/2018 22:28
E18-06141-08 [AA8207]	08/06/2018 23:02
E18-06141-09 [AA8208]	08/06/2018 23:35
10 PPBV CCCVS [AA8213CCCVS]	08/07/2018 2:22

Compound	CAS #	Reporting Limit (ppbv)	Concentration (ppbv)
Methyl ethyl ketone	78-93-3	0.13	ND
Methyl isobutyl ketone	108-10-1	0.20	ND
Methyl tert-butyl ether	1634-04-4	0.11	ND
Styrene	100-42-5	0.14	ND
Tert-butyl alcohol	75-65-0	0.18	ND
1,1,2,2-Tetrachloroethane	79-34-5	0.17	ND
Tetrachloroethene	127-18-4	0.20	ND
Toluene	108-88-3	0.17	ND
1,2,4-Trichlorobenzene	120-82-1	0.20	ND
1,1,1-Trichloroethane	71-55-6	0.20	ND
1,1,2-Trichloroethane	79-00-5	0.20	ND
Trichloroethene	79-01-6	0.20	ND
Trichlorofluoromethane	75-69-4	0.20	ND
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	0.20	ND
1,2,4-Trimethylbenzene	95-63-6	0.12	ND
1,3,5-Trimethylbenzene	108-67-8	0.17	ND
2,2,4-Trimethylpentane	540-84-1	0.20	ND
Vinyl bromide	593-60-2	0.20	ND
Vinyl chloride	75-01-4	0.20	ND
Xylenes (m&p)	179601-23-1	0.20	ND
Xylenes (o)	95-47-6	0.20	ND

Method Blank must be less than the Practical Quantitation Limit (PQL).

Data Path : C:\DATA\08-06-18\
 Data File : aa8193blk.D
 Acq On : 6 Aug 2018 1:42 pm
 Operator : jls
 Sample : Method Blank
 Misc : 1127, 500cc
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Aug 06 14:08:12 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

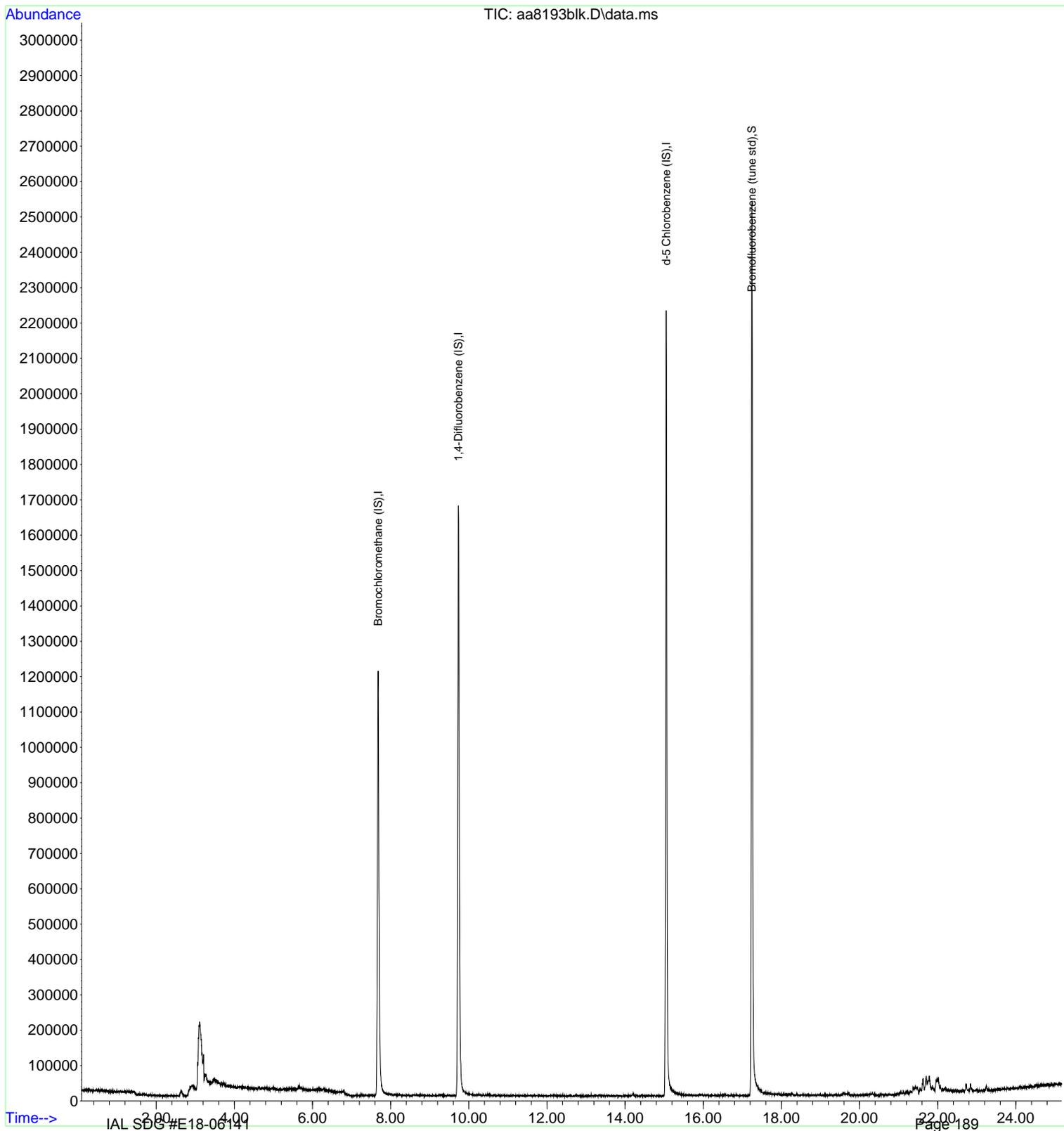
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Bromochloromethane (IS)	7.679	130	418368	10.00	ppbV	0.00
38) 1,4-Difluorobenzene (IS)	9.731	114	1550382	10.00	ppbV #	0.00
55) d-5 Chlorobenzene (IS)	15.052	117	1251713	10.00	ppbV	0.00
System Monitoring Compounds						
64) Bromofluorobenzene (tu...	17.245	95	1143580	10.02	ppbV	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\08-06-18\
Data File : aa8193blk.D
Acq On : 6 Aug 2018 1:42 pm
Operator : jls
Sample : Method Blank
Misc : 1127, 500cc
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Aug 06 14:08:12 2018
Quant Method : C:\msdchem\1\METHODS\0725.M
Quant Title : TO-15 on the Agilent 7890A / 5975C
QLast Update : Wed Jul 25 14:15:57 2018
Response via : Initial Calibration



Method Blank Report

Lab Sample Name: METHOD BLANK
Field Sample Name: METHOD BLANK
Matrix: Air
Dilution Factor: 1

Data File: AA8223BLK
Date Analyzed: 8/7/2018
Sample Volume: 500ml
GC/MS Column: RTX-1, 0.32 mmID

Runs with this Method Blank:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA8221BFB]	08/07/2018 9:33
10 PPBV DCVS [AA8222DCVS]	08/07/2018 10:17
METHOD BLANK [AA8223BLK]	08/07/2018 11:01
02 PPBV RLLCS [AA8224RLLCS]	08/07/2018 11:47
E18-06141-08 [AA8232]	08/07/2018 16:57
E18-06141-09 [AA8233]	08/07/2018 17:31
10 PPBV CCCVS [AA8242CCCVS]	08/07/2018 22:31

Compound	CAS #	Reporting Limit (ppbv)	Concentration (ppbv)
Acetone	67-64-1	0.20	ND
Benzene	71-43-2	0.15	ND
Bromodichloromethane	75-27-4	0.20	ND
Bromoform	75-25-2	0.20	ND
Bromomethane	74-83-9	0.20	ND
1,3-Butadiene	106-99-0	0.20	ND
Chlorobenzene	108-90-7	0.20	ND
Chloroethane	75-00-3	0.20	ND
Chloroform	67-66-3	0.20	ND
Chloromethane	74-87-3	0.20	ND
Carbon disulfide	75-15-0	0.20	ND
Carbon tetrachloride	56-23-5	0.20	ND
Cyclohexane	110-82-7	0.12	ND
Dibromochloromethane	124-48-1	0.20	ND
1,2-Dibromoethane	106-93-4	0.20	ND
1,2-Dichlorobenzene	95-50-1	0.20	ND
1,3-Dichlorobenzene	541-73-1	0.20	ND
1,4-Dichlorobenzene	106-46-7	0.20	ND
Dichlorodifluoromethane	75-71-8	0.20	ND
1,1-Dichloroethane	75-34-3	0.20	ND
1,2-Dichloroethane	107-06-2	0.20	ND
1,1-Dichloroethene	75-35-4	0.20	ND
1,2-Dichloroethene (cis)	156-59-2	0.19	ND
1,2-Dichloroethene (trans)	156-60-5	0.20	ND
1,2-Dichloropropane	78-87-5	0.20	ND
1,3-Dichloropropene (cis)	10061-01-5	0.19	ND
1,3-Dichloropropene (trans)	10061-02-6	0.18	ND
1,2-Dichlorotetrafluoroethane	76-14-2	0.20	ND
1,4-Dioxane	123-91-1	0.17	ND
Ethylbenzene	100-41-4	0.20	ND
n-Heptane	142-82-5	0.20	ND
1,3-Hexachlorobutadiene	87-68-3	0.20	ND
n-Hexane	110-54-3	0.20	ND
Methylene chloride	75-09-2	0.20	ND
Methyl ethyl ketone	78-93-3	0.13	ND

Method Blank must be less than the Practical Quantitation Limit (PQL).

Method Blank Report

Lab Sample Name: METHOD BLANK
Field Sample Name: METHOD BLANK
Matrix: Air
Dilution Factor: 1

Data File: AA8223BLK
Date Analyzed: 8/7/2018
Sample Volume: 500ml
GC/MS Column: RTX-1, 0.32 mmID

Runs with this Method Blank:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA8221BFB]	08/07/2018 9:33
10 PPBV DCVS [AA8222DCVS]	08/07/2018 10:17
METHOD BLANK [AA8223BLK]	08/07/2018 11:01
02 PPBV RLLCS [AA8224RLLCS]	08/07/2018 11:47
E18-06141-08 [AA8232]	08/07/2018 16:57
E18-06141-09 [AA8233]	08/07/2018 17:31
10 PPBV CCCVS [AA8242CCCVS]	08/07/2018 22:31

Compound	CAS #	Reporting Limit (ppbv)	Concentration (ppbv)
Methyl isobutyl ketone	108-10-1	0.20	ND
Methyl tert-butyl ether	1634-04-4	0.11	ND
Styrene	100-42-5	0.14	ND
Tert-butyl alcohol	75-65-0	0.18	ND
1,1,2,2-Tetrachloroethane	79-34-5	0.17	ND
Tetrachloroethene	127-18-4	0.20	ND
Toluene	108-88-3	0.17	ND
1,2,4-Trichlorobenzene	120-82-1	0.20	ND
1,1,1-Trichloroethane	71-55-6	0.20	ND
1,1,2-Trichloroethane	79-00-5	0.20	ND
Trichloroethene	79-01-6	0.20	ND
Trichlorofluoromethane	75-69-4	0.20	ND
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	0.20	ND
1,2,4-Trimethylbenzene	95-63-6	0.12	ND
1,3,5-Trimethylbenzene	108-67-8	0.17	ND
2,2,4-Trimethylpentane	540-84-1	0.20	ND
Vinyl bromide	593-60-2	0.20	ND
Vinyl chloride	75-01-4	0.20	ND
Xylenes (m&p)	179601-23-1	0.20	ND
Xylenes (o)	95-47-6	0.20	ND

Method Blank must be less than the Practical Quantitation Limit (PQL).

Data Path : C:\DATA\08-07-18\
 Data File : aa8223blk.D
 Acq On : 7 Aug 2018 11:01 am
 Operator : jjw
 Sample : Method Blank
 Misc : 1127, 500cc
 ALS Vial : 3 Sample Multiplier: 1

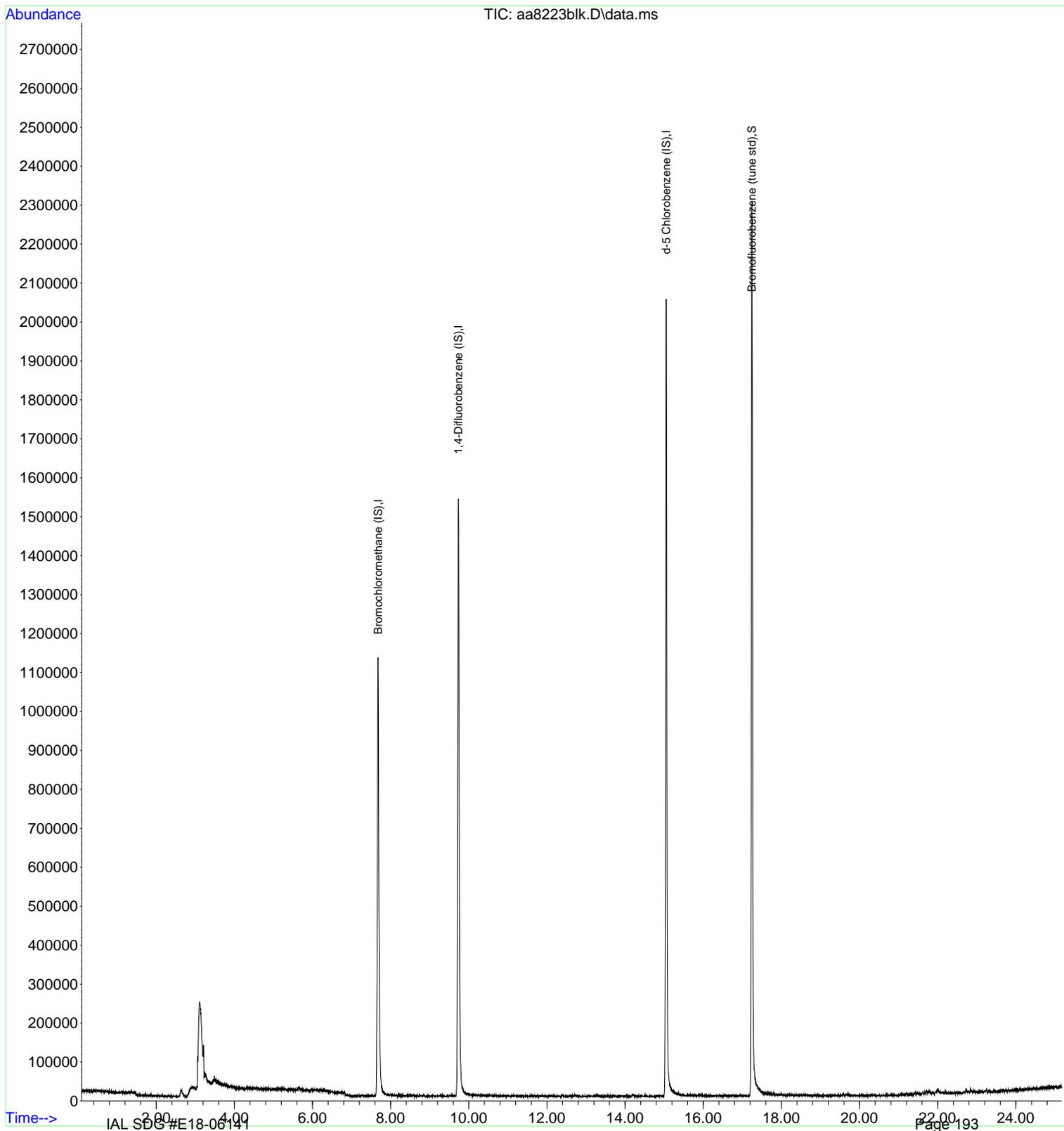
Quant Time: Aug 07 11:31:16 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Bromochloromethane (IS)	7.682	130	386755	10.00	ppbV	0.00
38) 1,4-Difluorobenzene (IS)	9.734	114	1385734	10.00	ppbV #	0.00
55) d-5 Chlorobenzene (IS)	15.052	117	1124467	10.00	ppbV	0.00
System Monitoring Compounds						
64) Bromofluorobenzene (tu...	17.245	95	997109	9.72	ppbV	0.00
Target Compounds						Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\08-07-18\
 Data File : aa8223blk.D
 Acq On : 7 Aug 2018 11:01 am
 Operator : jjw
 Sample : Method Blank
 Misc : 1127, 500cc
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Aug 07 11:31:16 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration



Integrated Analytical Laboratories, LLC

Volatile Organic Compounds by EPA Method TO-15

Laboratory Sample Duplicate Report

SDG Number: E18-04189
 IAL Sample ID: E18-04189-01
 Matrix: Air
 Summa ID: 5096

Date Received: 6/8/18
 Date Analyzed: 6/13/18,6/13/18
 Lab Data File#: AA7478,AA7479
 Dilution Factor: 1
 Injection Volume: 500ml
 GC/MS Column: RTX-1, 0.32 mmID

Compound	CAS #	Sample E18-04189-01 Concentration Reported		Sample Dup E18-04189-21 Concentration Reported		Reporting Limits ppbv	RPD
		ppbv	Q	ppbv	Q		
Acetone	67-64-1	3.1		3.1		0.40	0.00%
Allyl Chloride	107-05-1		0.40 U		0.40 U	0.40	0.00%
Benzene	71-43-2		0.40 U		0.40 U	0.40	0.00%
Bromodichloromethane	75-27-4		0.40 U		0.40 U	0.40	0.00%
Bromoform	75-25-2		0.40 U		0.40 U	0.40	0.00%
Bromomethane	74-83-9		0.40 U		0.40 U	0.40	0.00%
1,3-Butadiene	106-99-0		0.40 U		0.40 U	0.40	0.00%
Chlorobenzene	108-90-7		0.40 U		0.40 U	0.40	0.00%
Chloroethane	75-00-3		0.40 U		0.40 U	0.40	0.00%
Chloroform	67-66-3		0.40 U		0.40 U	0.40	0.00%
Chloromethane	74-87-3		0.40 U		0.40 U	0.40	0.00%
Carbon disulfide	75-15-0		0.40 U		0.40 U	0.40	0.00%
Carbon tetrachloride	56-23-5		0.40 U		0.40 U	0.40	0.00%
2-Chlorotoluene	95-49-8		0.40 U		0.40 U	0.40	0.00%
Cyclohexane	110-82-7		0.40 U		0.40 U	0.40	0.00%
Dibromochloromethane	124-48-1		0.40 U		0.40 U	0.40	0.00%
1,2-Dibromoethane	106-93-4		0.40 U		0.40 U	0.40	0.00%
1,2-Dichlorobenzene	95-50-1		0.40 U		0.40 U	0.40	0.00%
1,3-Dichlorobenzene	541-73-1		0.40 U		0.40 U	0.40	0.00%
1,4-Dichlorobenzene	106-46-7		0.40 U		0.40 U	0.40	0.00%
Dichlorodifluoromethane	75-71-8		0.40 U		0.40 U	0.40	0.00%
1,1-Dichloroethane	75-34-3		0.40 U		0.40 U	0.40	0.00%
1,2-Dichloroethane	107-06-2		0.40 U		0.40 U	0.40	0.00%
1,1-Dichloroethene	75-35-4		0.40 U		0.40 U	0.40	0.00%
1,2-Dichloroethene (cis)	156-59-2		0.40 U		0.40 U	0.40	0.00%
1,2-Dichloroethene (trans)	156-60-5		0.40 U		0.40 U	0.40	0.00%
1,2-Dichloropropane	78-87-5		0.40 U		0.40 U	0.40	0.00%
1,3-Dichloropropene (cis)	10061-01-5		0.40 U		0.40 U	0.40	0.00%
1,3-Dichloropropene (trans)	10061-02-6		0.40 U		0.40 U	0.40	0.00%
1,2-Dichlorotetrafluoroethane	76-14-2		0.40 U		0.40 U	0.40	0.00%
Ethylbenzene	100-41-4		0.40 U		0.40 U	0.40	0.00%
4-Ethyltoluene	622-96-8		0.40 U		0.40 U	0.40	0.00%
n-Heptane	142-82-5		0.40 U		0.40 U	0.40	0.00%
1,3-Hexachlorobutadiene	87-68-3		0.40 U		0.40 U	0.40	0.00%
n-Hexane	110-54-3		0.40 U		0.40 U	0.40	0.00%
Methylene chloride	75-09-2	2.3		2.6		0.40	-12.24%
Methyl ethyl ketone	78-93-3		0.40 U		0.40 U	0.40	0.00%
Methyl isobutyl ketone	108-10-1		0.40 U		0.40 U	0.40	0.00%

Qualifir:

E=Concentration exceeds upper level of calibration range for instrument.

D=Extra dilution required for this compound. J=Duplicate samples do not met RPD criteria.

U=Compound ND or under reporting limit.

Integrated Analytical Laboratories, LLC

Volatile Organic Compounds by EPA Method TO-15

Laboratory Sample Duplicate Report

SDG Number: E18-04189
 IAL Sample ID: E18-04189-01
 Matrix: Air
 Summa ID: 5096

Date Received: 6/8/18
 Date Analyzed: 6/13/18,6/13/18
 Lab Data File#: AA7478,AA7479
 Dilution Factor: 1
 Injection Volume: 500ml
 GC/MS Column: RTX-1, 0.32 mmID

Compound	CAS #	Sample E18-04189-01 Concentration Reported		Sample Dup E18-04189-21 Concentration Reported		Reporting Limits ppbv	RPD
		ppbv	Q	ppbv	Q		
Methyl tert-butyl ether	1634-04-4	0.40	U	0.40	U	0.40	0.00%
Styrene	100-42-5	0.40	U	0.40	U	0.40	0.00%
Tert-butyl alcohol	75-65-0	0.40	U	0.40	U	0.40	0.00%
1,1,2,2-Tetrachloroethane	79-34-5	0.40	U	0.40	U	0.40	0.00%
Tetrachloroethene	127-18-4	0.40	U	0.40	U	0.40	0.00%
Toluene	108-88-3	0.40	U	0.40	U	0.40	0.00%
1,2,4-Trichlorobenzene	120-82-1	0.40	U	0.40	U	0.40	0.00%
1,1,1-Trichloroethane	71-55-6	0.40	U	0.40	U	0.40	0.00%
1,1,2-Trichloroethane	79-00-5	0.40	U	0.40	U	0.40	0.00%
Trichloroethene	79-01-6	0.40	U	0.40	U	0.40	0.00%
Trichlorofluoromethane	75-69-4	0.40	U	0.40	U	0.40	0.00%
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	0.40	U	0.40	U	0.40	0.00%
1,2,4-Trimethylbenzene	95-63-6	0.40	U	0.40	U	0.40	0.00%
1,3,5-Trimethylbenzene	108-67-8	0.40	U	0.40	U	0.40	0.00%
2,2,4-Trimethylpentane	540-84-1	0.40	U	0.40	U	0.40	0.00%
Vinyl bromide	593-60-2	0.40	U	0.40	U	0.40	0.00%
Vinyl chloride	75-01-4	0.40	U	0.40	U	0.40	0.00%
Xylenes (m&p)	179601-23-1	0.40	U	0.40	U	0.40	0.00%
Xylenes (o)	95-47-6	0.40	U	0.40	U	0.40	0.00%

RPD must be <25% for all laboratory duplicate samples. Laboratory duplicate samples are run once daily.
NC = The RPD could not be calculated since the compound was only detected in either the parent or duplicate sample.

Qualifir:

E=Concentration exceeds upper level of calibration range for instrument.
 D=Extra dilution required for this compound. J=Duplicate samples do not met RPD criteria.
 U=Compound ND or under reporting limit.

Data Path : C:\DATA\06-13-18\
 Data File : aa7478.D
 Acq On : 13 Jun 2018 4:48 pm
 Operator : jls
 Sample : E18-04189-01
 Misc : 5096, 500cc
 ALS Vial : 8 Sample Multiplier: 1

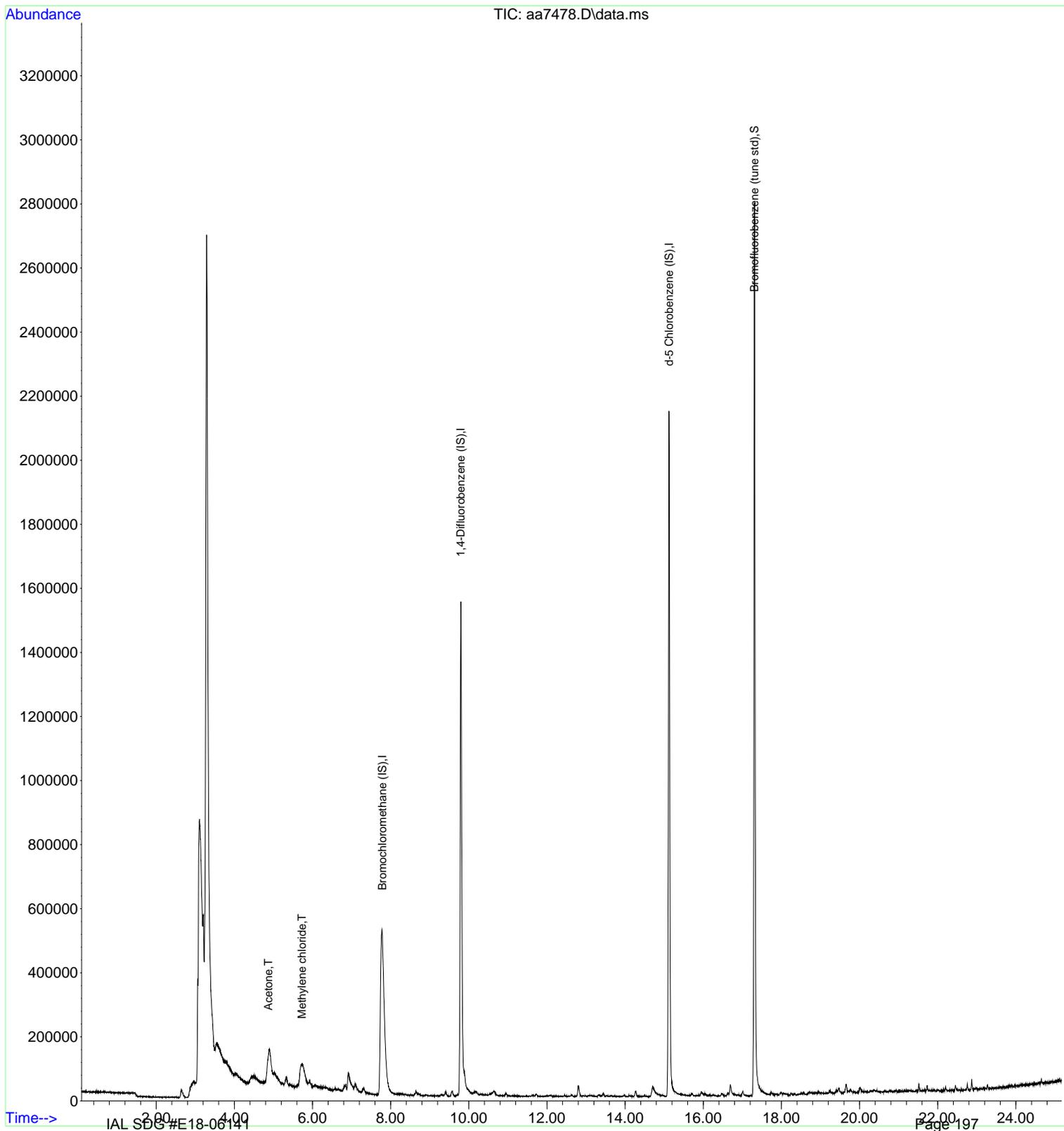
Quant Time: Jun 14 10:45:36 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 13:51:08 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
Internal Standards						
1) Bromochloromethane (IS)	7.788	130	409872	10.00	ppbV	0.06
38) 1,4-Difluorobenzene (IS)	9.795	114	1446308	10.00	ppbV #	0.01
55) d-5 Chlorobenzene (IS)	15.122	117	1304566	10.00	ppbV	0.01
System Monitoring Compounds						
64) Bromofluorobenzene (tu...	17.309	95	1183822	9.86	ppbV	0.00
Target Compounds						
14) Acetone	4.872	58	102946	3.11	ppbV	70
19) Methylene chloride	5.731	49	217471	2.27	ppbV	94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\06-13-18\
 Data File : aa7478.D
 Acq On : 13 Jun 2018 4:48 pm
 Operator : jls
 Sample : E18-04189-01
 Misc : 5096, 500cc
 ALS Vial : 8 Sample Multiplier: 1

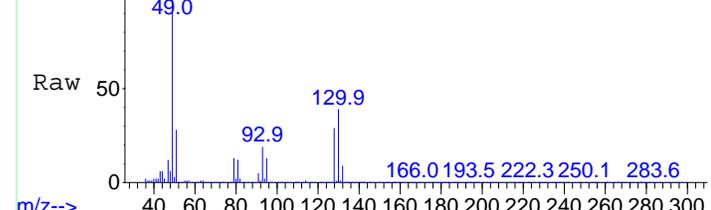
Quant Time: Jun 14 10:45:36 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 13:51:08 2018
 Response via : Initial Calibration



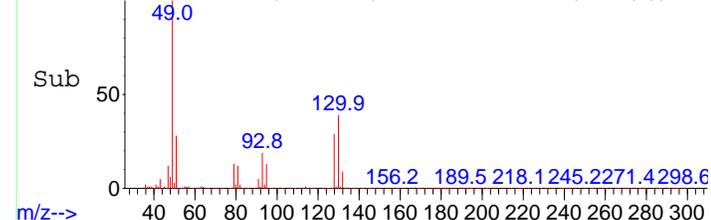
Abundance Scan 2378 (7.733 min): aa6994std03.D\data.ms (-2353) (-)



Abundance Scan 2395 (7.788 min): aa7478.D\data.ms

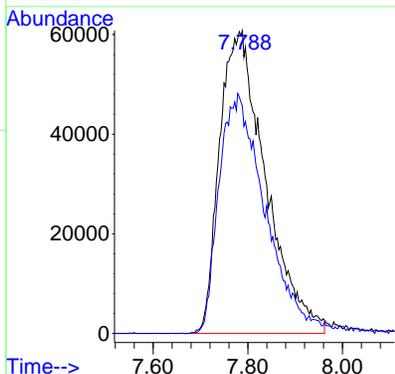


Abundance Scan 2395 (7.788 min): aa7478.D\data.ms (-2344) (-)

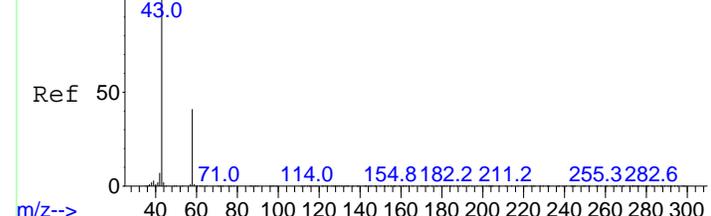


#1
Bromochloromethane (IS)
Concen: 10.00 ppbV
RT: 7.788 min Scan# 2395
Delta R.T. 0.064 min
Lab File: aa7478.D
Acq: 13 Jun 2018 4:48 pm

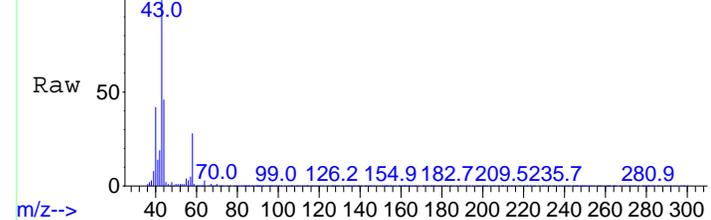
Tgt Ion: 130 Resp: 409872
Ion Ratio Lower Upper
130 100
128 78.8 62.6 94.0



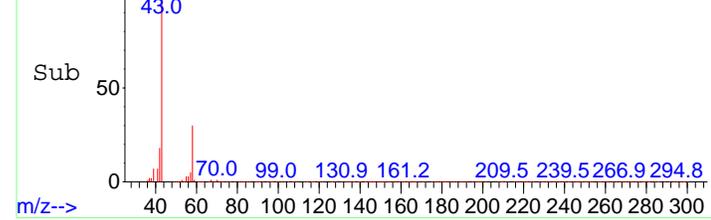
Abundance Scan 1470 (4.814 min): aa6994std03.D\data.ms (-1454) (-)



Abundance Scan 1488 (4.872 min): aa7478.D\data.ms

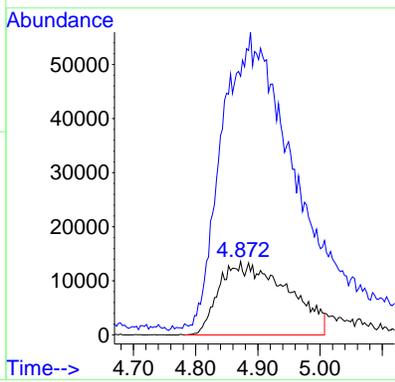


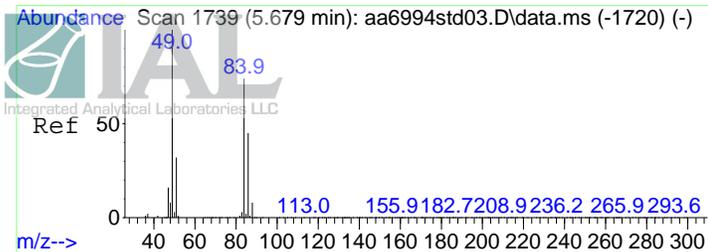
Abundance Scan 1488 (4.872 min): aa7478.D\data.ms (-1441) (-)



#14
Acetone
Concen: 3.11 ppbV
RT: 4.872 min Scan# 1488
Delta R.T. 0.052 min
Lab File: aa7478.D
Acq: 13 Jun 2018 4:48 pm

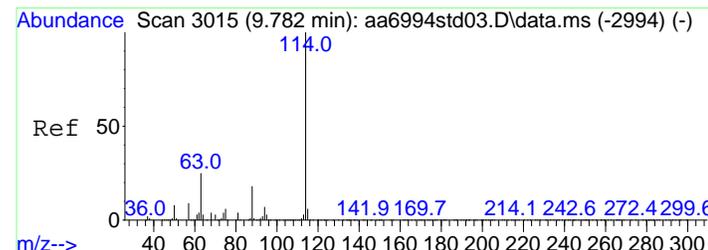
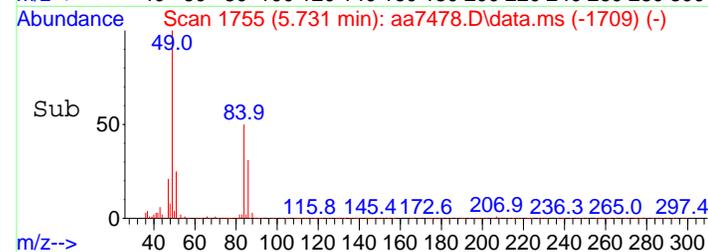
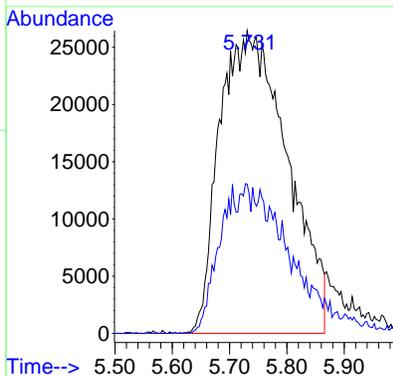
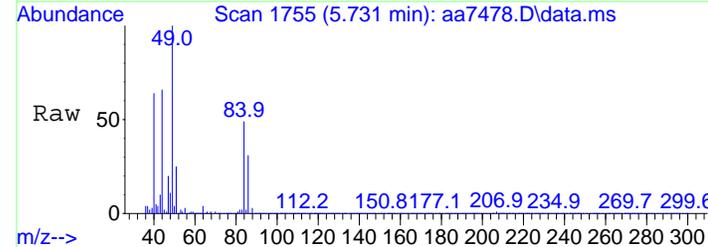
Tgt Ion: 58 Resp: 102946
Ion Ratio Lower Upper
58 100
43 392.3 263.2 394.8





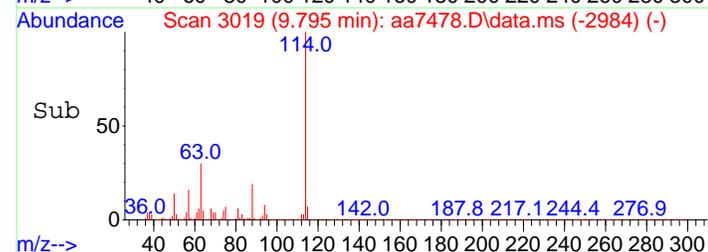
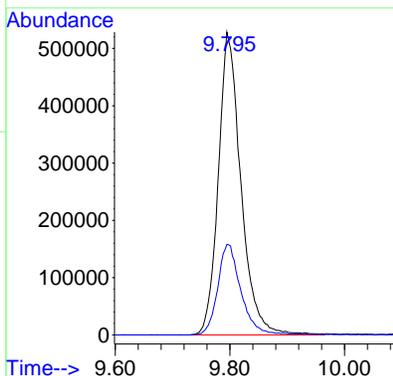
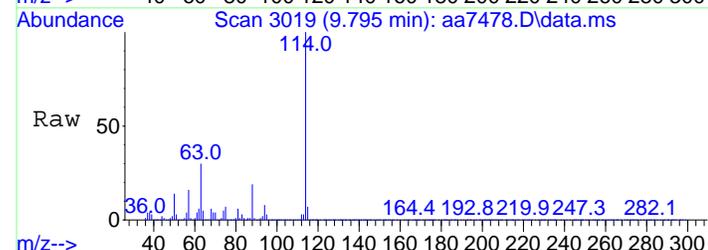
#19
 Methylene chloride
 Concen: 2.27 ppbV
 RT: 5.731 min Scan# 1755
 Delta R.T. 0.049 min
 Lab File: aa7478.D
 Acq: 13 Jun 2018 4:48 pm

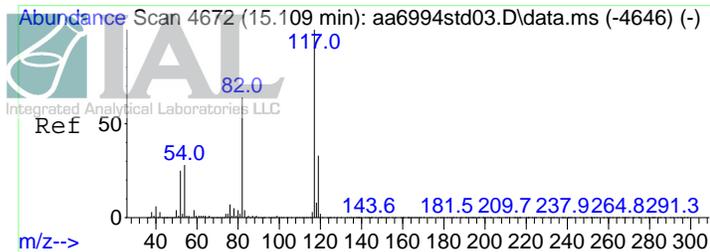
Tgt Ion: 49 Resp: 217471
 Ion Ratio Lower Upper
 49 100
 84 47.8 41.6 62.4



#38
 1,4-Difluorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 9.795 min Scan# 3019
 Delta R.T. 0.013 min
 Lab File: aa7478.D
 Acq: 13 Jun 2018 4:48 pm

Tgt Ion: 114 Resp: 1446308
 Ion Ratio Lower Upper
 114 100
 63 30.6 20.0 30.0#

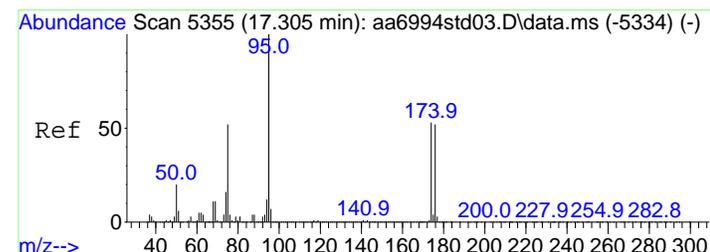
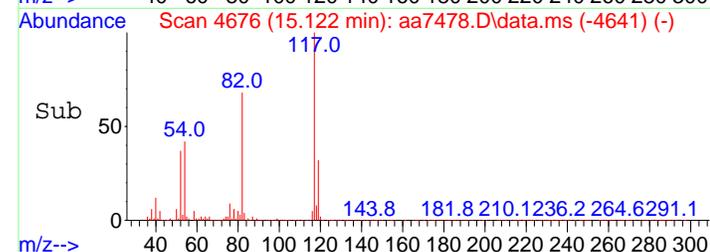
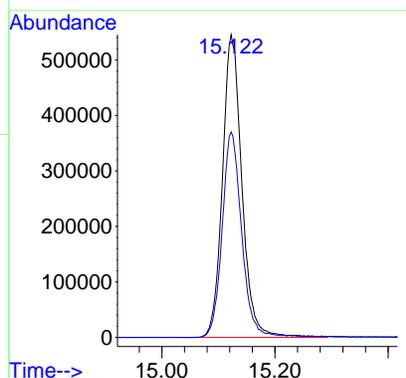
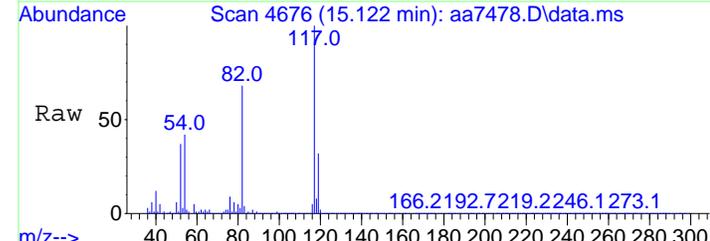




#55
 d-5 Chlorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 15.122 min Scan# 4676
 Delta R.T. 0.013 min
 Lab File: aa7478.D
 Acq: 13 Jun 2018 4:48 pm

Tgt Ion: 117 Resp: 1304566

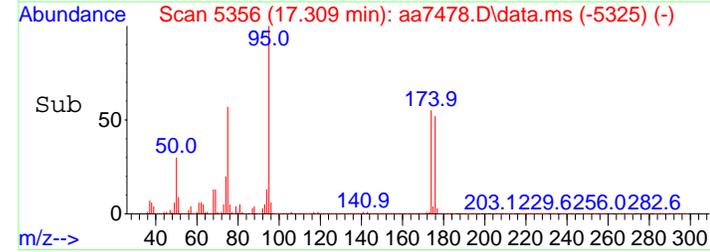
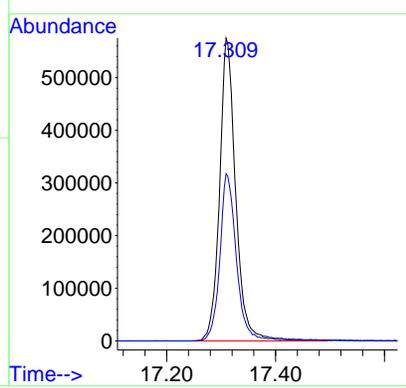
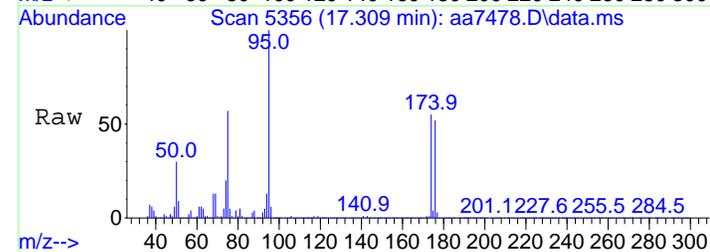
Ion	Ratio	Lower	Upper
117	100		
82	68.1	56.0	84.0



#64
 Bromofluorobenzene (tune std)
 Concen: 9.86 ppbV
 RT: 17.309 min Scan# 5356
 Delta R.T. 0.000 min
 Lab File: aa7478.D
 Acq: 13 Jun 2018 4:48 pm

Tgt Ion: 95 Resp: 1183822

Ion	Ratio	Lower	Upper
95	100		
174	56.7	61.5	92.3#



Data Path : C:\DATA\06-13-18\
 Data File : aa7479.D
 Acq On : 13 Jun 2018 5:27 pm
 Operator : jls
 Sample : E18-04189-21
 Misc : dup of E18-4189-01, can # 5096
 ALS Vial : 9 Sample Multiplier: 1

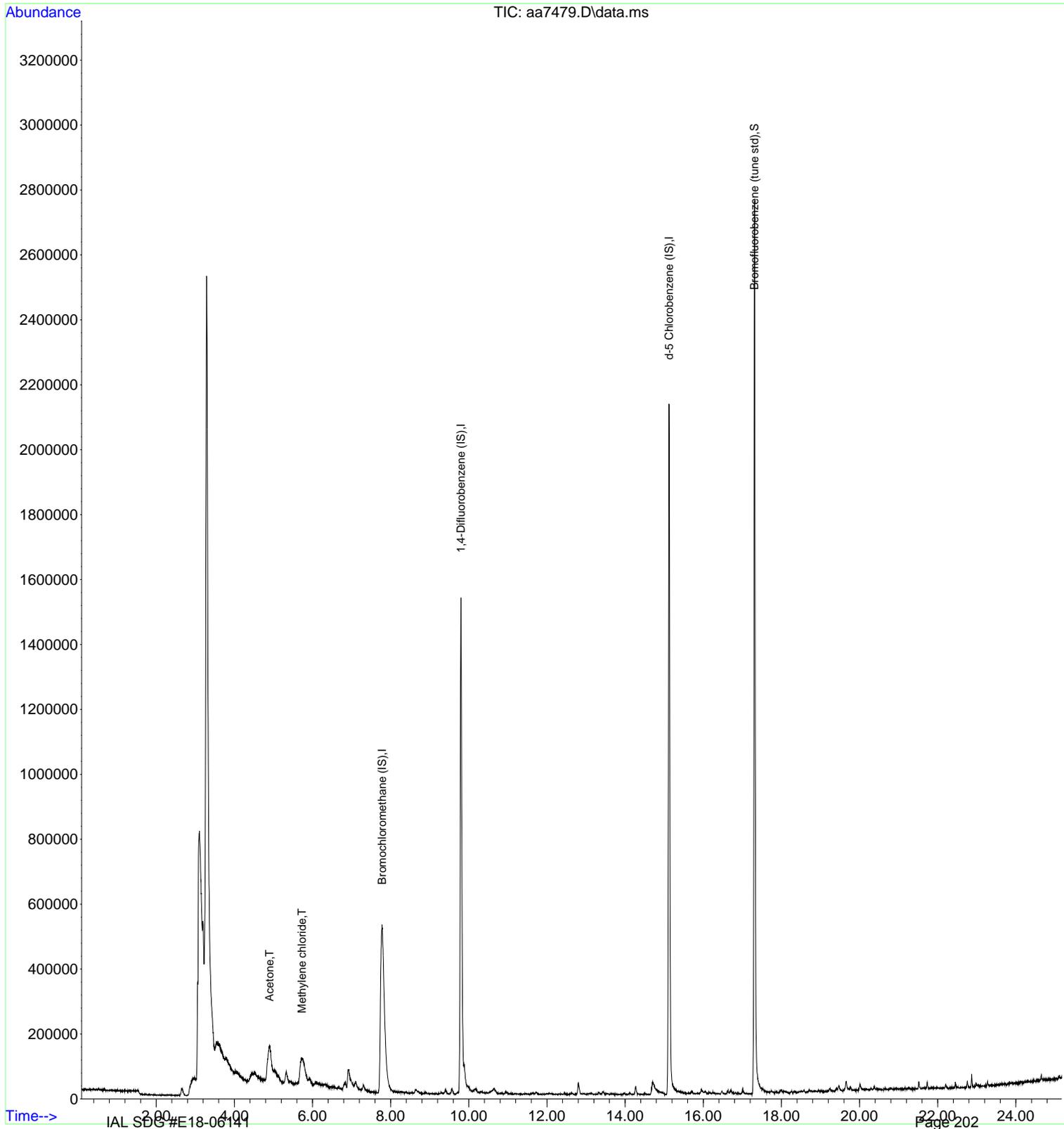
Quant Time: Jun 14 11:13:05 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 13:51:08 2018
 Response via : Initial Calibration

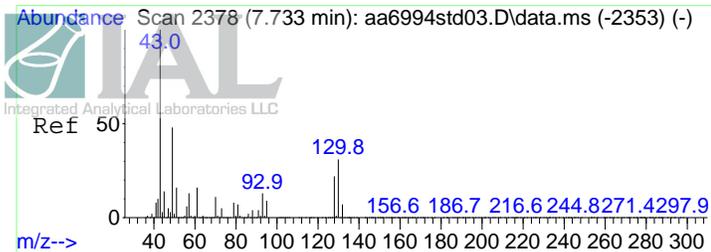
Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
Internal Standards						
1) Bromochloromethane (IS)	7.782	130	417414	10.00	ppbV	0.06
38) 1,4-Difluorobenzene (IS)	9.798	114	1445638	10.00	ppbV #	0.02
55) d-5 Chlorobenzene (IS)	15.126	117	1308945	10.00	ppbV	0.02
System Monitoring Compounds						
64) Bromofluorobenzene (tu...	17.312	95	1173967	9.74	ppbV	0.00
Target Compounds						
14) Acetone	4.901	58	105050	3.11	ppbV #	59
19) Methylene chloride	5.731	49	254918	2.62	ppbV	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\06-13-18\
Data File : aa7479.D
Acq On : 13 Jun 2018 5:27 pm
Operator : jls
Sample : E18-04189-21
Misc : dup of E18-4189-01, can # 5096
ALS Vial : 9 Sample Multiplier: 1

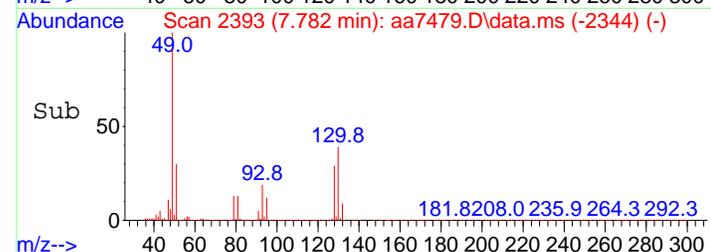
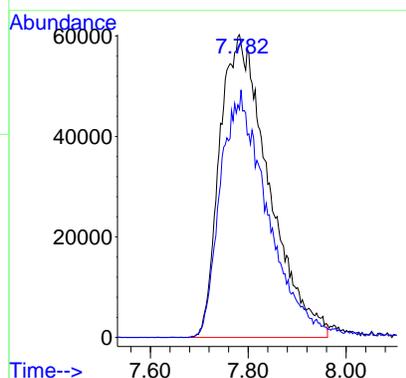
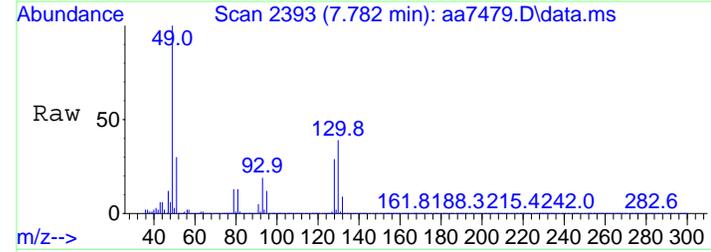
Quant Time: Jun 14 11:13:05 2018
Quant Method : C:\msdchem\1\METHODS\0518.M
Quant Title : TO-15 on the Agilent 7890A / 5975C
QLast Update : Fri May 18 13:51:08 2018
Response via : Initial Calibration





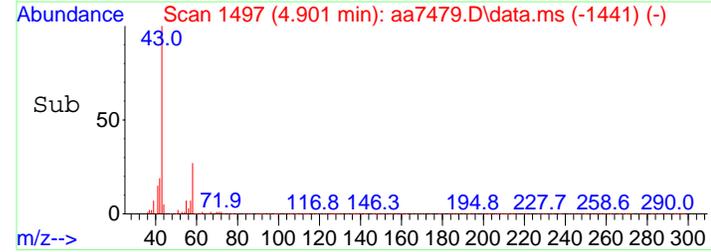
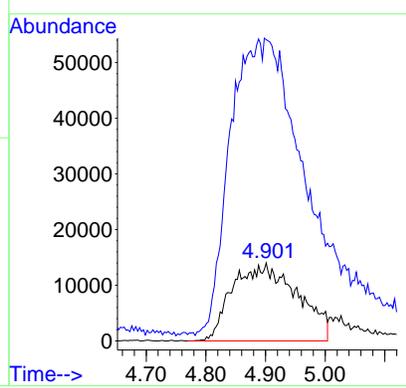
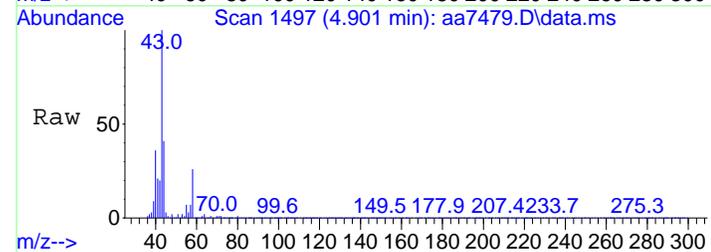
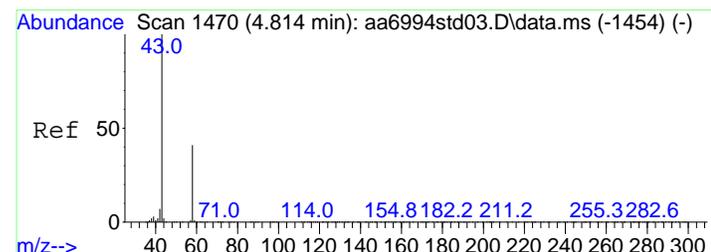
#1
 Bromochloromethane (IS)
 Concen: 10.00 ppbV
 RT: 7.782 min Scan# 2393
 Delta R.T. 0.058 min
 Lab File: aa7479.D
 Acq: 13 Jun 2018 5:27 pm

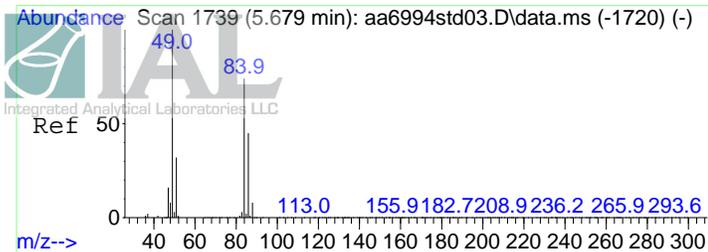
Tgt Ion: 130 Resp: 417414
 Ion Ratio Lower Upper
 130 100
 128 77.9 62.6 94.0



#14
 Acetone
 Concen: 3.11 ppbV
 RT: 4.901 min Scan# 1497
 Delta R.T. 0.081 min
 Lab File: aa7479.D
 Acq: 13 Jun 2018 5:27 pm

Tgt Ion: 58 Resp: 105050
 Ion Ratio Lower Upper
 58 100
 43 413.6 263.2 394.8#

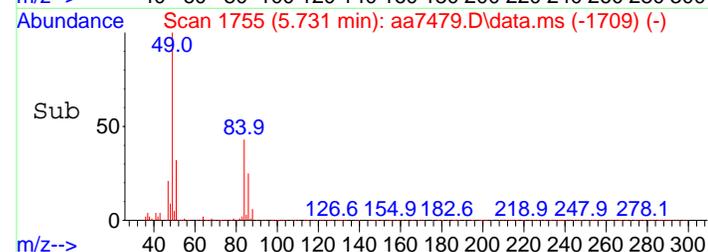
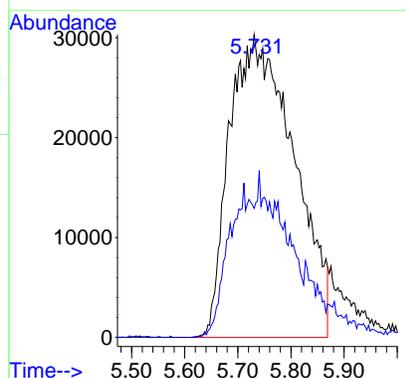
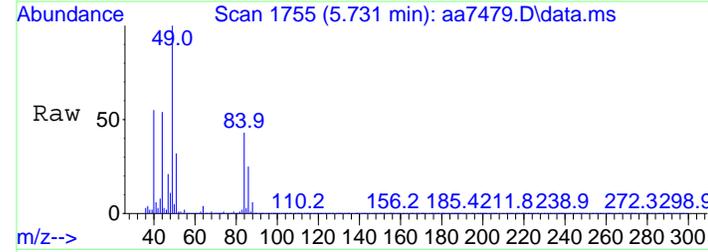




#19
 Methylene chloride
 Concen: 2.62 ppbV
 RT: 5.731 min Scan# 1755
 Delta R.T. 0.049 min
 Lab File: aa7479.D
 Acq: 13 Jun 2018 5:27 pm

Tgt Ion: 49 Resp: 254918

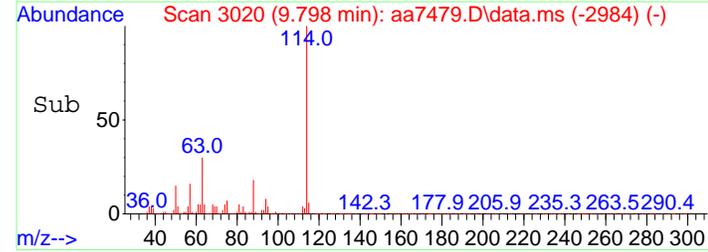
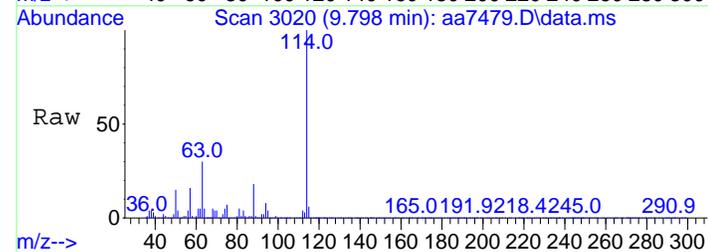
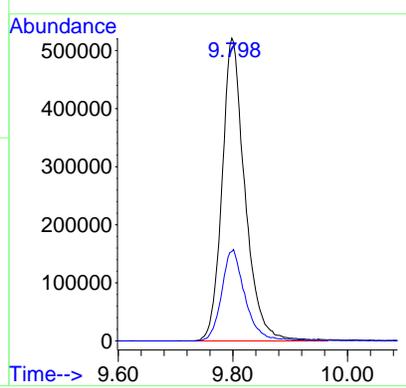
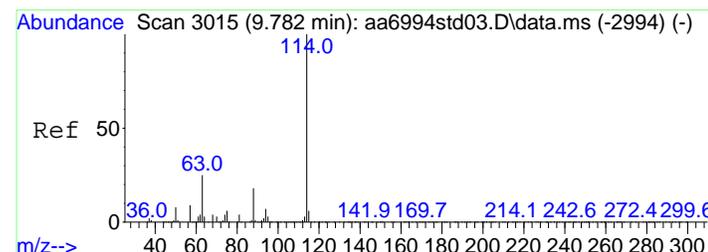
Ion	Ratio	Lower	Upper
49	100		
84	48.4	41.6	62.4

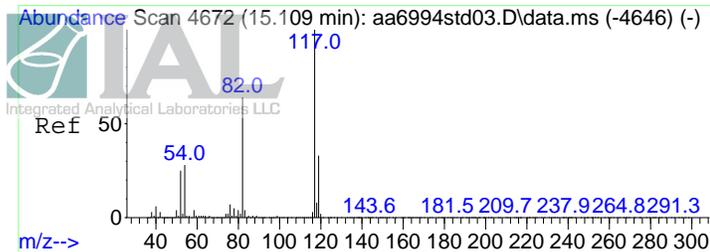


#38
 1,4-Difluorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 9.798 min Scan# 3020
 Delta R.T. 0.016 min
 Lab File: aa7479.D
 Acq: 13 Jun 2018 5:27 pm

Tgt Ion: 114 Resp: 1445638

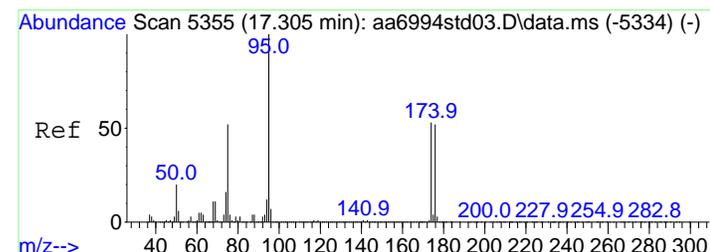
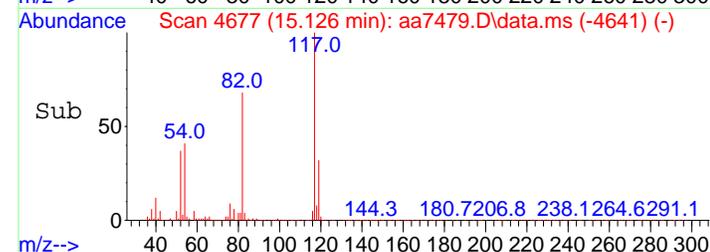
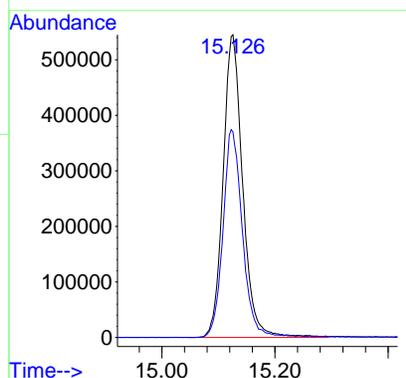
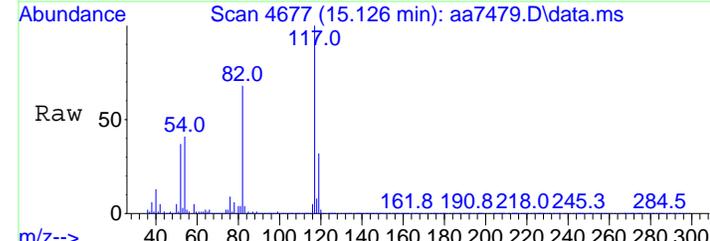
Ion	Ratio	Lower	Upper
114	100		
63	30.7	20.0	30.0#





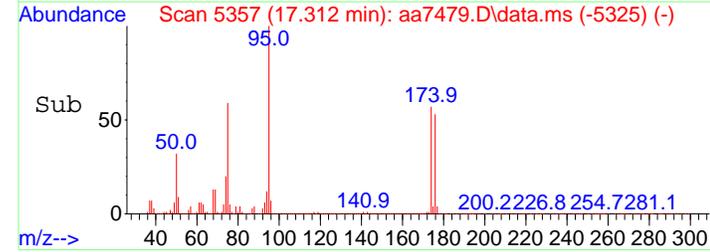
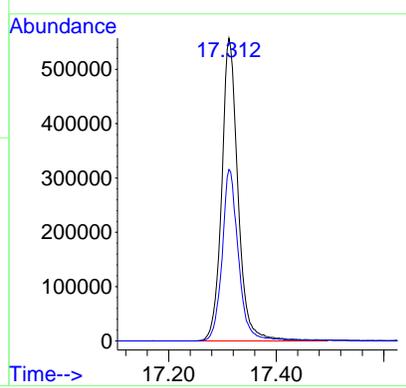
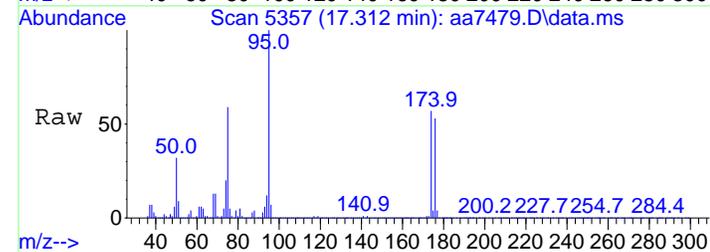
#55
 d-5 Chlorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 15.126 min Scan# 4677
 Delta R.T. 0.016 min
 Lab File: aa7479.D
 Acq: 13 Jun 2018 5:27 pm

Tgt Ion	Resp	Lower	Upper
117	1308945		
82	68.1	56.0	84.0



#64
 Bromofluorobenzene (tune std)
 Concen: 9.74 ppbV
 RT: 17.312 min Scan# 5357
 Delta R.T. 0.004 min
 Lab File: aa7479.D
 Acq: 13 Jun 2018 5:27 pm

Tgt Ion	Resp	Lower	Upper
95	1173967		
174	56.7	61.5	92.3#



Integrated Analytical Laboratories, LLC

Volatile Organic Compounds by EPA Method TO-15

Laboratory Sample Duplicate Report

SDG Number: E18-06204
 IAL Sample ID: E18-06204-01
 Matrix: Air
 Summa ID: 3280

Date Received: 8/6/18
 Date Analyzed: 8/6/18,8/6/18
 Lab Data File#: AA8195,AA8196
 Dilution Factor: 1
 Injection Volume: 500ml
 GC/MS Column: RTX-1, 0.32 mmID

Compound	CAS #	Sample E18-06204-01 Concentration Reported		Sample Dup E18-06204-21 Concentration Reported		Reporting Limits ppbv	RPD
		ppbv	Q	ppbv	Q		
Acetone	67-64-1	10		10.0		0.20	0.00%
Allyl Chloride	107-05-1		0.20 U		0.20 U	0.20	0.00%
Benzene	71-43-2	0.26		0.24		0.20	8.00%
Bromodichloromethane	75-27-4		0.20 U		0.20 U	0.20	0.00%
Bromoform	75-25-2		0.20 U		0.20 U	0.20	0.00%
Bromomethane	74-83-9		0.20 U		0.20 U	0.20	0.00%
1,3-Butadiene	106-99-0		0.20 U		0.20 U	0.20	0.00%
Chlorobenzene	108-90-7		0.20 U		0.20 U	0.20	0.00%
Chloroethane	75-00-3		0.20 U		0.20 U	0.20	0.00%
Chloroform	67-66-3		0.20 U		0.20 U	0.20	0.00%
Chloromethane	74-87-3		0.20 U		0.20 U	0.20	0.00%
Carbon disulfide	75-15-0		0.20 U		0.20 U	0.20	0.00%
Carbon tetrachloride	56-23-5		0.20 U		0.20 U	0.20	0.00%
2-Chlorotoluene	95-49-8		0.20 U		0.20 U	0.20	0.00%
Cyclohexane	110-82-7		0.20 U		0.20 U	0.20	0.00%
Dibromochloromethane	124-48-1		0.20 U		0.20 U	0.20	0.00%
1,2-Dibromoethane	106-93-4		0.20 U		0.20 U	0.20	0.00%
1,2-Dichlorobenzene	95-50-1		0.20 U		0.20 U	0.20	0.00%
1,3-Dichlorobenzene	541-73-1		0.20 U		0.20 U	0.20	0.00%
1,4-Dichlorobenzene	106-46-7		0.20 U		0.20 U	0.20	0.00%
Dichlorodifluoromethane	75-71-8		0.20 U		0.20 U	0.20	0.00%
1,1-Dichloroethane	75-34-3		0.20 U		0.20 U	0.20	0.00%
1,2-Dichloroethane	107-06-2		0.20 U		0.20 U	0.20	0.00%
1,1-Dichloroethene	75-35-4		0.20 U		0.20 U	0.20	0.00%
1,2-Dichloroethene (cis)	156-59-2		0.20 U		0.20 U	0.20	0.00%
1,2-Dichloroethene (trans)	156-60-5		0.20 U		0.20 U	0.20	0.00%
1,2-Dichloropropane	78-87-5		0.20 U		0.20 U	0.20	0.00%
1,3-Dichloropropene (cis)	10061-01-5		0.20 U		0.20 U	0.20	0.00%
1,3-Dichloropropene (trans)	10061-02-6		0.20 U		0.20 U	0.20	0.00%
1,2-Dichlorotetrafluoroethane	76-14-2		0.20 U		0.20 U	0.20	0.00%
1,4-Dioxane	123-91-1		0.20 U		0.20 U	0.20	0.00%
Ethanol	64-17-5	84	E	80	E	0.20	4.88%
Ethylbenzene	100-41-4		0.20 U		0.20 U	0.20	0.00%
4-Ethyltoluene	622-96-8		0.20 U		0.20 U	0.20	0.00%
n-Heptane	142-82-5		0.20 U		0.20 U	0.20	0.00%
1,3-Hexachlorobutadiene	87-68-3		0.20 U		0.20 U	0.20	0.00%
n-Hexane	110-54-3	0.20		0.21		0.20	-4.88%
Isopropanol	67-63-0	1.2		1.2		0.20	0.00%

Qualifir:

E=Concentration exceeds upper level of calibration range for instrument.
 D=Extra dilution required for this compound. J=Duplicate samples do not met RPD criteria.
 U=Compound ND or under reporting limit.

Integrated Analytical Laboratories, LLC

Volatile Organic Compounds by EPA Method TO-15

Laboratory Sample Duplicate Report

SDG Number: E18-06204
 IAL Sample ID: E18-06204-01
 Matrix: Air
 Summa ID: 3280

Date Received: 8/6/18
 Date Analyzed: 8/6/18,8/6/18
 Lab Data File#: AA8195,AA8196
 Dilution Factor: 1
 Injection Volume: 500ml
 GC/MS Column: RTX-1, 0.32 mmID

Compound	CAS #	Sample E18-06204-01 Concentration Reported		Sample Dup E18-06204-21 Concentration Reported		Reporting Limits ppbv	RPD
		ppbv	Q	ppbv	Q		
Methylene chloride	75-09-2	1.5		1.4		0.20	6.90%
Methyl ethyl ketone	78-93-3	1.7		1.8		0.20	-5.71%
Methyl isobutyl ketone	108-10-1		0.20 U		0.20 U	0.20	0.00%
Methyl methacrylate	80-62-6		0.20 U		0.20 U	0.20	0.00%
Methyl tert-butyl ether	1634-04-4		0.20 U		0.20 U	0.20	0.00%
Styrene	100-42-5		0.20 U		0.20 U	0.20	0.00%
Tert-butyl alcohol	75-65-0		0.20 U		0.20 U	0.20	0.00%
1,1,2,2-Tetrachloroethane	79-34-5		0.20 U		0.20 U	0.20	0.00%
Tetrachloroethene	127-18-4		0.20 U		0.20 U	0.20	0.00%
Tetrahydrofuran	109-99-9		0.20 U		0.20 U	0.20	0.00%
Toluene	108-88-3	0.23		0.23		0.20	0.00%
1,2,4-Trichlorobenzene	120-82-1		0.20 U		0.20 U	0.20	0.00%
1,1,1-Trichloroethane	71-55-6		0.20 U		0.20 U	0.20	0.00%
1,1,2-Trichloroethane	79-00-5		0.20 U		0.20 U	0.20	0.00%
Trichloroethene	79-01-6		0.20 U		0.20 U	0.20	0.00%
Trichlorofluoromethane	75-69-4	0.57		0.55		0.20	3.57%
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1		0.20 U		0.20 U	0.20	0.00%
1,2,4-Trimethylbenzene	95-63-6		0.20 U		0.20 U	0.20	0.00%
1,3,5-Trimethylbenzene	108-67-8		0.20 U		0.20 U	0.20	0.00%
2,2,4-Trimethylpentane	540-84-1		0.20 U		0.20 U	0.20	0.00%
Vinyl bromide	593-60-2		0.20 U		0.20 U	0.20	0.00%
Vinyl chloride	75-01-4		0.20 U		0.20 U	0.20	0.00%
Xylenes (m&p)	179601-23-1		0.40 U		0.40 U	0.40	0.00%
Xylenes (o)	95-47-6		0.20 U		0.20 U	0.20	0.00%

RPD must be <25% for all laboratory duplicate samples. Laboratory duplicate samples are run once daily.

NC = The RPD could not be calculated since the compound was only detected in either the parent or duplicate sample.

Qualifir:

E=Concentration exceeds upper level of calibration range for instrument.

D=Extra dilution required for this compound. J=Duplicate samples do not met RPD criteria.

U=Compound ND or under reporting limit.

Data Path : C:\DATA\08-06-18\
 Data File : aa8195.D
 Acq On : 6 Aug 2018 4:17 pm
 Operator : jls
 Sample : E18-06204-01
 Misc : 3280, 500cc
 ALS Vial : 5 Sample Multiplier: 1

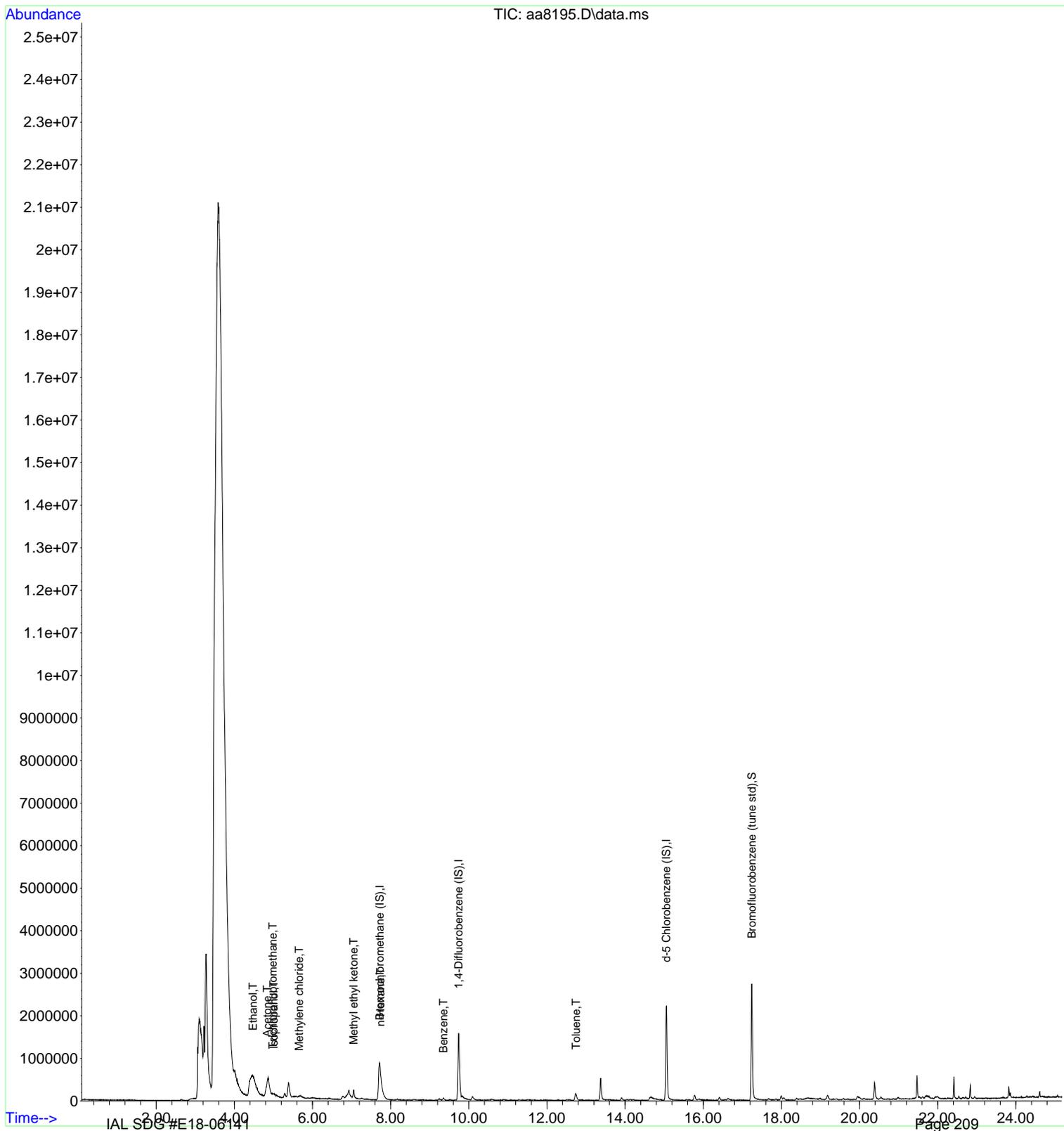
Quant Time: Aug 07 09:36:00 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) Bromochloromethane (IS)	7.724	130	395950	10.00	ppbV	0.04	
38) 1,4-Difluorobenzene (IS)	9.740	114	1565245	10.00	ppbV	# 0.00	
55) d-5 Chlorobenzene (IS)	15.055	117	1357197	10.00	ppbV	0.00	
System Monitoring Compounds							
64) Bromofluorobenzene (tu...	17.242	95	1243769	10.05	ppbV	0.00	
Target Compounds							
11) Ethanol	4.474	45	2609452	83.69	ppbV	#	35
14) Acetone	4.837	58	306980	10.18	ppbV	#	1
15) Trichlorofluoromethane	4.985	101	77186	0.57	ppbV		84
16) Isopropanol	5.011	45	114485	1.16	ppbV	#	53
19) Methylene chloride	5.657	49	116112	1.53	ppbV	#	40
27) Methyl ethyl ketone	7.052	43	191738	1.72	ppbV	#	76
30) n-Hexane	7.737	57	19229	0.20	ppbV	#	22
35) Benzene	9.351	78	44737	0.26	ppbV	#	88
50) Toluene	12.743	91	45918	0.23	ppbV		92

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\08-06-18\
 Data File : aa8195.D
 Acq On : 6 Aug 2018 4:17 pm
 Operator : jls
 Sample : E18-06204-01
 Misc : 3280, 500cc
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 07 09:36:00 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

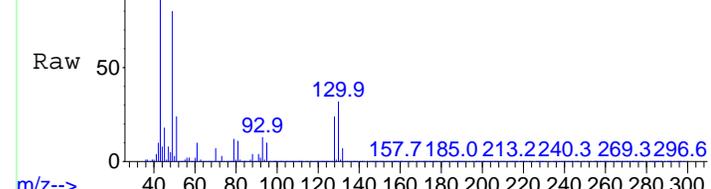


Abundance Scan 2378 (7.733 min): aa6994std03.D\data.ms (-2353) (-)



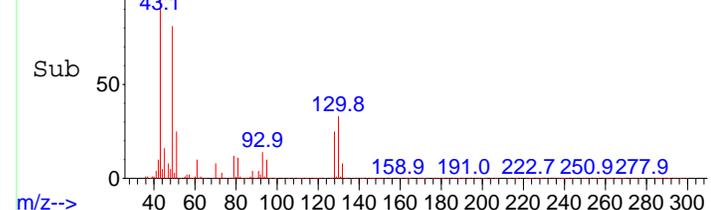
m/z-->

Abundance Scan 2375 (7.724 min): aa8195.D\data.ms



m/z-->

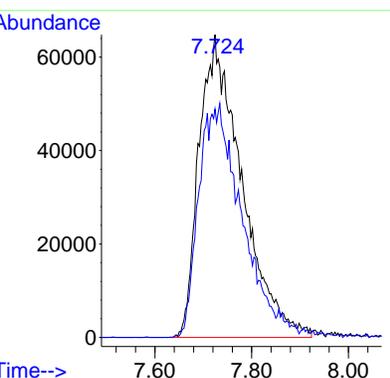
Abundance Scan 2375 (7.724 min): aa8195.D\data.ms (-2333) (-)



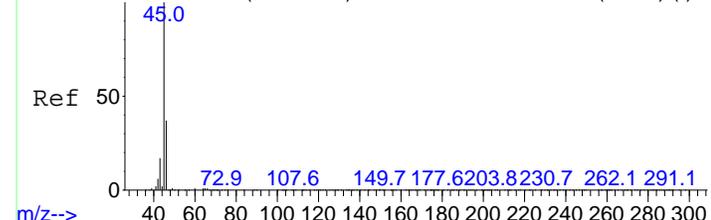
m/z-->

#1
Bromochloromethane (IS)
Concen: 10.00 ppbV
RT: 7.724 min Scan# 2375
Delta R.T. 0.035 min
Lab File: aa8195.D
Acq: 6 Aug 2018 4:17 pm

Tgt Ion	Resp	Lower	Upper
130	395950		
130	100		
128	78.4	62.6	94.0

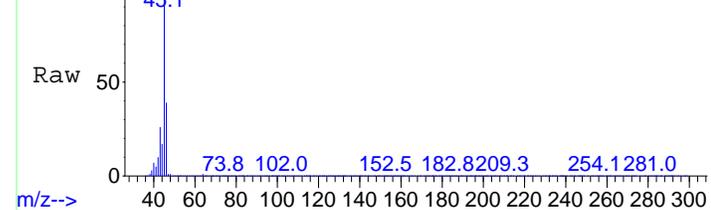


Abundance Scan 1341 (4.399 min): aa6994std03.D\data.ms (-1322) (-)



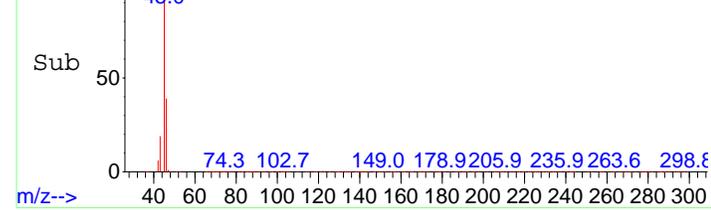
m/z-->

Abundance Scan 1364 (4.474 min): aa8195.D\data.ms



m/z-->

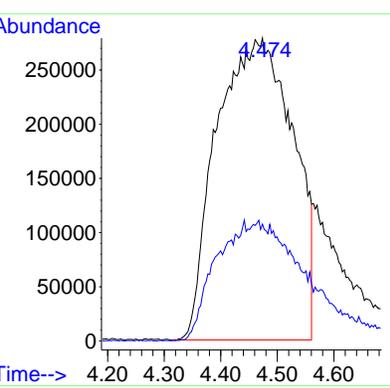
Abundance Scan 1364 (4.474 min): aa8195.D\data.ms (-1303) (-)



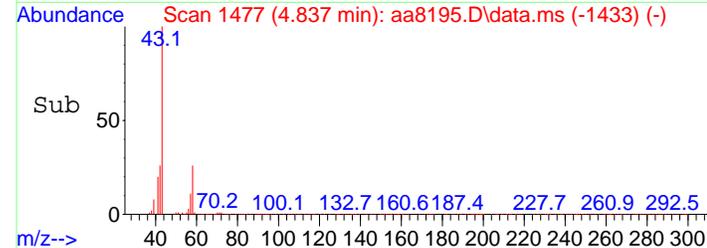
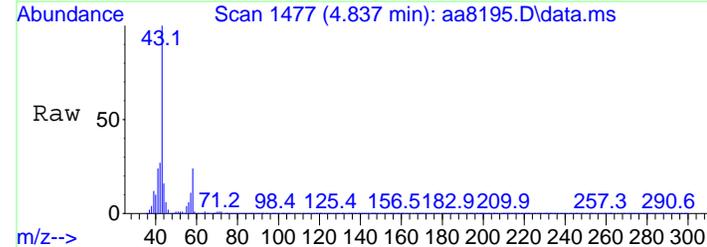
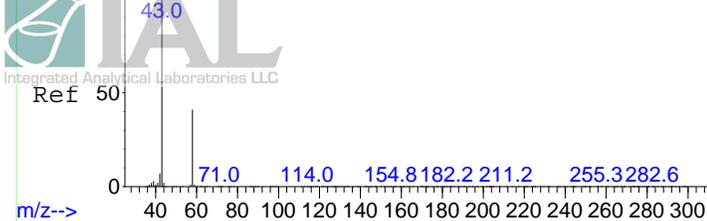
m/z-->

#11
Ethanol
Concen: 83.69 ppbV
RT: 4.474 min Scan# 1364
Delta R.T. 0.097 min
Lab File: aa8195.D
Acq: 6 Aug 2018 4:17 pm

Tgt Ion	Resp	Lower	Upper
45	2609452		
45	100		
46	0.0	32.6	48.8#

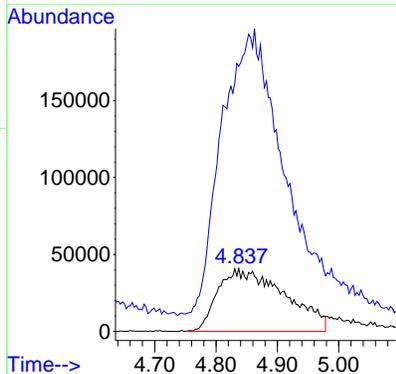


Abundance Scan 1470 (4.814 min): aa6994std03.D\data.ms (-1454) (-)

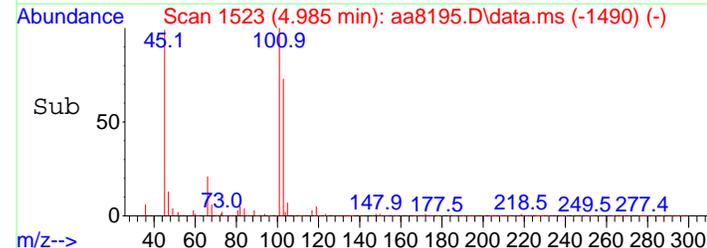
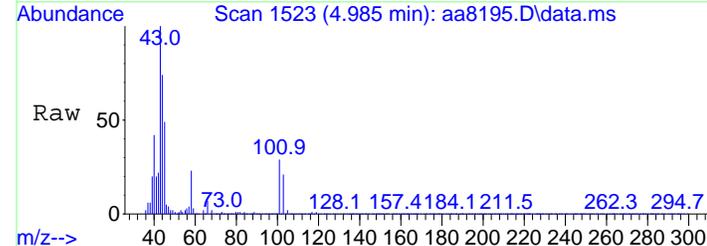
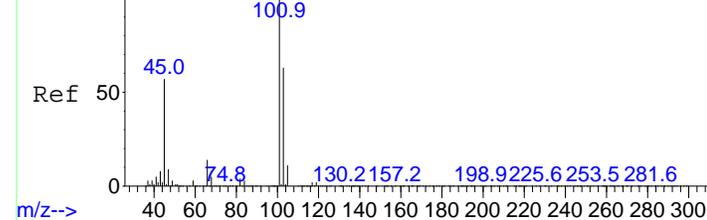


#14
Acetone
Concen: 10.18 ppbV
RT: 4.837 min Scan# 1477
Delta R.T. 0.042 min
Lab File: aa8195.D
Acq: 6 Aug 2018 4:17 pm

Tgt Ion: 58 Resp: 306980
Ion Ratio Lower Upper
58 100
43 0.0 263.2 394.8#

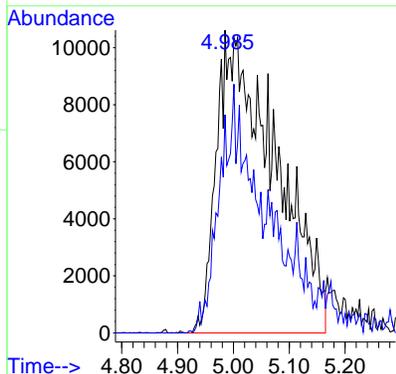


Abundance Scan 1528 (5.000 min): aa6994std03.D\data.ms (-1502) (-)



#15
Trichlorofluoromethane
Concen: 0.57 ppbV
RT: 4.985 min Scan# 1523
Delta R.T. 0.007 min
Lab File: aa8195.D
Acq: 6 Aug 2018 4:17 pm

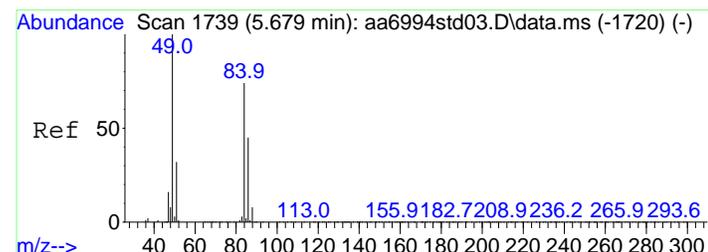
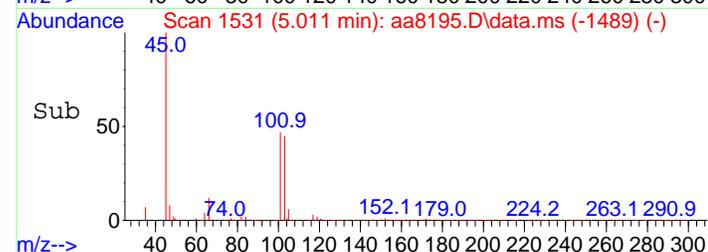
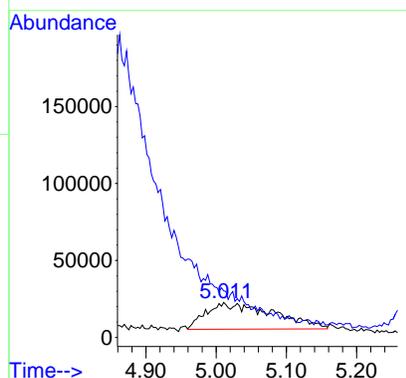
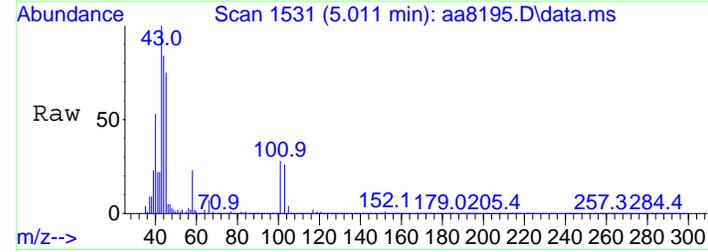
Tgt Ion: 101 Resp: 77186
Ion Ratio Lower Upper
101 100
103 52.3 52.0 78.0





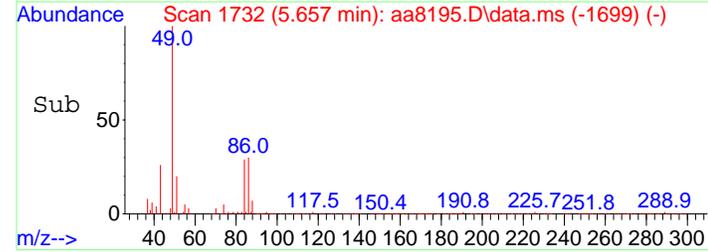
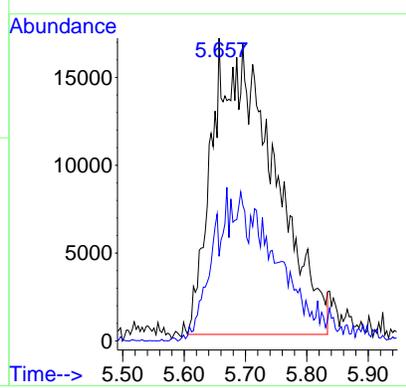
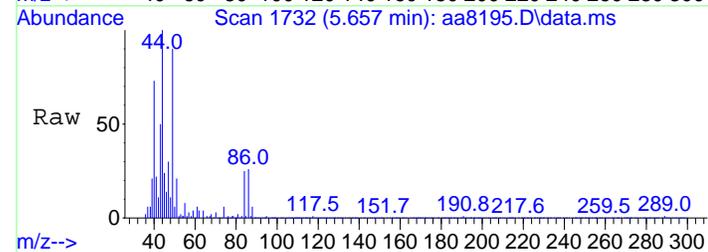
#16
 Isopropanol
 Concen: 1.16 ppbV
 RT: 5.011 min Scan# 1531
 Delta R.T. 0.036 min
 Lab File: aa8195.D
 Acq: 6 Aug 2018 4:17 pm

Tgt Ion	Resp	Lower	Upper
45	114485		
43	100	0.0	18.4
43	0.0	18.4	27.6#



#19
 Methylene chloride
 Concen: 1.53 ppbV
 RT: 5.657 min Scan# 1732
 Delta R.T. 0.007 min
 Lab File: aa8195.D
 Acq: 6 Aug 2018 4:17 pm

Tgt Ion	Resp	Lower	Upper
49	116112		
49	100		
84	9.6	41.6	62.4#

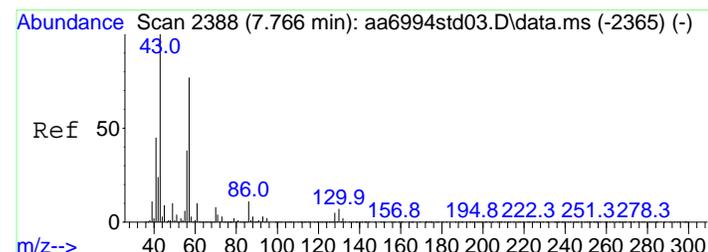
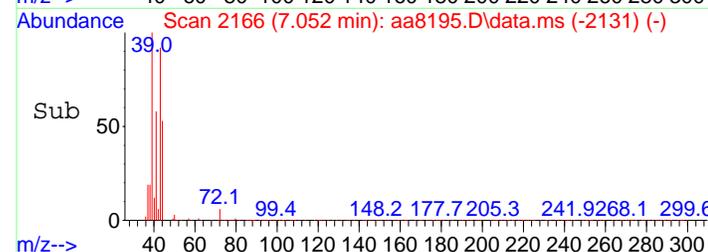
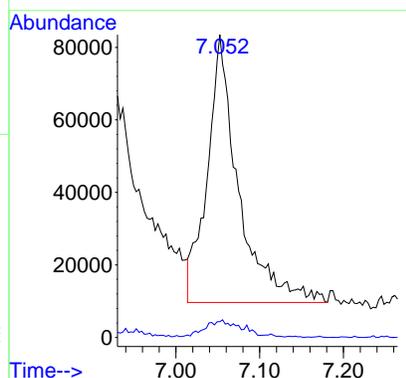
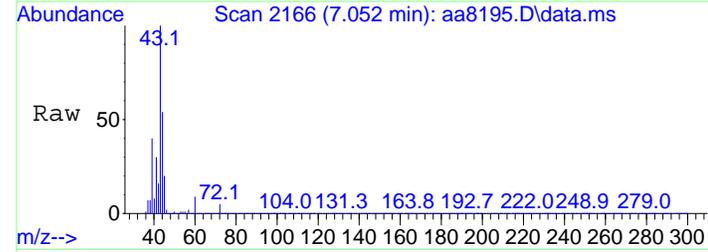




#27
 Methyl ethyl ketone
 Concen: 1.72 ppbV
 RT: 7.052 min Scan# 2166
 Delta R.T. 0.013 min
 Lab File: aa8195.D
 Acq: 6 Aug 2018 4:17 pm

Tgt Ion: 43 Resp: 191738

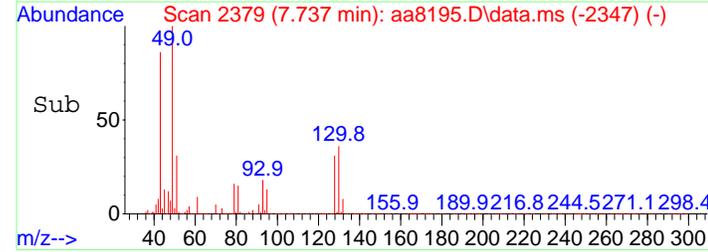
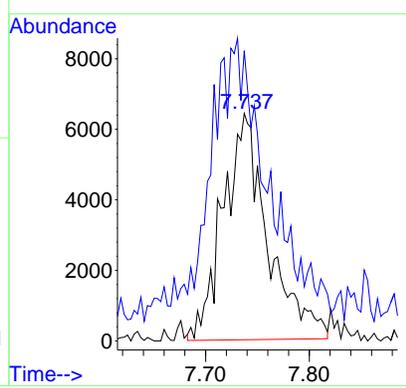
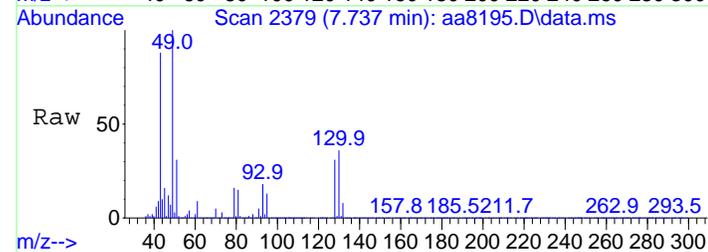
Ion	Ratio	Lower	Upper
43	100		
72	7.4	14.4	21.6#

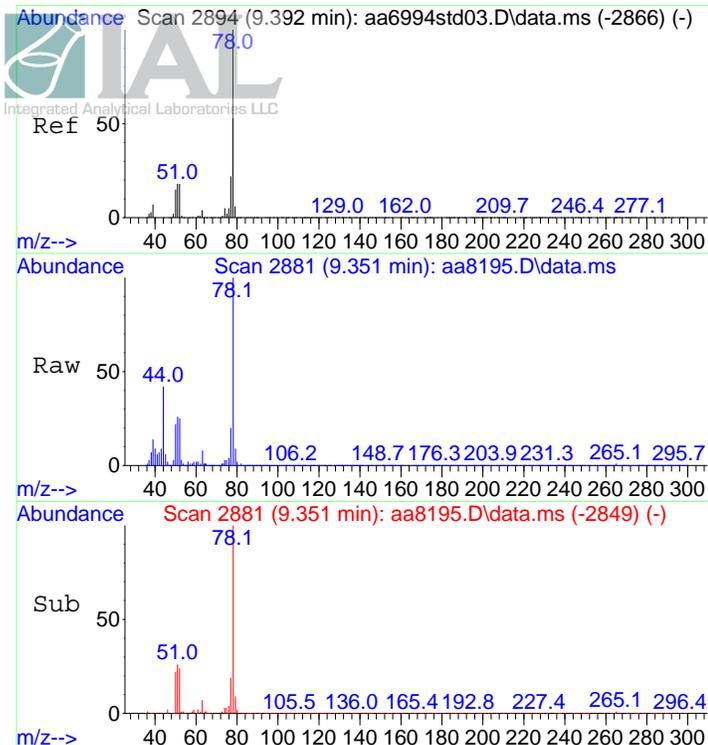


#30
 n-Hexane
 Concen: 0.20 ppbV
 RT: 7.737 min Scan# 2379
 Delta R.T. 0.003 min
 Lab File: aa8195.D
 Acq: 6 Aug 2018 4:17 pm

Tgt Ion: 57 Resp: 19229

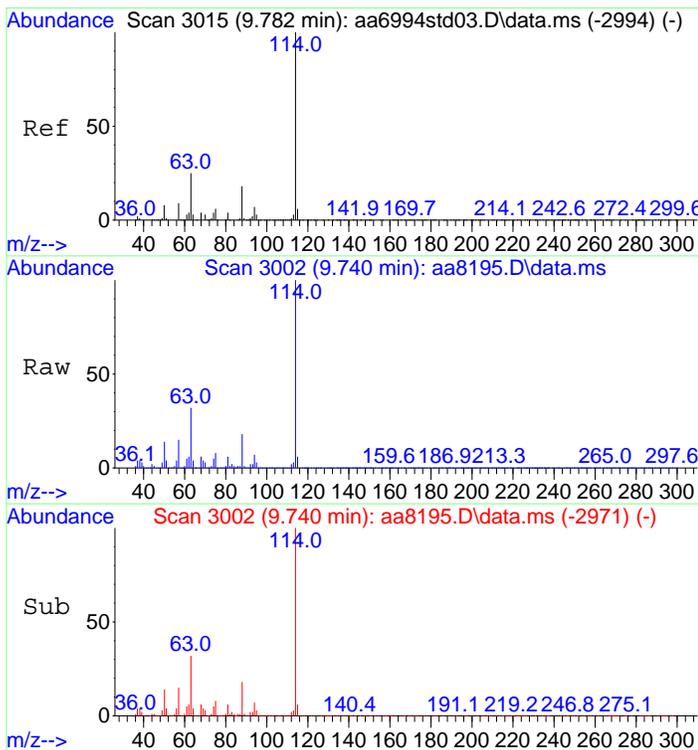
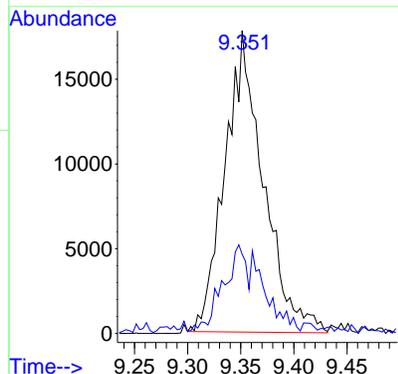
Ion	Ratio	Lower	Upper
57	100		
41	154.6	67.2	100.8#





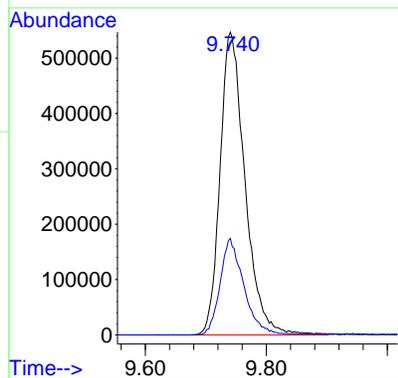
#35
Benzene
Concen: 0.26 ppbV
RT: 9.351 min Scan# 2881
Delta R.T. 0.003 min
Lab File: aa8195.D
Acq: 6 Aug 2018 4:17 pm

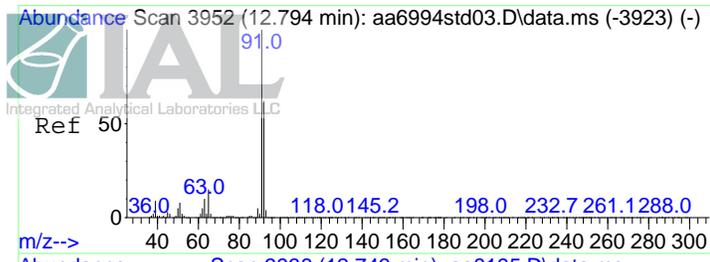
Tgt Ion	Resp	Lower	Upper
78	100		
51	17.8	19.2	28.8#



#38
1,4-Difluorobenzene (IS)
Concen: 10.00 ppbV
RT: 9.740 min Scan# 3002
Delta R.T. 0.000 min
Lab File: aa8195.D
Acq: 6 Aug 2018 4:17 pm

Tgt Ion	Resp	Lower	Upper
114	100		
63	30.5	20.0	30.0#

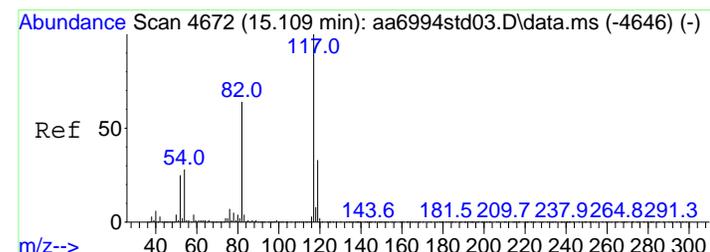
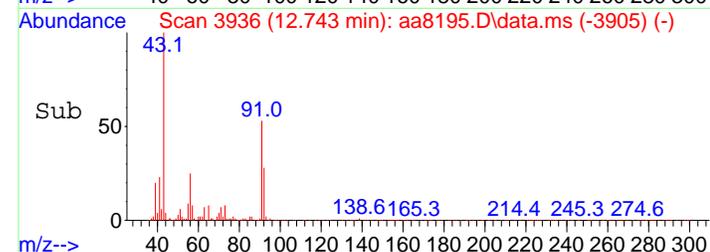
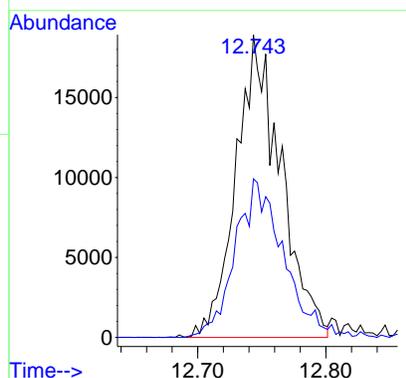
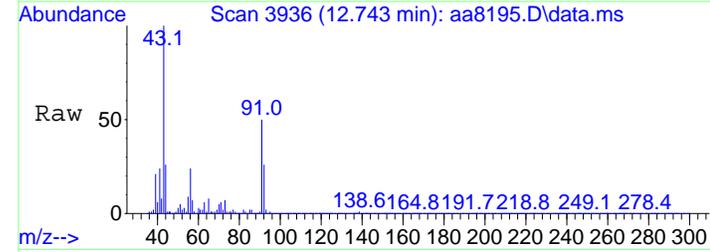




#50
 Toluene
 Concen: 0.23 ppbV
 RT: 12.743 min Scan# 3936
 Delta R.T. 0.000 min
 Lab File: aa8195.D
 Acq: 6 Aug 2018 4:17 pm

Tgt Ion: 91 Resp: 45918

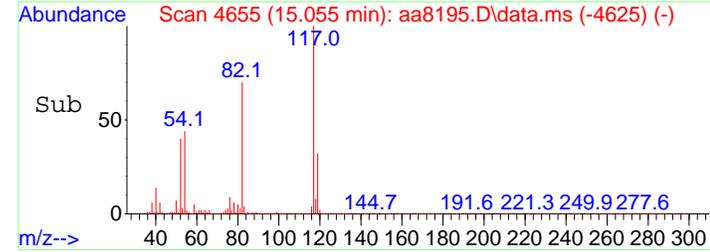
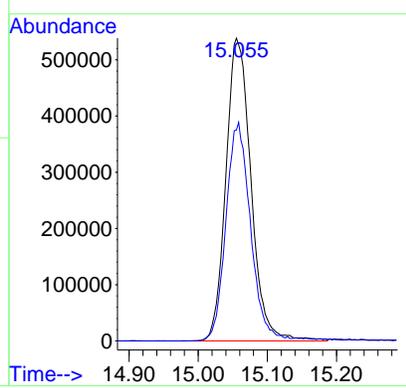
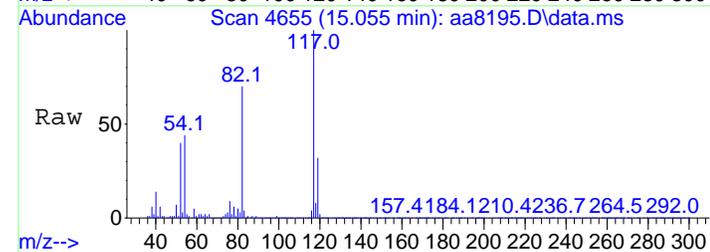
Ion	Ratio	Lower	Upper
91	100		
92	56.1	50.0	75.0



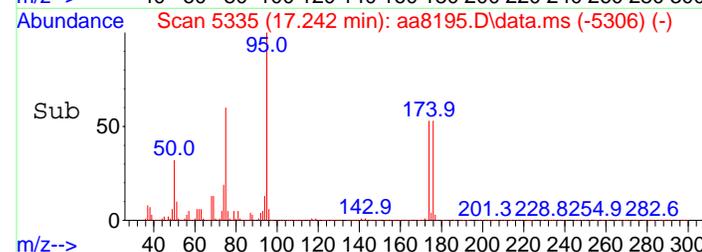
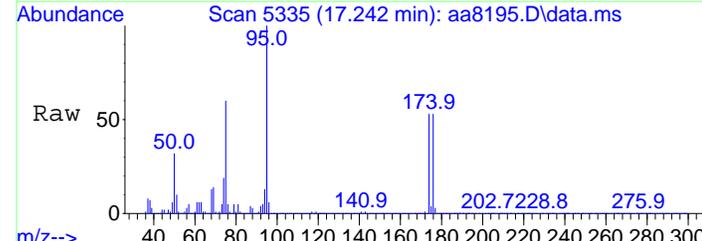
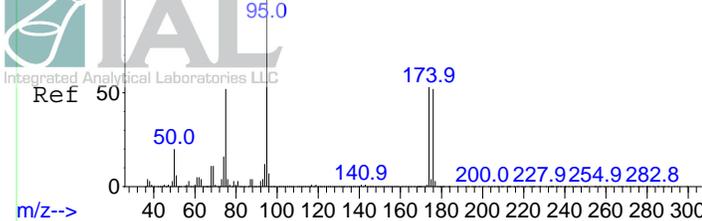
#55
 d-5 Chlorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 15.055 min Scan# 4655
 Delta R.T. -0.003 min
 Lab File: aa8195.D
 Acq: 6 Aug 2018 4:17 pm

Tgt Ion: 117 Resp: 1357197

Ion	Ratio	Lower	Upper
117	100		
82	71.2	56.0	84.0

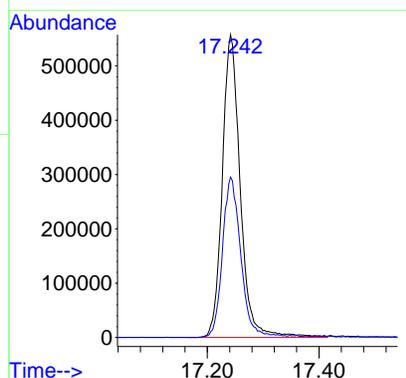


Abundance Scan 5355 (17.305 min): aa6994std03.D\data.ms (-5334) (-)



#64
Bromofluorobenzene (tune std)
Concen: 10.05 ppbV
RT: 17.242 min Scan# 5335
Delta R.T. -0.006 min
Lab File: aa8195.D
Acq: 6 Aug 2018 4:17 pm

Tgt Ion	Resp	Lower	Upper
95	1243769		
95	100		
174	54.0	61.5	92.3#



Data Path : C:\DATA\08-06-18\
 Data File : aa8196.D
 Acq On : 6 Aug 2018 4:50 pm
 Operator : jls
 Sample : E18-06204-21
 Misc : dup of E18-06204-01, can # 3280
 ALS Vial : 6 Sample Multiplier: 1

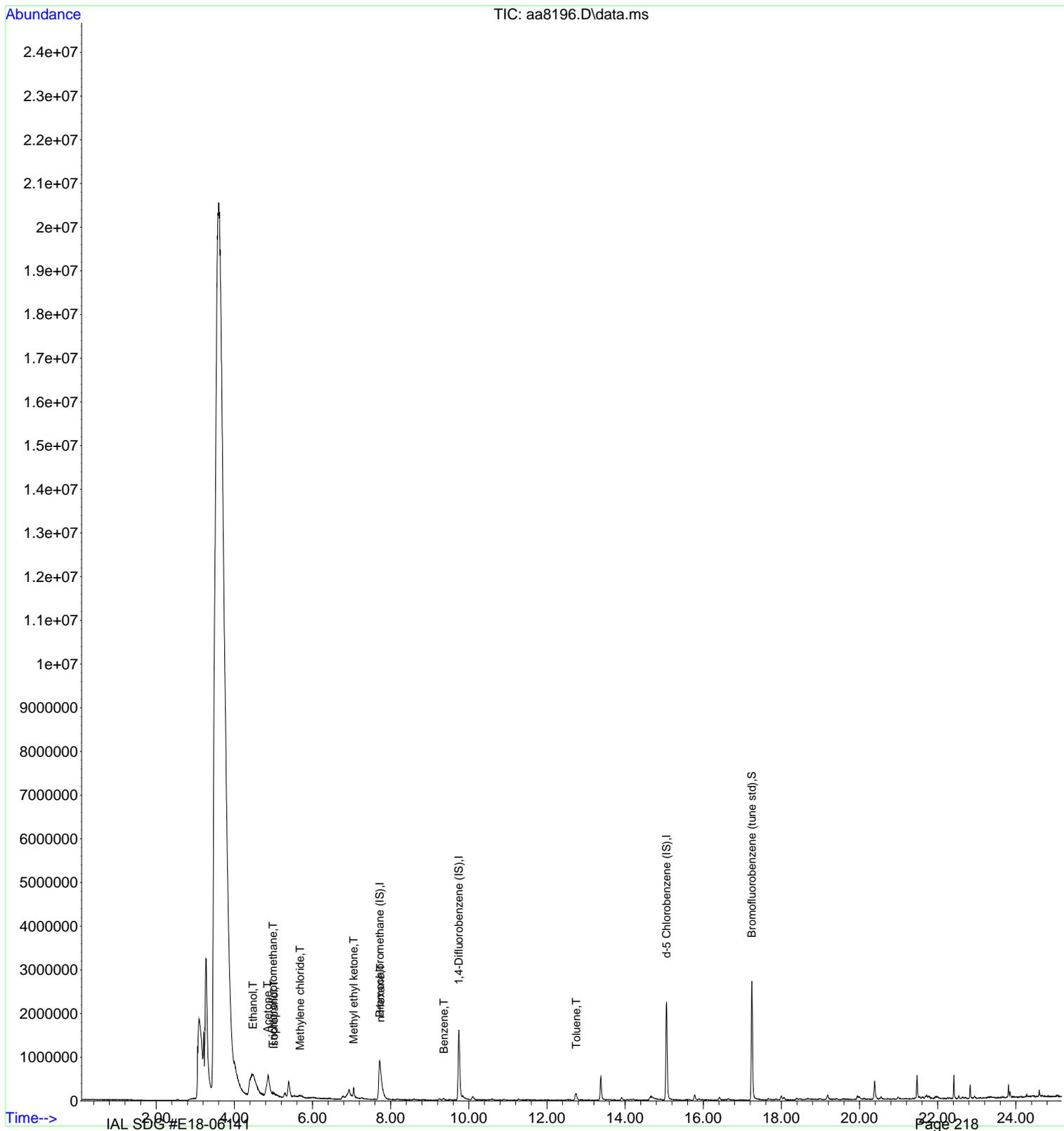
Quant Time: Aug 07 09:42:13 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

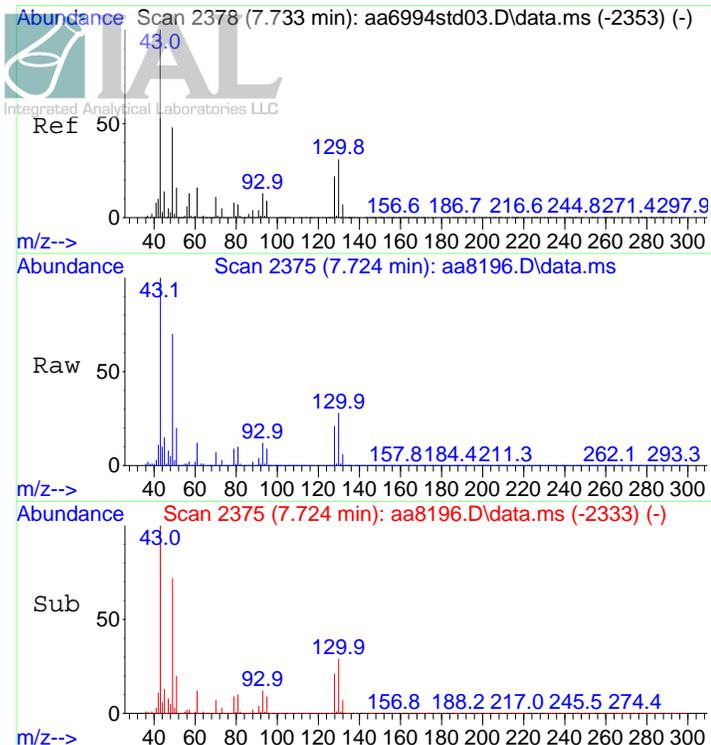
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Bromochloromethane (IS)	7.724	130	410636	10.00	ppbV	0.04
38) 1,4-Difluorobenzene (IS)	9.743	114	1600414	10.00	ppbV	0.00
55) d-5 Chlorobenzene (IS)	15.061	117	1372203	10.00	ppbV	0.00
System Monitoring Compounds						
64) Bromofluorobenzene (tu...	17.241	95	1243695	9.94	ppbV	0.00
Target Compounds						
11) Ethanol	4.480	45	2591610	80.15	ppbV #	35
14) Acetone	4.853	58	312445	9.99	ppbV #	61
15) Trichlorofluoromethane	4.991	101	77441	0.55	ppbV	89
16) Isopropanol	5.014	45	125232	1.23	ppbV #	53
19) Methylene chloride	5.682	49	113349	1.44	ppbV	87
27) Methyl ethyl ketone	7.052	43	207474	1.79	ppbV #	73
30) n-Hexane	7.737	57	20657	0.21	ppbV #	77
35) Benzene	9.361	78	43169	0.24	ppbV	94
50) Toluene	12.750	91	46925	0.23	ppbV	89

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\08-06-18\
 Data File : aa8196.D
 Acq On : 6 Aug 2018 4:50 pm
 Operator : jls
 Sample : E18-06204-21
 Misc : dup of E18-06204-01, can # 3280
 ALS Vial : 6 Sample Multiplier: 1

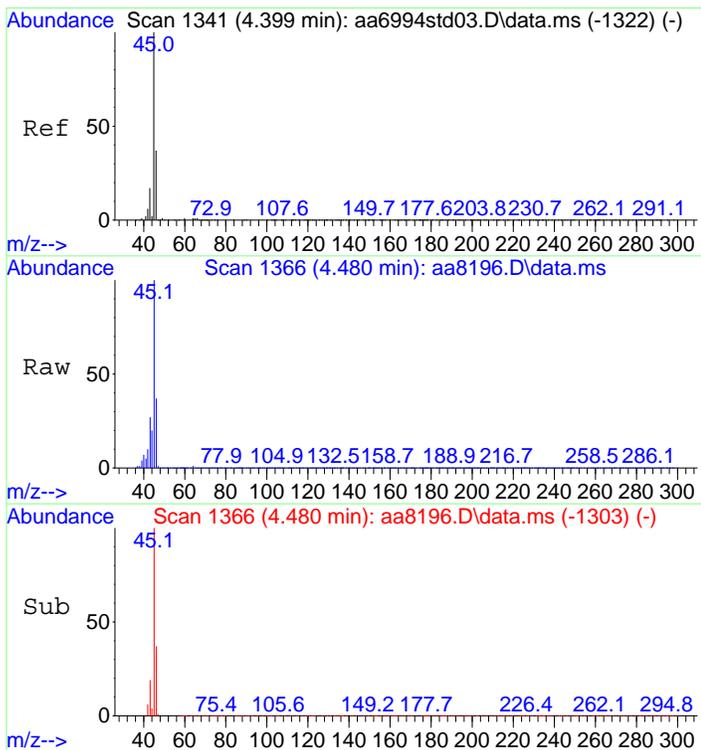
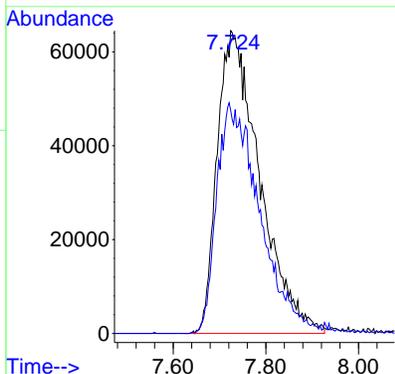
Quant Time: Aug 07 09:42:13 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration





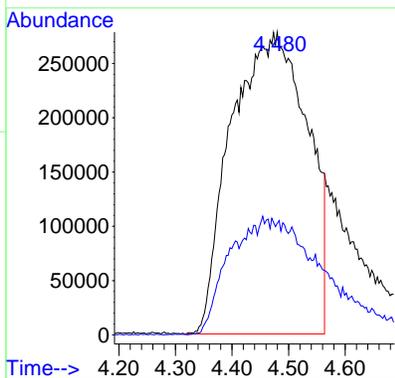
#1
 Bromochloromethane (IS)
 Concen: 10.00 ppbV
 RT: 7.724 min Scan# 2375
 Delta R.T. 0.035 min
 Lab File: aa8196.D
 Acq: 6 Aug 2018 4:50 pm

Tgt Ion	Resp	Lower	Upper
130	410636		
130	100		
128	76.8	62.6	94.0

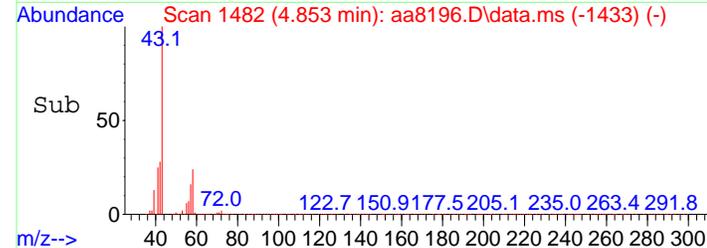
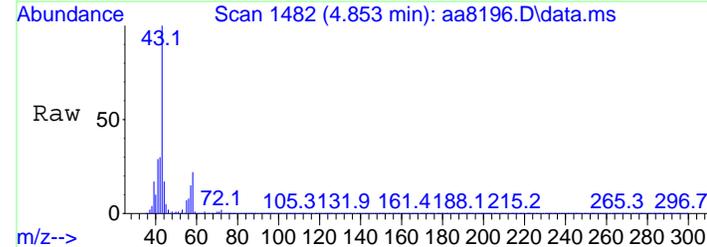
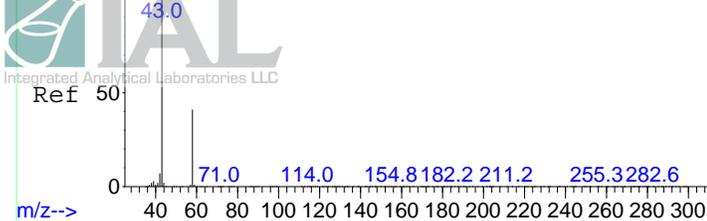


#11
 Ethanol
 Concen: 80.15 ppbV
 RT: 4.480 min Scan# 1366
 Delta R.T. 0.103 min
 Lab File: aa8196.D
 Acq: 6 Aug 2018 4:50 pm

Tgt Ion	Resp	Lower	Upper
45	2591610		
45	100		
46	0.0	32.6	48.8#

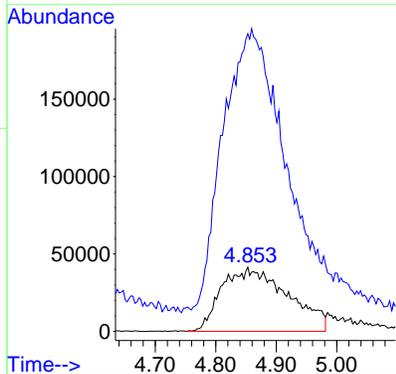


Abundance Scan 1470 (4.814 min): aa6994std03.D\data.ms (-1454) (-)

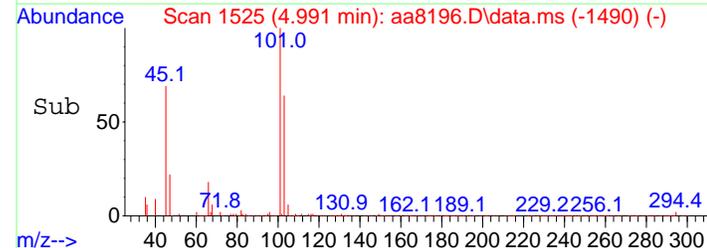
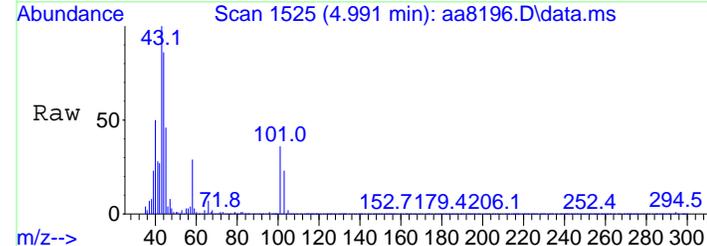
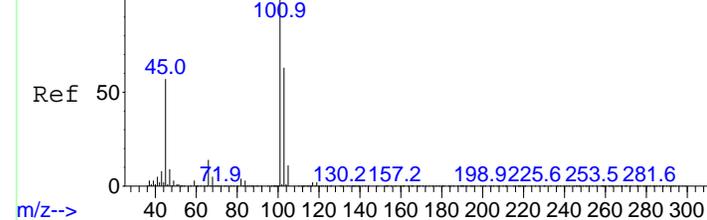


#14
Acetone
Concen: 9.99 ppbV
RT: 4.853 min Scan# 1482
Delta R.T. 0.058 min
Lab File: aa8196.D
Acq: 6 Aug 2018 4:50 pm

Tgt Ion: 58 Resp: 312445
Ion Ratio Lower Upper
58 100
43 410.7 263.2 394.8#

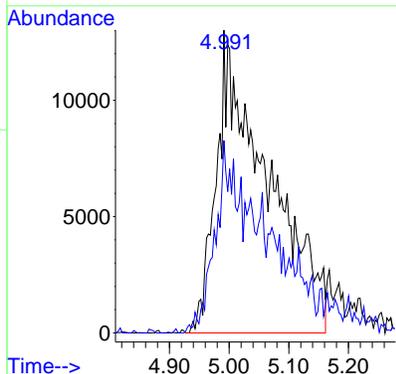


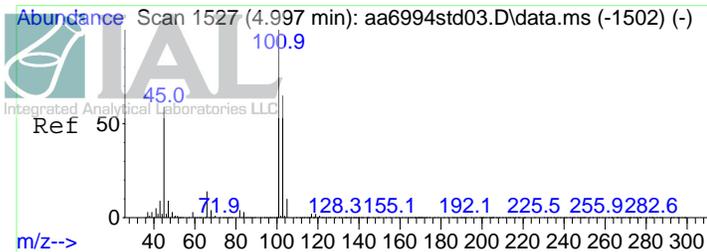
Abundance Scan 1528 (5.000 min): aa6994std03.D\data.ms (-1502) (-)



#15
Trichlorofluoromethane
Concen: 0.55 ppbV
RT: 4.991 min Scan# 1525
Delta R.T. 0.013 min
Lab File: aa8196.D
Acq: 6 Aug 2018 4:50 pm

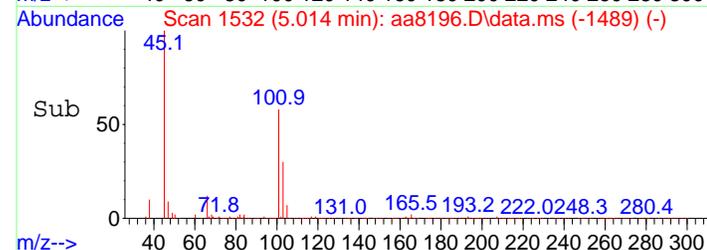
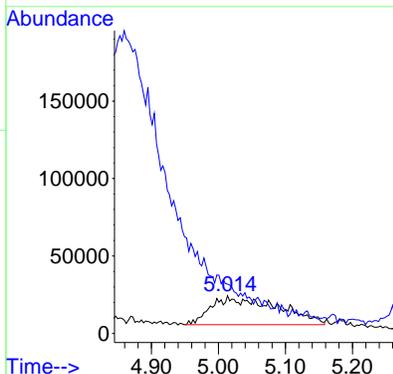
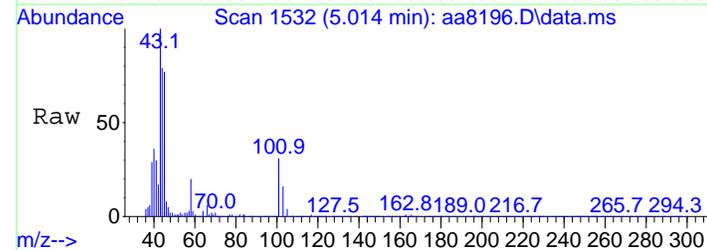
Tgt Ion: 101 Resp: 77441
Ion Ratio Lower Upper
101 100
103 56.1 52.0 78.0





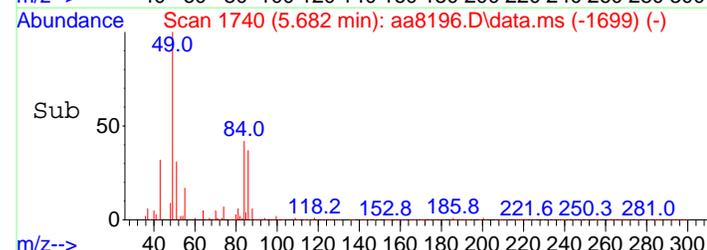
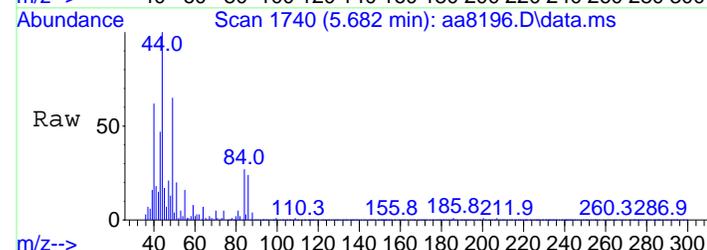
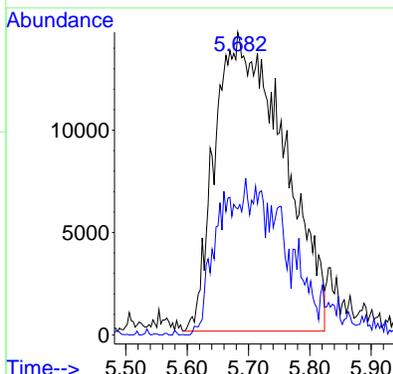
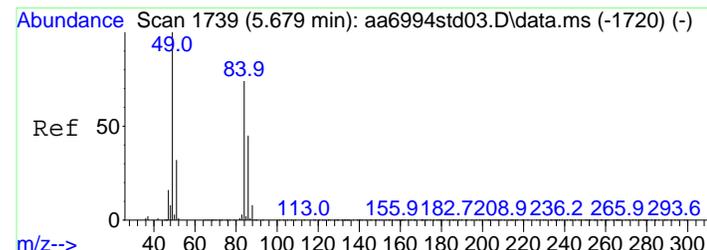
#16
 Isopropanol
 Concen: 1.23 ppbV
 RT: 5.014 min Scan# 1532
 Delta R.T. 0.039 min
 Lab File: aa8196.D
 Acq: 6 Aug 2018 4:50 pm

Tgt Ion: 45 Resp: 125232
 Ion Ratio Lower Upper
 45 100
 43 0.0 18.4 27.6#



#19
 Methylene chloride
 Concen: 1.44 ppbV
 RT: 5.682 min Scan# 1740
 Delta R.T. 0.032 min
 Lab File: aa8196.D
 Acq: 6 Aug 2018 4:50 pm

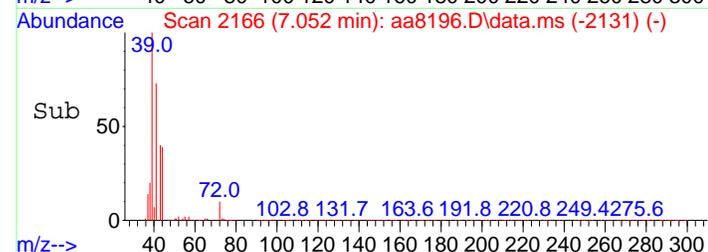
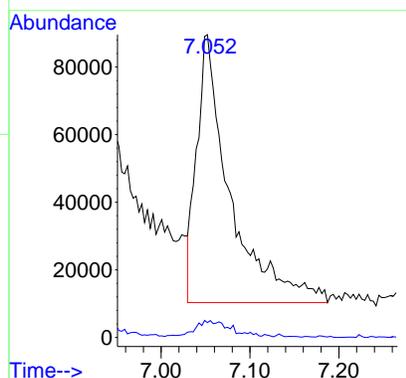
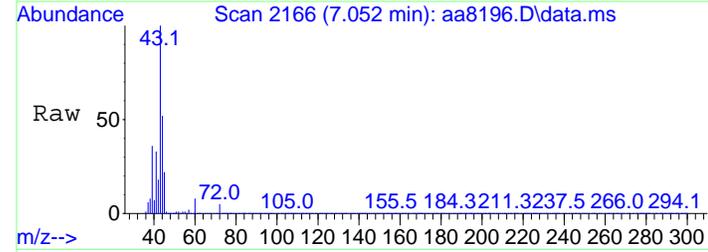
Tgt Ion: 49 Resp: 113349
 Ion Ratio Lower Upper
 49 100
 84 43.1 41.6 62.4





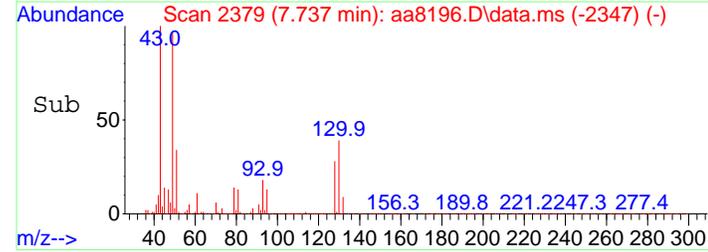
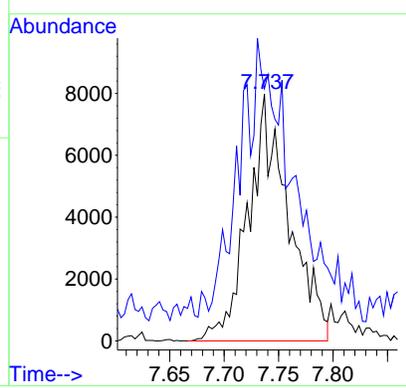
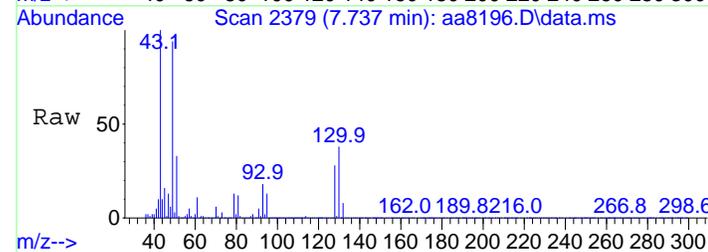
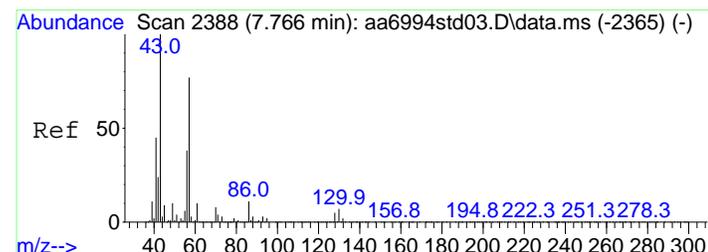
#27
Methyl ethyl ketone
Concen: 1.79 ppbV
RT: 7.052 min Scan# 2166
Delta R.T. 0.013 min
Lab File: aa8196.D
Acq: 6 Aug 2018 4:50 pm

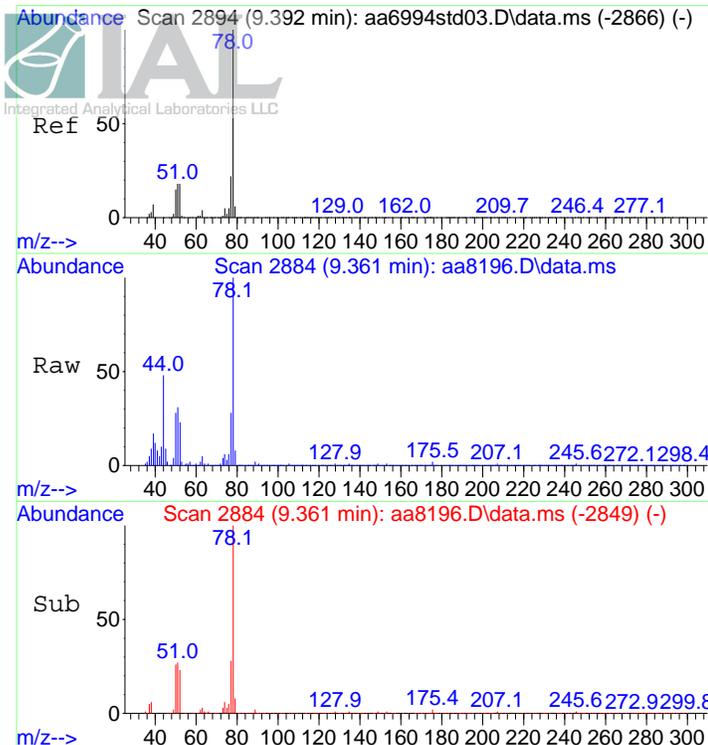
Tgt Ion: 43 Resp: 207474
Ion Ratio Lower Upper
43 100
72 6.0 14.4 21.6#



#30
n-Hexane
Concen: 0.21 ppbV
RT: 7.737 min Scan# 2379
Delta R.T. 0.003 min
Lab File: aa8196.D
Acq: 6 Aug 2018 4:50 pm

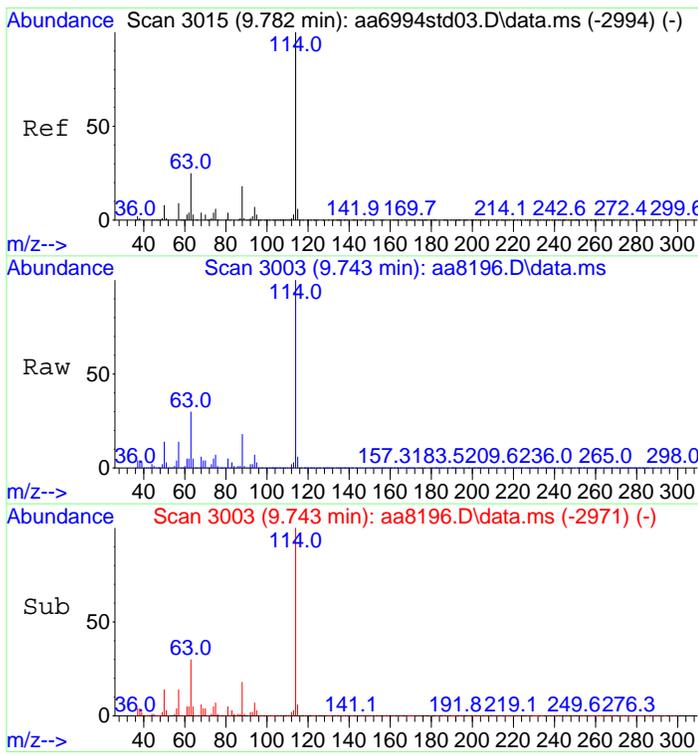
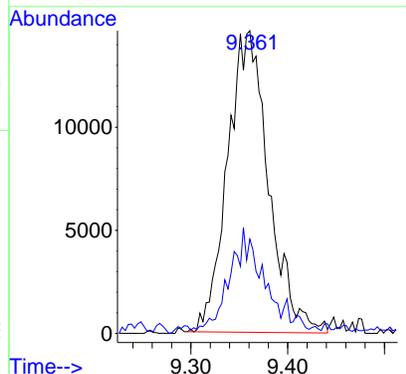
Tgt Ion: 57 Resp: 20657
Ion Ratio Lower Upper
57 100
41 105.1 67.2 100.8#





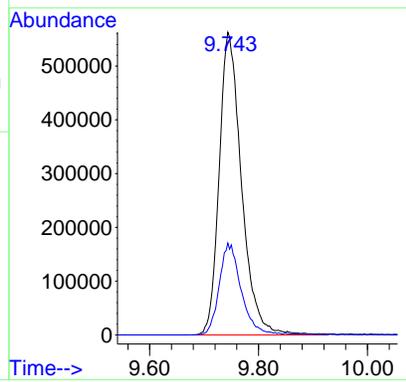
#35
Benzene
Concen: 0.24 ppbV
RT: 9.361 min Scan# 2884
Delta R.T. 0.013 min
Lab File: aa8196.D
Acq: 6 Aug 2018 4:50 pm

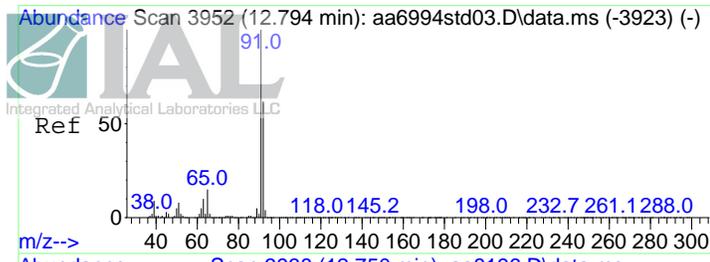
Tgt Ion	Resp	Lower	Upper
78	43169		
78	100		
51	27.1	19.2	28.8



#38
1,4-Difluorobenzene (IS)
Concen: 10.00 ppbV
RT: 9.743 min Scan# 3003
Delta R.T. 0.003 min
Lab File: aa8196.D
Acq: 6 Aug 2018 4:50 pm

Tgt Ion	Resp	Lower	Upper
114	1600414		
114	100		
63	29.6	20.0	30.0

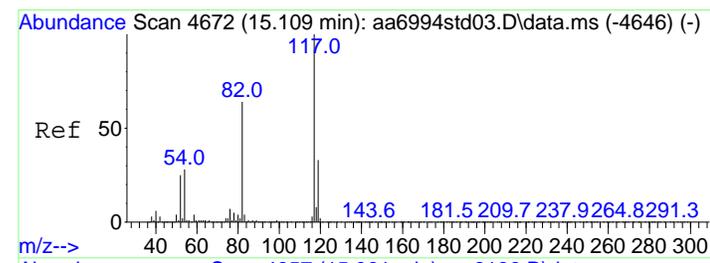
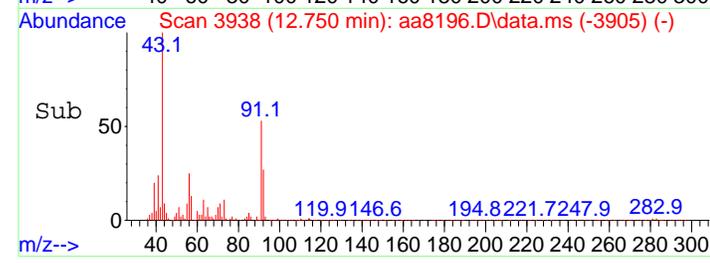
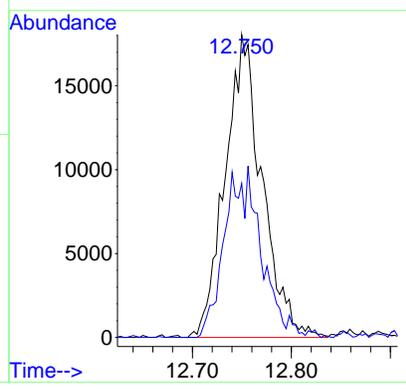
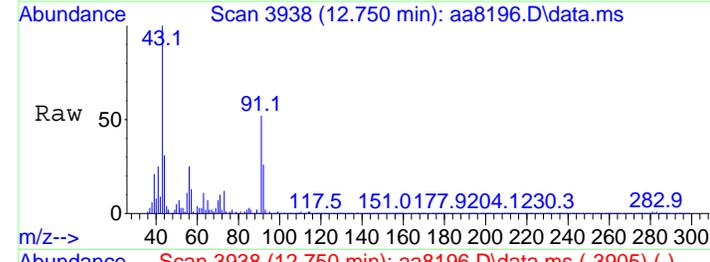




#50
 Toluene
 Concen: 0.23 ppbV
 RT: 12.750 min Scan# 3938
 Delta R.T. 0.006 min
 Lab File: aa8196.D
 Acq: 6 Aug 2018 4:50 pm

Tgt Ion: 91 Resp: 46925

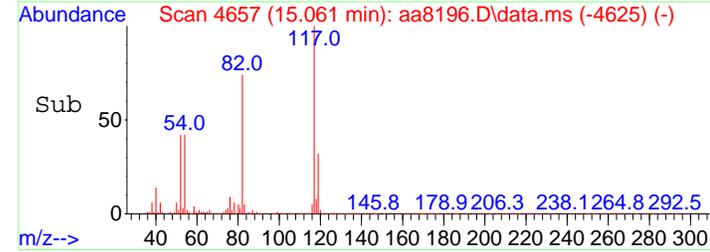
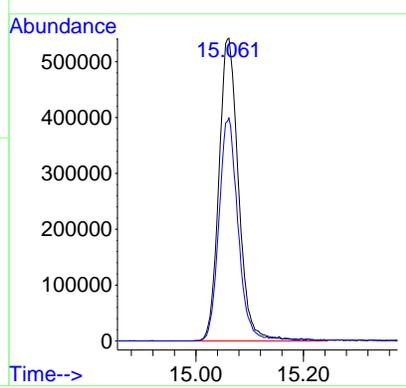
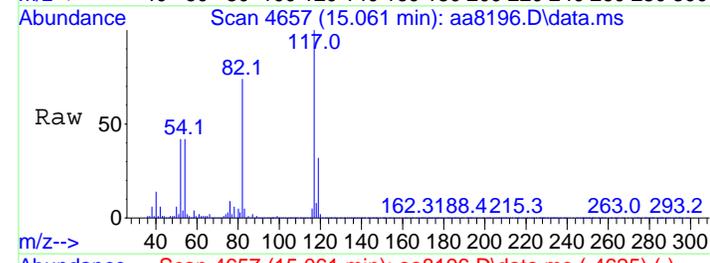
Ion	Ratio	Lower	Upper
91	100		
92	54.0	50.0	75.0



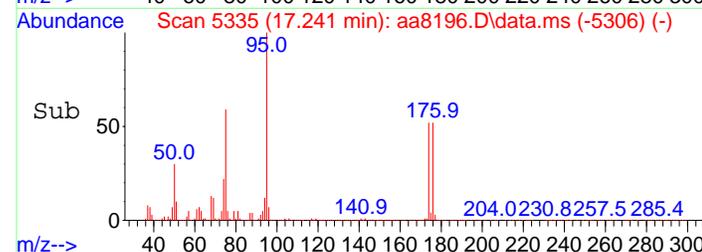
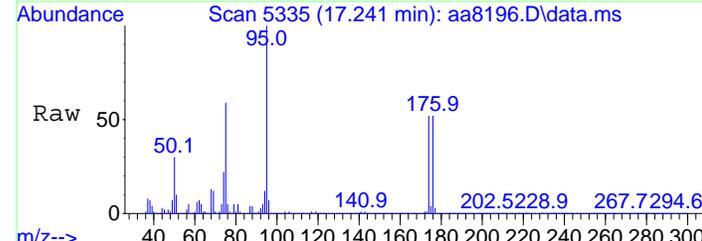
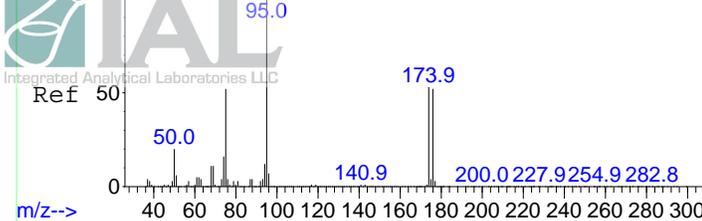
#55
 d-5 Chlorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 15.061 min Scan# 4657
 Delta R.T. 0.003 min
 Lab File: aa8196.D
 Acq: 6 Aug 2018 4:50 pm

Tgt Ion: 117 Resp: 1372203

Ion	Ratio	Lower	Upper
117	100		
82	72.0	56.0	84.0

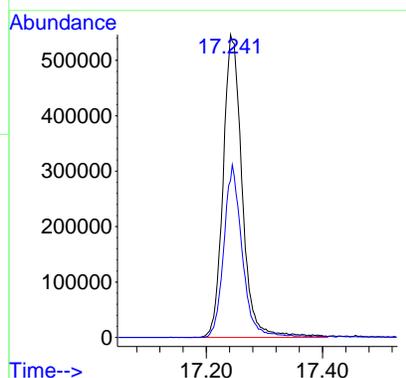


Abundance Scan 5355 (17.305 min): aa6994std03.D\data.ms (-5334) (-)



#64
Bromofluorobenzene (tune std)
Concen: 9.94 ppbV
RT: 17.241 min Scan# 5335
Delta R.T. -0.006 min
Lab File: aa8196.D
Acq: 6 Aug 2018 4:50 pm

Tgt Ion	Resp	Lower	Upper
95	1243695		
95	100		
174	55.0	61.5	92.3#



Integrated Analytical Laboratories, LLC

Volatile Organic Compounds by EPA Method TO-15

Laboratory Sample Duplicate Report

SDG Number: E18-06173
 IAL Sample ID: E18-06173-01
 Matrix: Air
 Summa ID: 1070

Date Received: 8/3/18
 Date Analyzed: 8/7/18,8/7/18
 Lab Data File#: AA8228,AA8229
 Dilution Factor: 1
 Injection Volume: 50ml
 GC/MS Column: RTX-1, 0.32 mmID

Compound	CAS #	Sample E18-06173-01 Concentration Reported		Sample Dup E18-06173-21 Concentration Reported		Reporting Limits ppbv	RPD
		ppbv	Q	ppbv	Q		
Acetone	67-64-1	30		27		2.0	10.53%
Allyl Chloride	107-05-1		2.0 U		2.0 U	2.0	0.00%
Benzene	71-43-2		2.0 U		2.0 U	2.0	0.00%
Bromodichloromethane	75-27-4		2.0 U		2.0 U	2.0	0.00%
Bromoform	75-25-2		2.0 U		2.0 U	2.0	0.00%
Bromomethane	74-83-9		2.0 U		2.0 U	2.0	0.00%
1,3-Butadiene	106-99-0		2.0 U		2.0 U	2.0	0.00%
Chlorobenzene	108-90-7		2.0 U		2.0 U	2.0	0.00%
Chloroethane	75-00-3		2.0 U		2.0 U	2.0	0.00%
Chloroform	67-66-3		2.0 U		2.0 U	2.0	0.00%
Chloromethane	74-87-3		2.0 U		2.0 U	2.0	0.00%
Carbon disulfide	75-15-0		2.0 U		2.0 U	2.0	0.00%
Carbon tetrachloride	56-23-5		2.0 U		2.0 U	2.0	0.00%
2-Chlorotoluene	95-49-8		2.0 U		2.0 U	2.0	0.00%
Cyclohexane	110-82-7		2.0 U		2.0 U	2.0	0.00%
Dibromochloromethane	124-48-1		2.0 U		2.0 U	2.0	0.00%
1,2-Dibromoethane	106-93-4		2.0 U		2.0 U	2.0	0.00%
1,2-Dichlorobenzene	95-50-1		2.0 U		2.0 U	2.0	0.00%
1,3-Dichlorobenzene	541-73-1		2.0 U		2.0 U	2.0	0.00%
1,4-Dichlorobenzene	106-46-7		2.0 U		2.0 U	2.0	0.00%
Dichlorodifluoromethane	75-71-8		2.0 U		2.0 U	2.0	0.00%
1,1-Dichloroethane	75-34-3		2.0 U		2.0 U	2.0	0.00%
1,2-Dichloroethane	107-06-2		2.0 U		2.0 U	2.0	0.00%
1,1-Dichloroethene	75-35-4		2.0 U		2.0 U	2.0	0.00%
1,2-Dichloroethene (cis)	156-59-2		2.0 U		2.0 U	2.0	0.00%
1,2-Dichloroethene (trans)	156-60-5		2.0 U		2.0 U	2.0	0.00%
1,2-Dichloropropane	78-87-5		2.0 U		2.0 U	2.0	0.00%
1,3-Dichloropropene (cis)	10061-01-5		2.0 U		2.0 U	2.0	0.00%
1,3-Dichloropropene (trans)	10061-02-6		2.0 U		2.0 U	2.0	0.00%
1,2-Dichlorotetrafluoroethane	76-14-2		2.0 U		2.0 U	2.0	0.00%
1,4-Dioxane	123-91-1		2.0 U		2.0 U	2.0	0.00%
Ethanol	64-17-5	380		370		2.0	2.67%
Ethylbenzene	100-41-4		2.0 U		2.0 U	2.0	0.00%
4-Ethyltoluene	622-96-8		2.0 U		2.0 U	2.0	0.00%
n-Heptane	142-82-5		2.0 U		2.0 U	2.0	0.00%
1,3-Hexachlorobutadiene	87-68-3		2.0 U		2.0 U	2.0	0.00%
n-Hexane	110-54-3		2.0 U		2.0 U	2.0	0.00%
Isopropanol	67-63-0	5.2		5.5		2.0	-5.61%

Qualifir:

E=Concentration exceeds upper level of calibration range for instrument.

D=Extra dilution required for this compound. J=Duplicate samples do not met RPD criteria.

U=Compound ND or under reporting limit.

Integrated Analytical Laboratories, LLC

Volatile Organic Compounds by EPA Method TO-15

Laboratory Sample Duplicate Report

SDG Number: E18-06173
 IAL Sample ID: E18-06173-01
 Matrix: Air
 Summa ID: 1070

Date Received: 8/3/18
 Date Analyzed: 8/7/18,8/7/18
 Lab Data File#: AA8228,AA8229
 Dilution Factor: 1
 Injection Volume: 50ml
 GC/MS Column: RTX-1, 0.32 mmID

Compound	CAS #	Sample E18-06173-01 Concentration Reported		Sample Dup E18-06173-21 Concentration Reported		Reporting Limits ppbv	RPD
		ppbv	Q	ppbv	Q		
Methylene chloride	75-09-2	99		100		2.0	-1.01%
Methyl ethyl ketone	78-93-3	2.8		2.2		2.0	24.00%
Methyl isobutyl ketone	108-10-1		2.0 U		2.0 U	2.0	0.00%
Methyl methacrylate	80-62-6		2.0 U		2.0 U	2.0	0.00%
Methyl tert-butyl ether	1634-04-4		2.0 U		2.0 U	2.0	0.00%
Styrene	100-42-5		2.0 U		2.0 U	2.0	0.00%
Tert-butyl alcohol	75-65-0	3.1		3.0		2.0	3.28%
1,1,2,2-Tetrachloroethane	79-34-5		2.0 U		2.0 U	2.0	0.00%
Tetrachloroethene	127-18-4		2.0 U		2.0 U	2.0	0.00%
Tetrahydrofuran	109-99-9		2.0 U		2.0 U	2.0	0.00%
Toluene	108-88-3	2.7		2.7		2.0	0.00%
1,2,4-Trichlorobenzene	120-82-1		2.0 U		2.0 U	2.0	0.00%
1,1,1-Trichloroethane	71-55-6		2.0 U		2.0 U	2.0	0.00%
1,1,2-Trichloroethane	79-00-5		2.0 U		2.0 U	2.0	0.00%
Trichloroethene	79-01-6		2.0 U		2.0 U	2.0	0.00%
Trichlorofluoromethane	75-69-4		2.0 U		2.0 U	2.0	0.00%
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1		2.0 U		2.0 U	2.0	0.00%
1,2,4-Trimethylbenzene	95-63-6		2.0 U		2.0 U	2.0	0.00%
1,3,5-Trimethylbenzene	108-67-8		2.0 U		2.0 U	2.0	0.00%
2,2,4-Trimethylpentane	540-84-1		2.0 U		2.0 U	2.0	0.00%
Vinyl bromide	593-60-2		2.0 U		2.0 U	2.0	0.00%
Vinyl chloride	75-01-4		2.0 U		2.0 U	2.0	0.00%
Xylenes (m&p)	179601-23-1		4.0 U		4.0 U	4.0	0.00%
Xylenes (o)	95-47-6		2.0 U		2.0 U	2.0	0.00%

RPD must be <25% for all laboratory duplicate samples. Laboratory duplicate samples are run once daily.

NC = The RPD could not be calculated since the compound was only detected in either the parent or duplicate sample.

Qualifir:

E=Concentration exceeds upper level of calibration range for instrument.
 D=Extra dilution required for this compound. J=Duplicate samples do not met RPD criteria.
 U=Compound ND or under reporting limit.

Data Path : C:\DATA\08-07-18\
 Data File : aa8228.D
 Acq On : 7 Aug 2018 2:44 pm
 Operator : jjw
 Sample : E18-06173-01 x 10 dil
 Misc : 1070, 50cc
 ALS Vial : 8 Sample Multiplier: 1

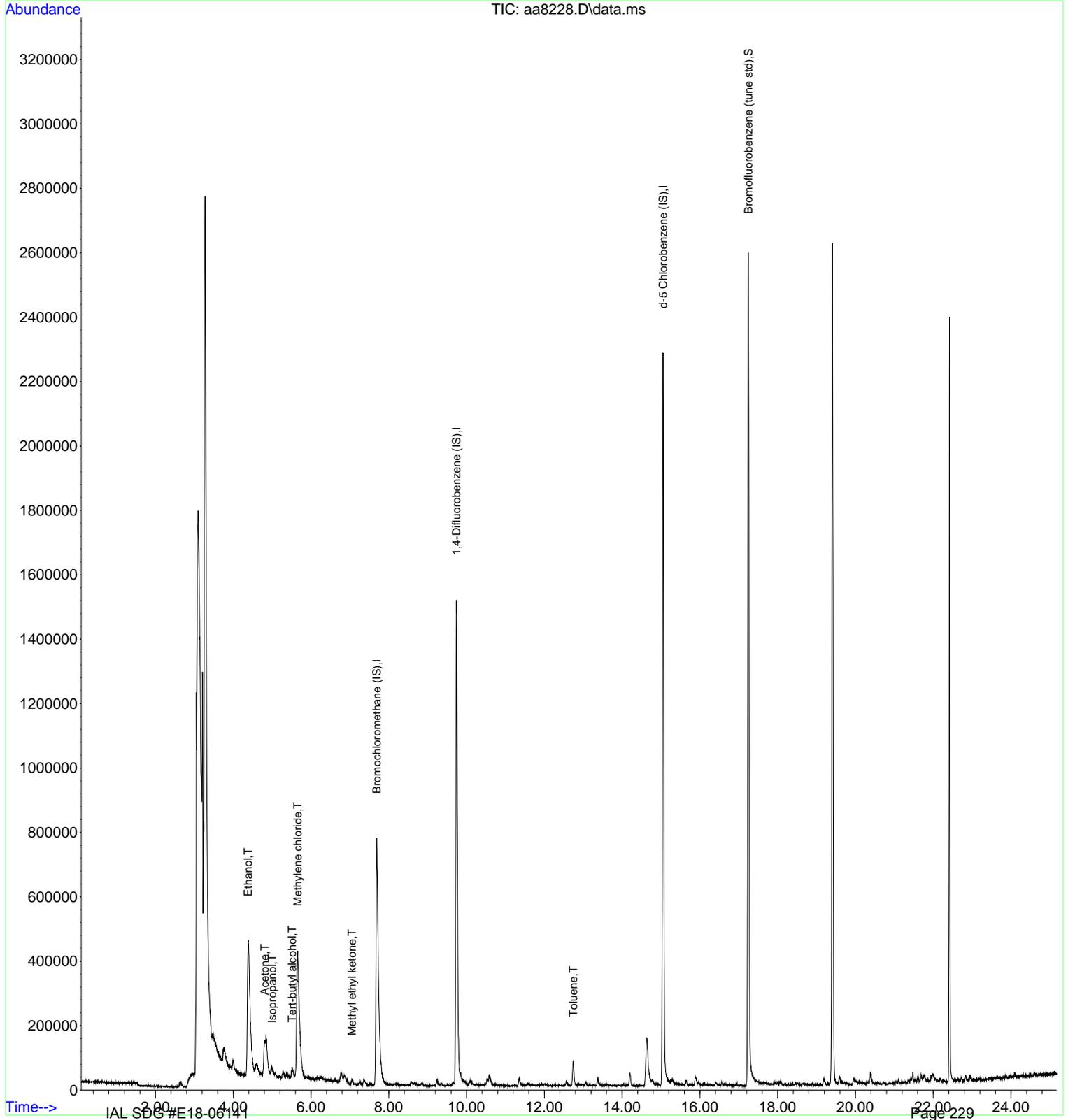
Quant Time: Aug 08 13:07:49 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

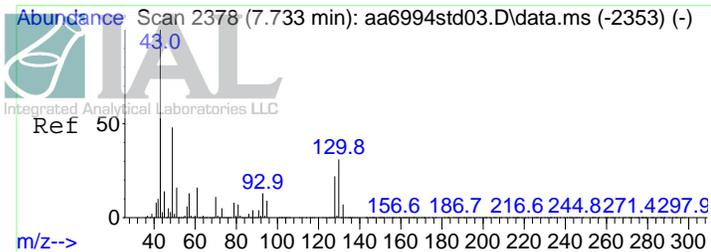
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Bromochloromethane (IS)	7.692	130	384714	10.00	ppbV	0.00
38) 1,4-Difluorobenzene (IS)	9.740	114	1413711	10.00	ppbV	0.00
55) d-5 Chlorobenzene (IS)	15.055	117	1261827	10.00	ppbV	0.00
System Monitoring Compounds						
64) Bromofluorobenzene (tu...	17.248	95	1085952	9.44	ppbV	0.00
Target Compounds						
11) Ethanol	4.383	45	1144225	37.77	ppbV	97
14) Acetone	4.808	58	88386	3.02	ppbV #	49
16) Isopropanol	5.001	45	49479	0.52	ppbV #	53
19) Methylene chloride	5.657	49	729142	9.89	ppbV	96
20) Tert-butyl alcohol	5.512	59	37800	0.31	ppbV	100
27) Methyl ethyl ketone	7.052	43	30133	0.28	ppbV #	74
50) Toluene	12.743	91	48498	0.27	ppbV	94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\08-07-18\
 Data File : aa8228.D
 Acq On : 7 Aug 2018 2:44 pm
 Operator : jjw
 Sample : E18-06173-01 x 10 dil
 Misc : 1070, 50cc
 ALS Vial : 8 Sample Multiplier: 1

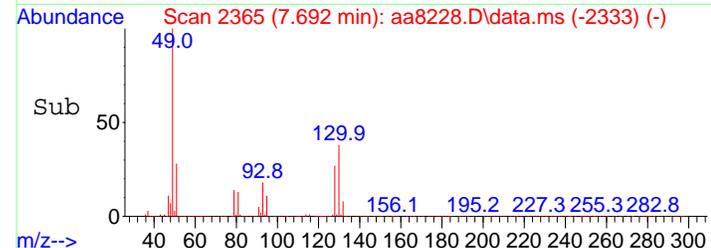
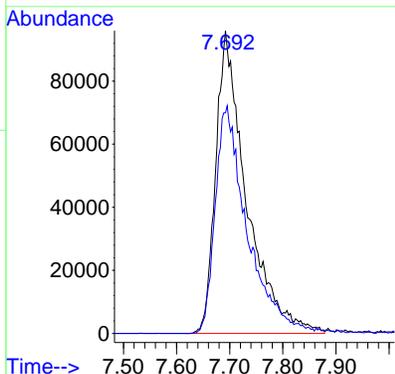
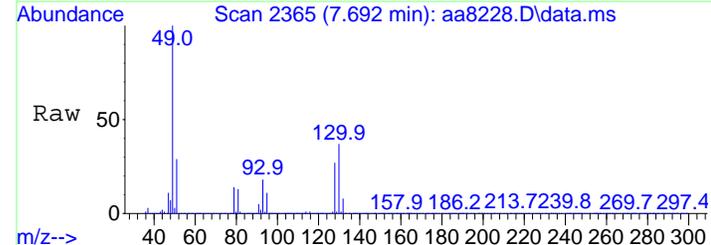
Quant Time: Aug 08 13:07:49 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration





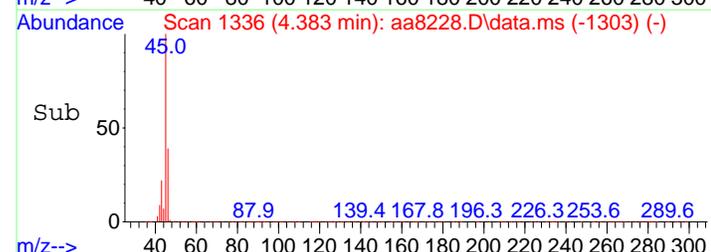
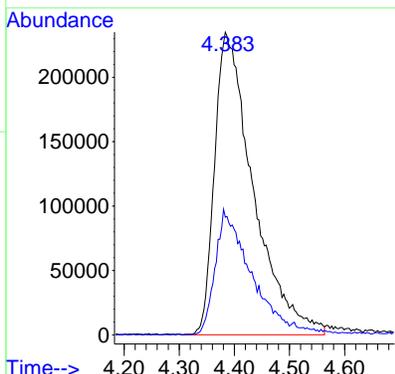
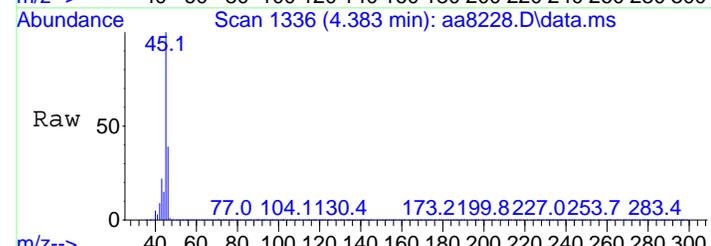
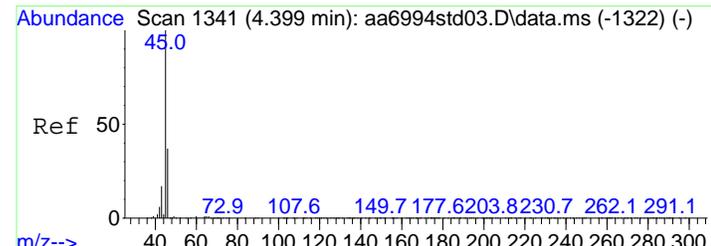
#1
 Bromochloromethane (IS)
 Concen: 10.00 ppbV
 RT: 7.692 min Scan# 2365
 Delta R.T. 0.003 min
 Lab File: aa8228.D
 Acq: 7 Aug 2018 2:44 pm

Tgt Ion:130 Resp: 384714
 Ion Ratio Lower Upper
 130 100
 128 77.9 62.6 94.0

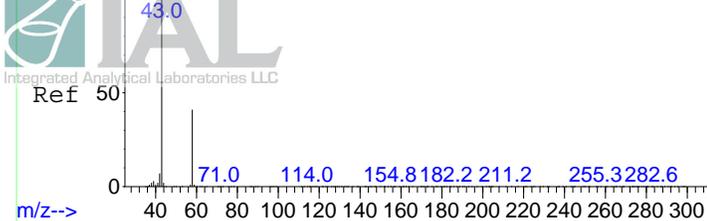


#11
 Ethanol
 Concen: 37.77 ppbV
 RT: 4.383 min Scan# 1336
 Delta R.T. 0.006 min
 Lab File: aa8228.D
 Acq: 7 Aug 2018 2:44 pm

Tgt Ion: 45 Resp: 1144225
 Ion Ratio Lower Upper
 45 100
 46 38.7 32.6 48.8

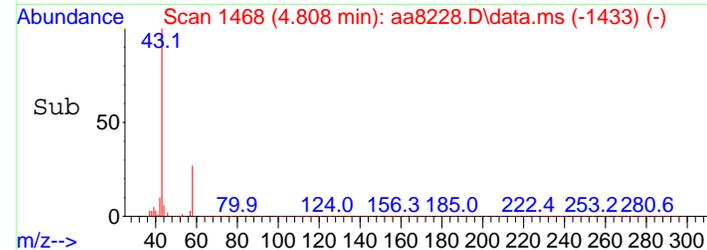
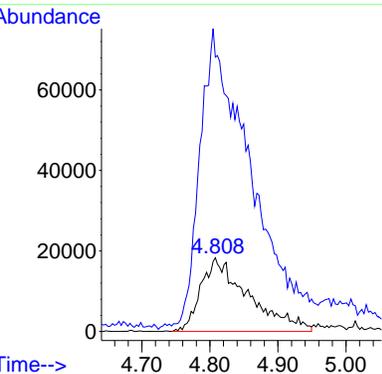
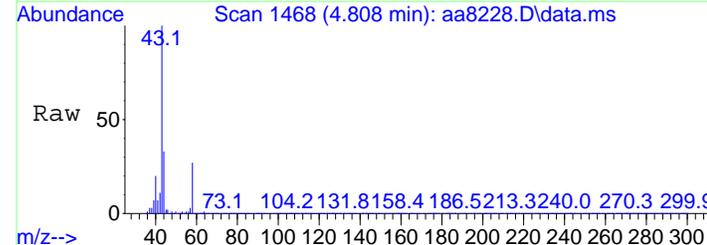


Abundance Scan 1470 (4.814 min): aa6994std03.D\data.ms (-1454) (-)



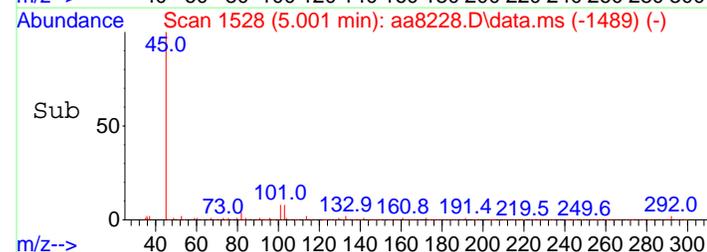
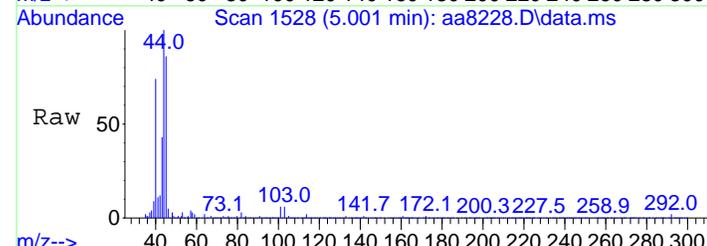
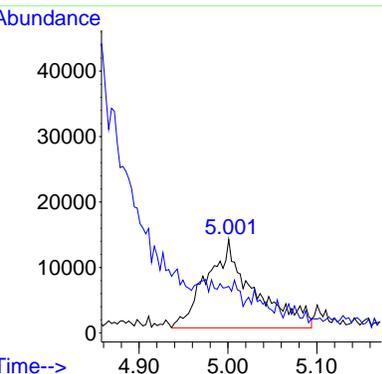
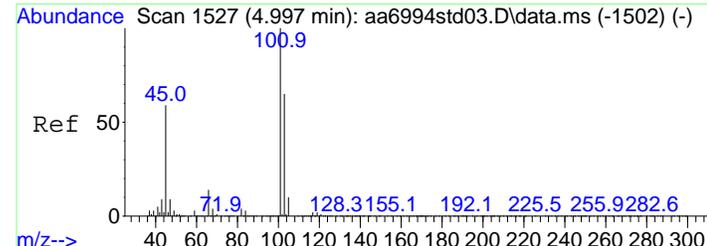
#14
Acetone
Concen: 3.02 ppbV
RT: 4.808 min Scan# 1468
Delta R.T. 0.013 min
Lab File: aa8228.D
Acq: 7 Aug 2018 2:44 pm

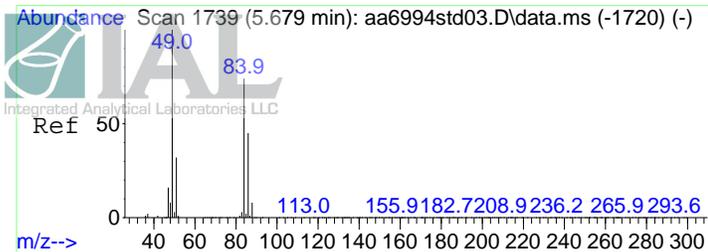
Tgt Ion: 58 Resp: 88386
Ion Ratio Lower Upper
58 100
43 436.1 263.2 394.8#



#16
Isopropanol
Concen: 0.52 ppbV
RT: 5.001 min Scan# 1528
Delta R.T. 0.026 min
Lab File: aa8228.D
Acq: 7 Aug 2018 2:44 pm

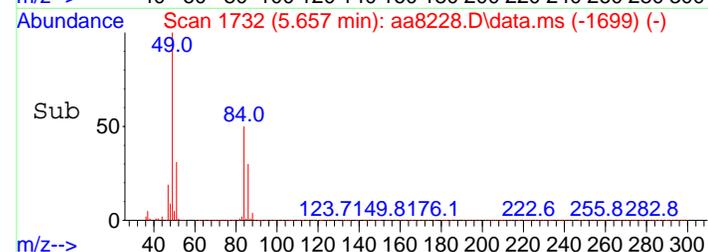
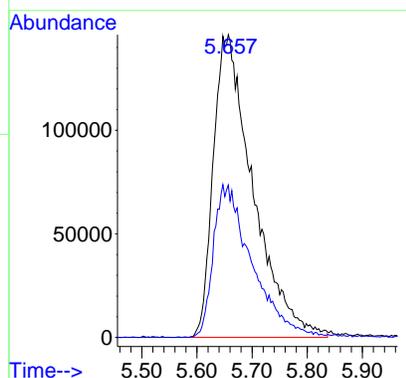
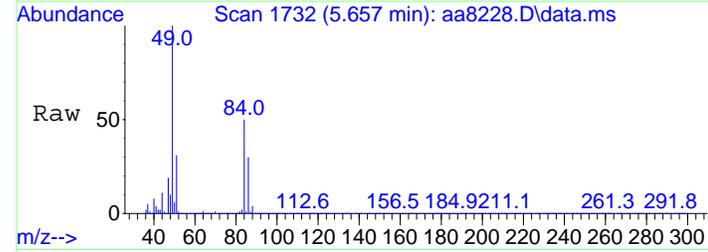
Tgt Ion: 45 Resp: 49479
Ion Ratio Lower Upper
45 100
43 0.0 18.4 27.6#





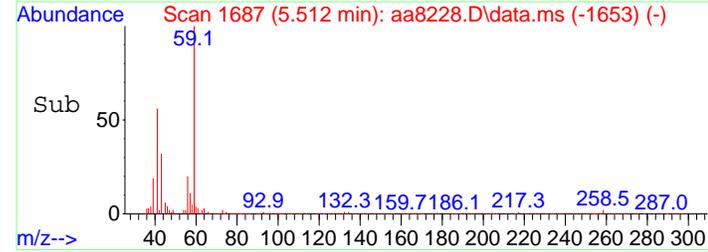
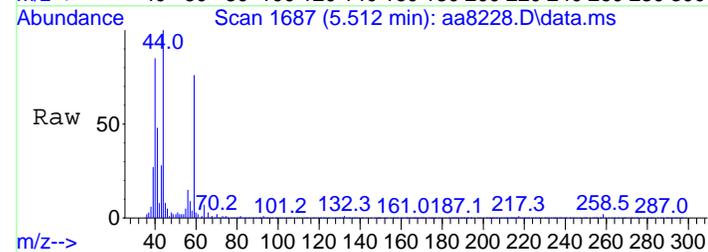
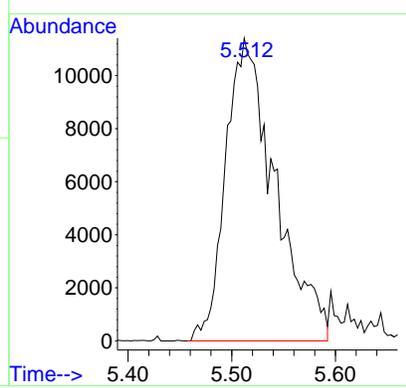
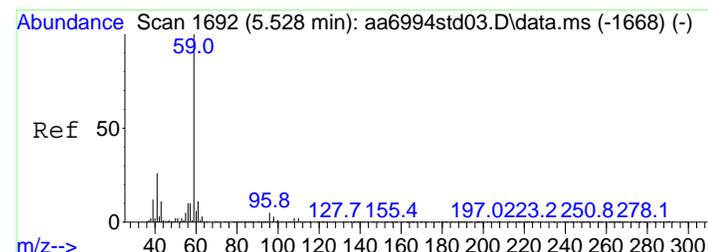
#19
 Methylene chloride
 Concen: 9.89 ppbV
 RT: 5.657 min Scan# 1732
 Delta R.T. 0.006 min
 Lab File: aa8228.D
 Acq: 7 Aug 2018 2:44 pm

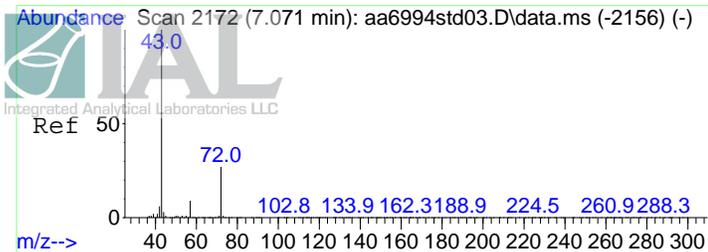
Tgt Ion: 49 Resp: 729142
 Ion Ratio Lower Upper
 49 100
 84 49.1 41.6 62.4



#20
 Tert-butyl alcohol
 Concen: 0.31 ppbV
 RT: 5.512 min Scan# 1687
 Delta R.T. 0.010 min
 Lab File: aa8228.D
 Acq: 7 Aug 2018 2:44 pm

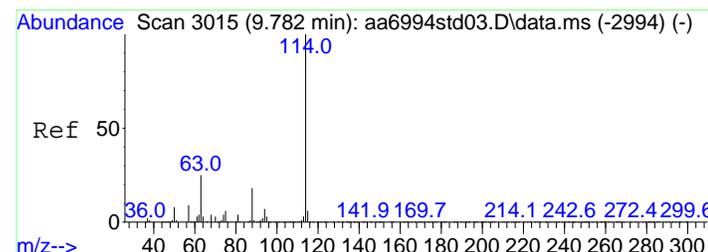
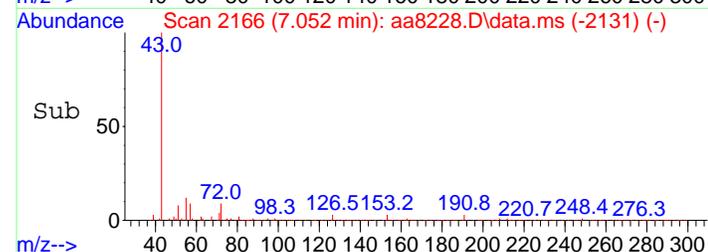
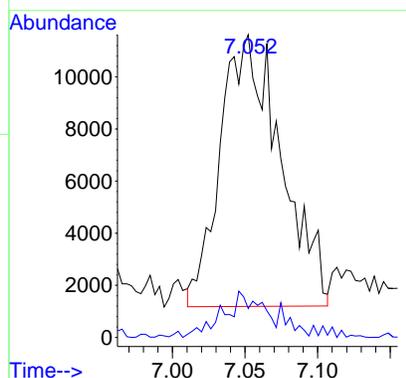
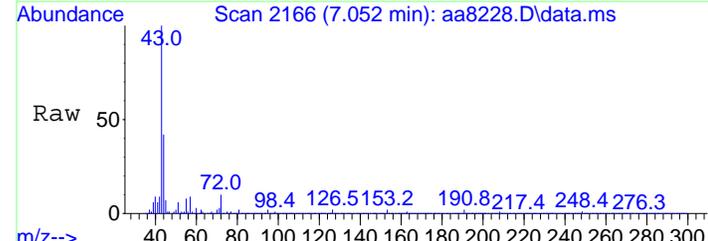
Tgt Ion: 59 Resp: 37800





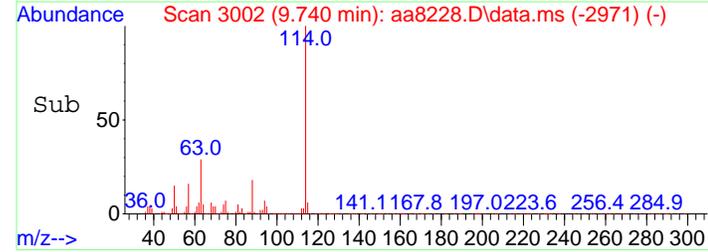
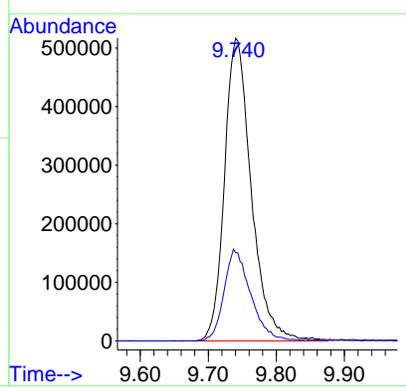
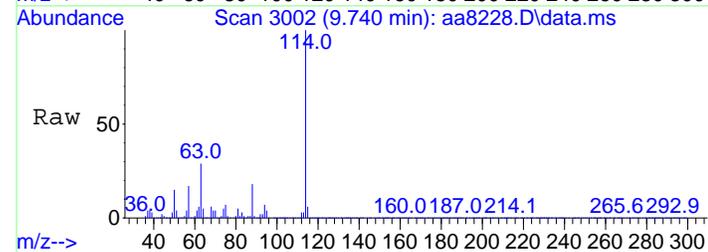
#27
 Methyl ethyl ketone
 Concen: 0.28 ppbV
 RT: 7.052 min Scan# 2166
 Delta R.T. 0.013 min
 Lab File: aa8228.D
 Acq: 7 Aug 2018 2:44 pm

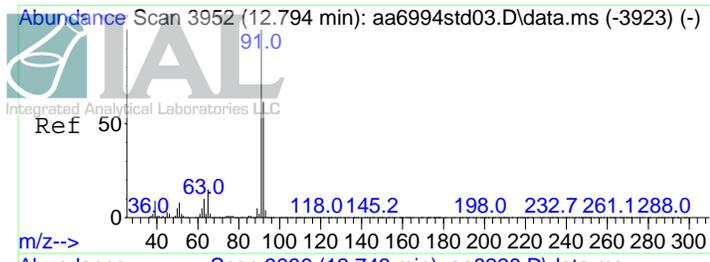
Tgt Ion	Resp	Lower	Upper
43	100		
72	6.5	14.4	21.6#



#38
 1,4-Difluorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 9.740 min Scan# 3002
 Delta R.T. 0.000 min
 Lab File: aa8228.D
 Acq: 7 Aug 2018 2:44 pm

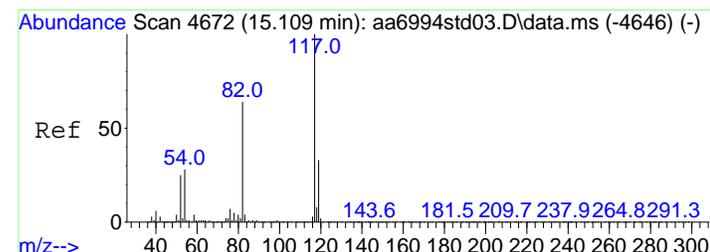
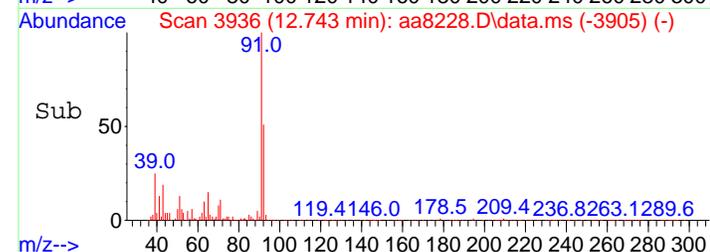
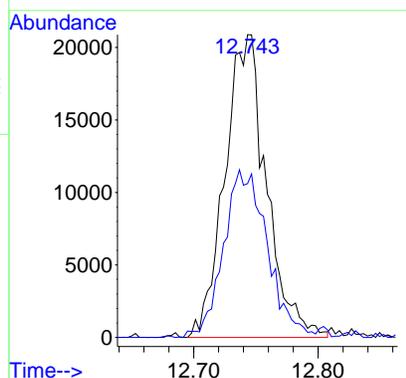
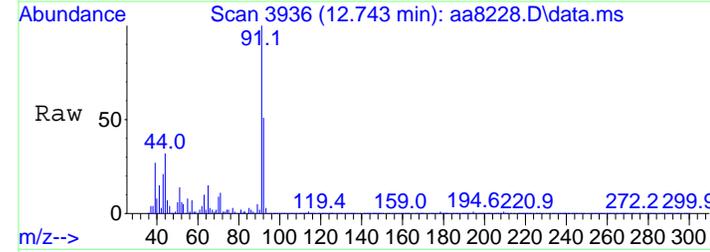
Tgt Ion	Resp	Lower	Upper
114	100		
63	29.9	20.0	30.0





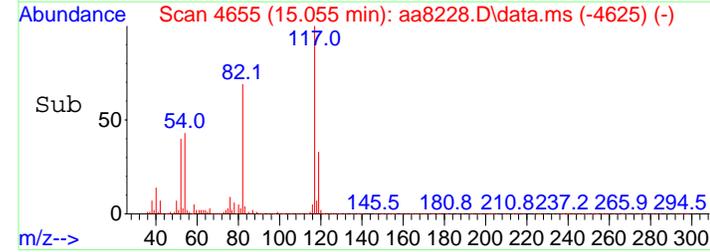
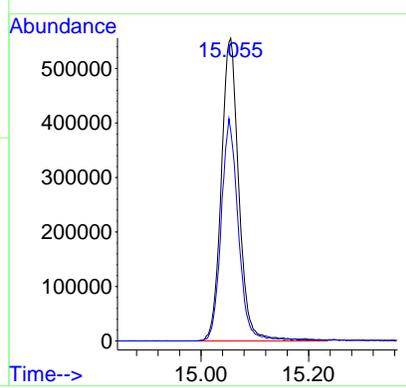
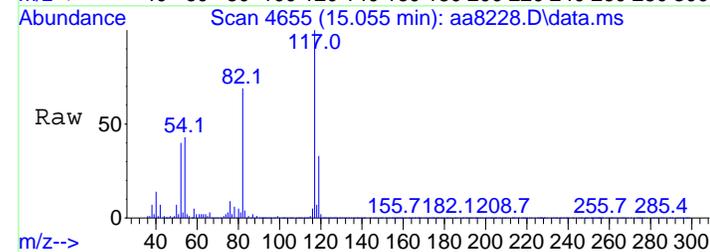
#50
 Toluene
 Concen: 0.27 ppbV
 RT: 12.743 min Scan# 3936
 Delta R.T. 0.000 min
 Lab File: aa8228.D
 Acq: 7 Aug 2018 2:44 pm

Tgt Ion	Resp	Lower	Upper
91	100		
92	57.9	50.0	75.0

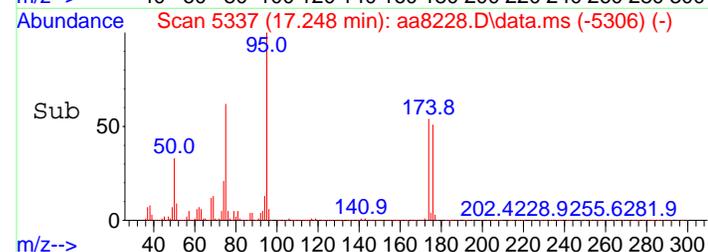
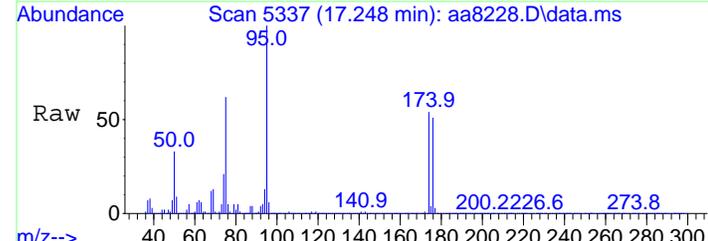
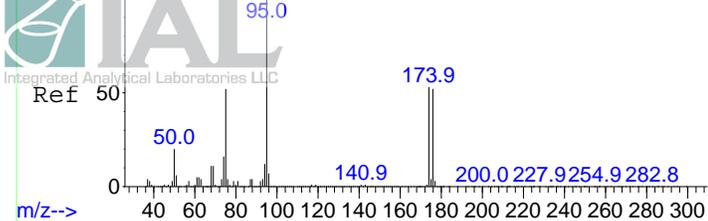


#55
 d-5 Chlorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 15.055 min Scan# 4655
 Delta R.T. -0.003 min
 Lab File: aa8228.D
 Acq: 7 Aug 2018 2:44 pm

Tgt Ion	Resp	Lower	Upper
117	100		
82	71.2	56.0	84.0

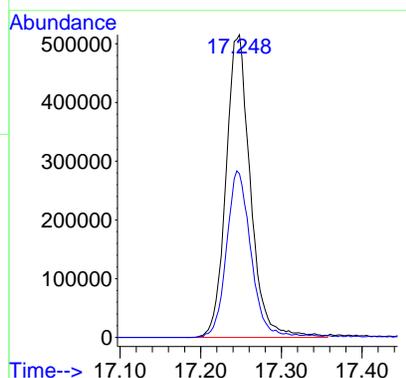


Abundance Scan 5355 (17.305 min): aa6994std03.D\data.ms (-5334) (-)



#64
Bromofluorobenzene (tune std)
Concen: 9.44 ppbV
RT: 17.248 min Scan# 5337
Delta R.T. 0.000 min
Lab File: aa8228.D
Acq: 7 Aug 2018 2:44 pm

Tgt Ion	Resp	Lower	Upper
95	1085952		
95	100		
174	55.3	61.5	92.3#



Data Path : C:\DATA\08-07-18\
 Data File : aa8229.D
 Acq On : 7 Aug 2018 3:17 pm
 Operator : jjw
 Sample : E18-06173-21 x 10 dil
 Misc : dup of E18-06173-01 x 10 dil, can 1070
 ALS Vial : 9 Sample Multiplier: 1

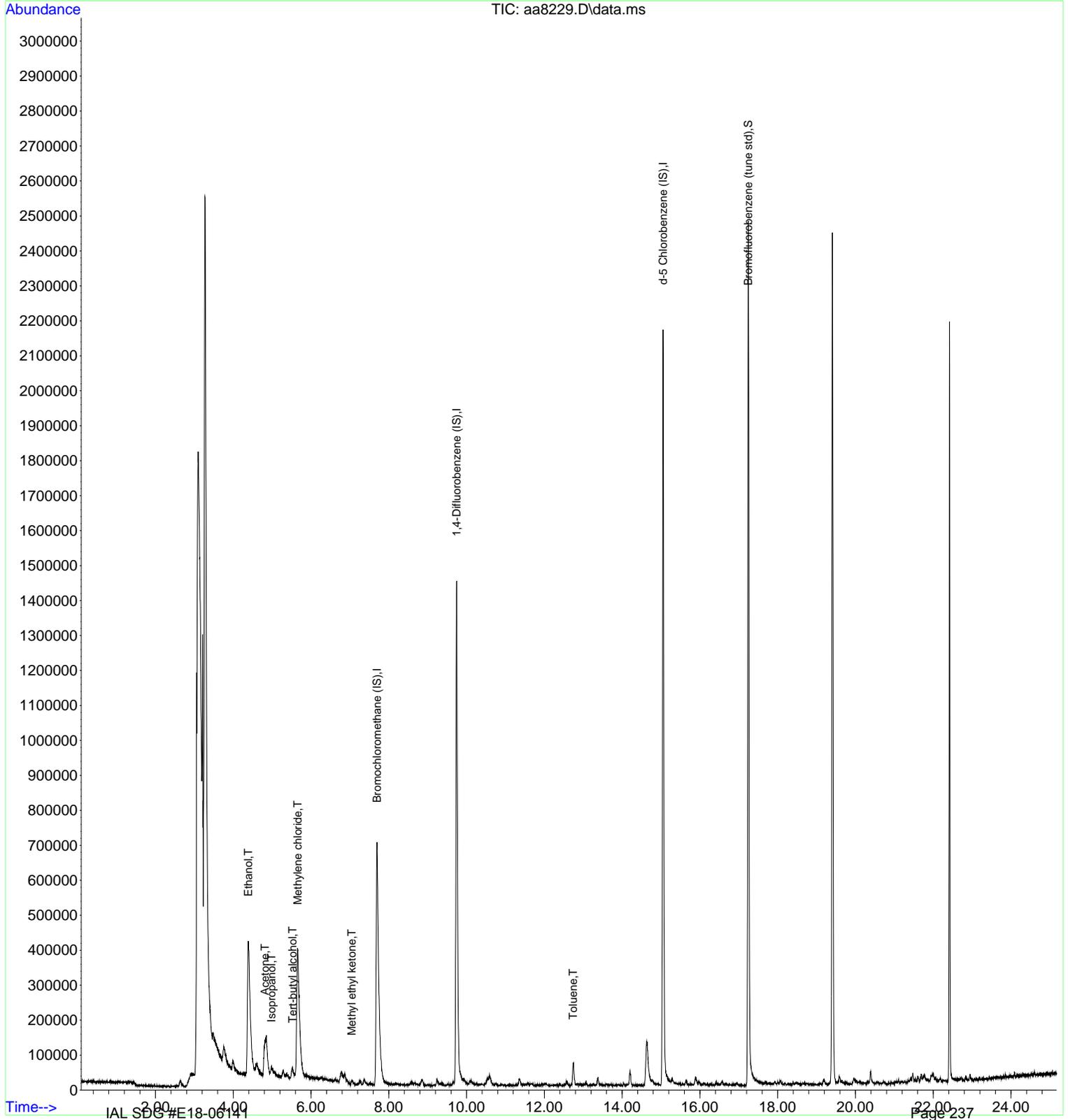
Quant Time: Aug 08 13:09:58 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration

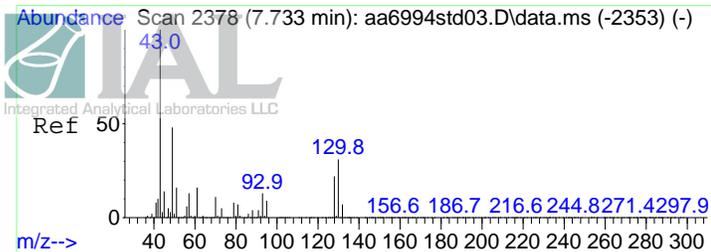
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Bromochloromethane (IS)	7.695	130	353353	10.00	ppbV	0.00
38) 1,4-Difluorobenzene (IS)	9.743	114	1303289	10.00	ppbV #	0.00
55) d-5 Chlorobenzene (IS)	15.055	117	1182462	10.00	ppbV	0.00
System Monitoring Compounds						
64) Bromofluorobenzene (tu...	17.245	95	1022916	9.49	ppbV	0.00
Target Compounds						
11) Ethanol	4.387	45	1026205	36.88	ppbV	97
14) Acetone	4.814	58	73001	2.71	ppbV #	40
16) Isopropanol	4.985	45	48384	0.55	ppbV #	71
19) Methylene chloride	5.660	49	680511	10.05	ppbV	94
20) Tert-butyl alcohol	5.525	59	33629	0.30	ppbV	100
27) Methyl ethyl ketone	7.046	43	21982	0.22	ppbV	98
50) Toluene	12.737	91	43258	0.26	ppbV	89

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\08-07-18\
 Data File : aa8229.D
 Acq On : 7 Aug 2018 3:17 pm
 Operator : jjw
 Sample : E18-06173-21 x 10 dil
 Misc : dup of E18-06173-01 x 10 dil, can 1070
 ALS Vial : 9 Sample Multiplier: 1

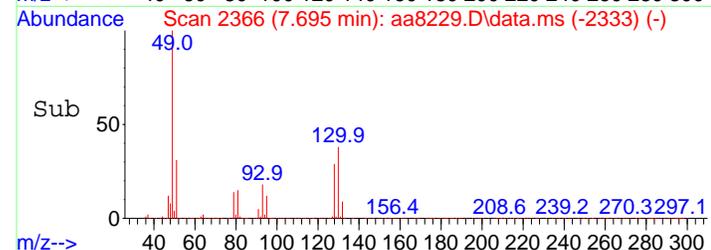
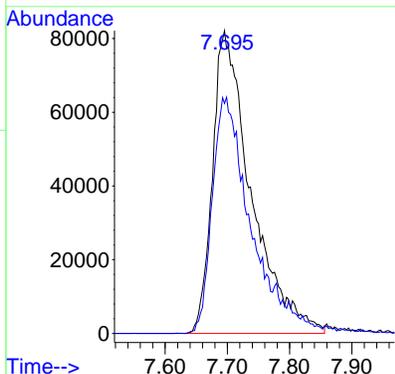
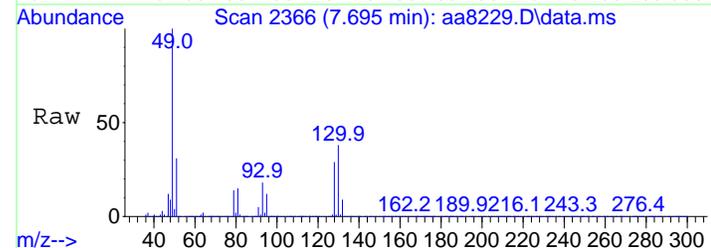
Quant Time: Aug 08 13:09:58 2018
 Quant Method : C:\msdchem\1\METHODS\0725.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Wed Jul 25 14:15:57 2018
 Response via : Initial Calibration





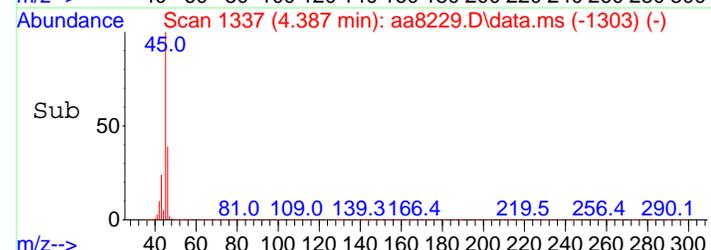
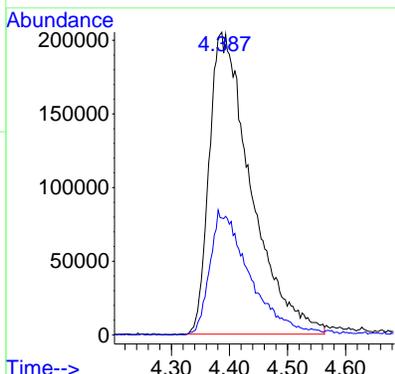
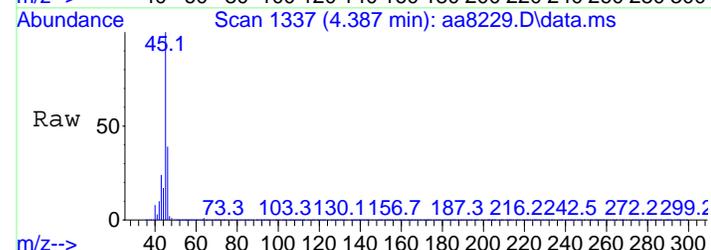
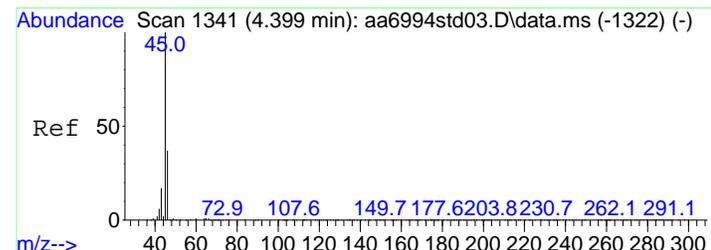
#1
 Bromochloromethane (IS)
 Concen: 10.00 ppbV
 RT: 7.695 min Scan# 2366
 Delta R.T. 0.006 min
 Lab File: aa8229.D
 Acq: 7 Aug 2018 3:17 pm

Tgt Ion:130 Resp: 353353
 Ion Ratio Lower Upper
 130 100
 128 77.9 62.6 94.0

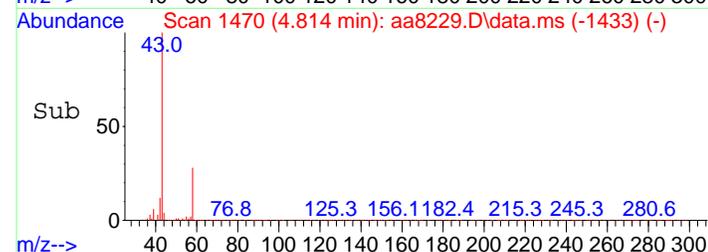
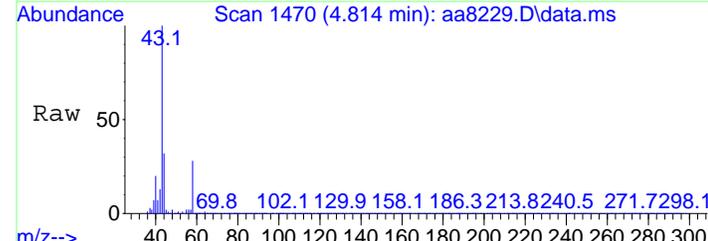
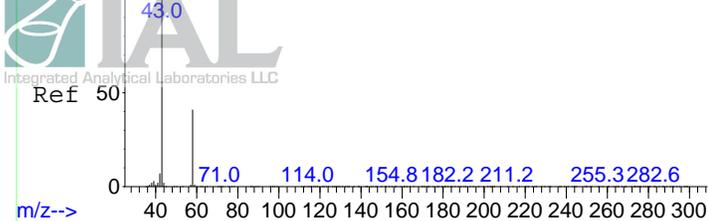


#11
 Ethanol
 Concen: 36.88 ppbV
 RT: 4.387 min Scan# 1337
 Delta R.T. 0.010 min
 Lab File: aa8229.D
 Acq: 7 Aug 2018 3:17 pm

Tgt Ion: 45 Resp: 1026205
 Ion Ratio Lower Upper
 45 100
 46 39.0 32.6 48.8

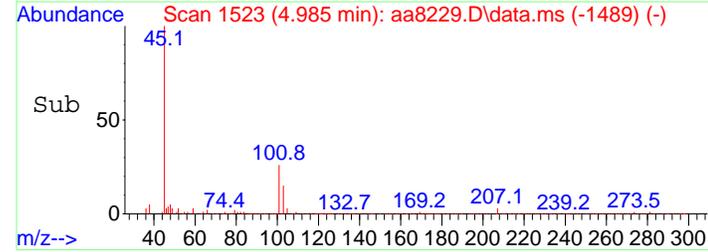
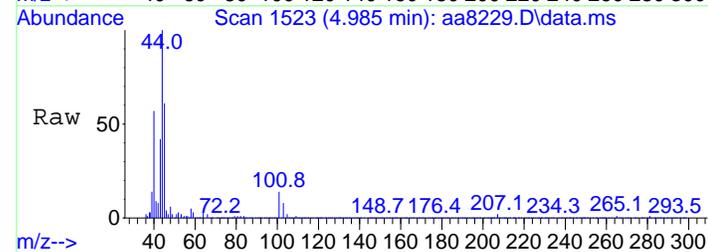
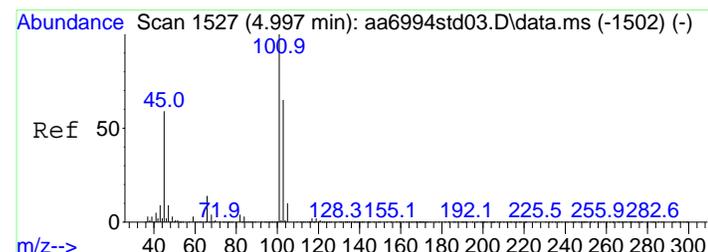
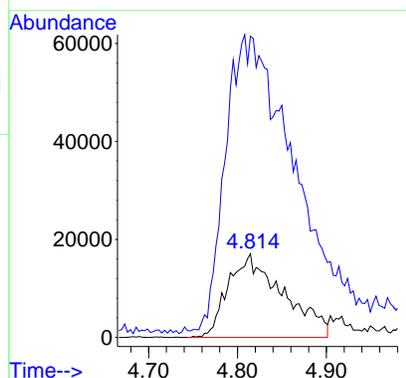


Abundance Scan 1470 (4.814 min): aa6994std03.D\data.ms (-1454) (-)



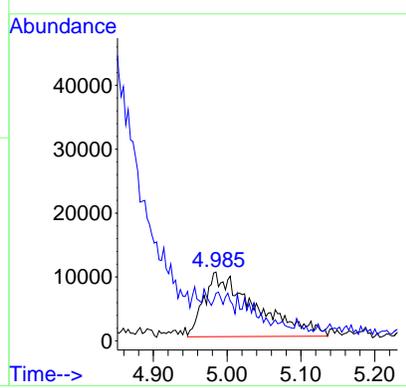
#14
Acetone
Concen: 2.71 ppbV
RT: 4.814 min Scan# 1470
Delta R.T. 0.019 min
Lab File: aa8229.D
Acq: 7 Aug 2018 3:17 pm

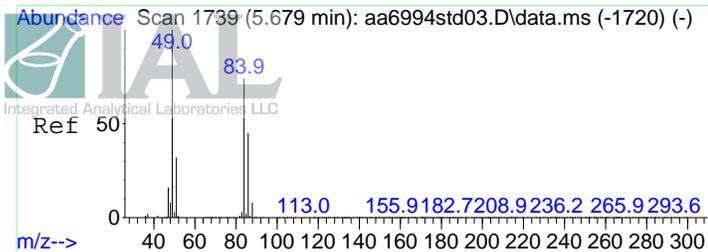
Tgt Ion	Resp	Lower	Upper
58	73001		
58	100		
43	454.8	263.2	394.8#



#16
Isopropanol
Concen: 0.55 ppbV
RT: 4.985 min Scan# 1523
Delta R.T. 0.010 min
Lab File: aa8229.D
Acq: 7 Aug 2018 3:17 pm

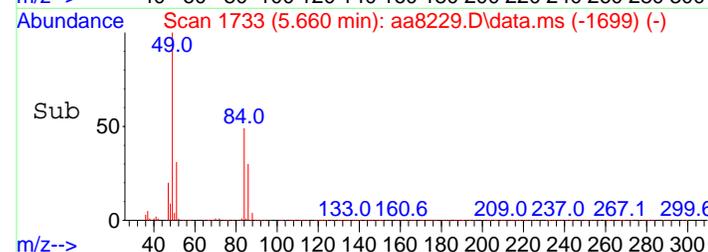
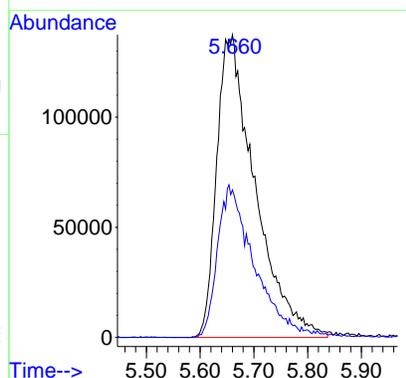
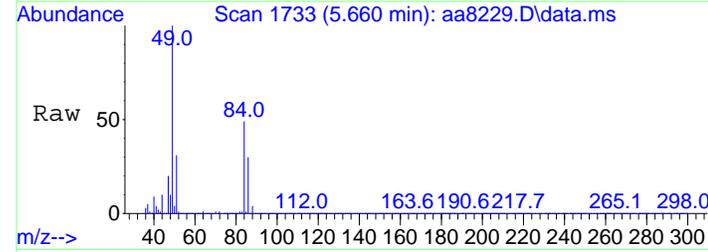
Tgt Ion	Resp	Lower	Upper
45	48384		
45	100		
43	8.8	18.4	27.6#





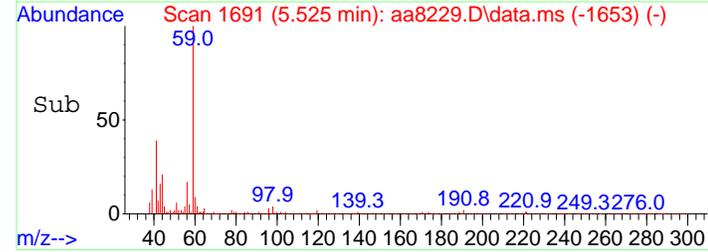
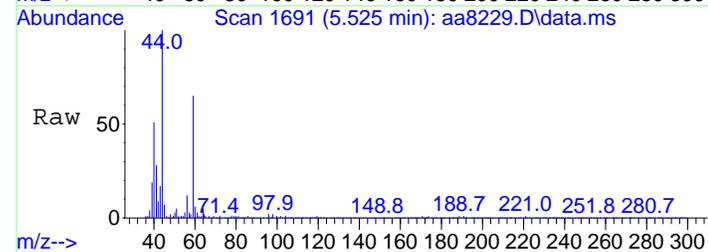
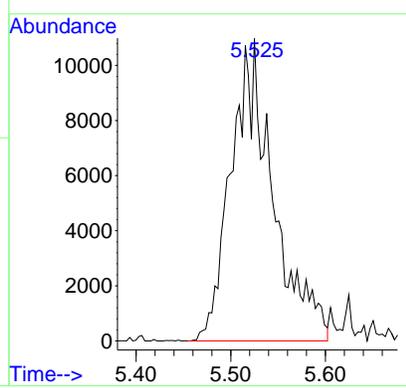
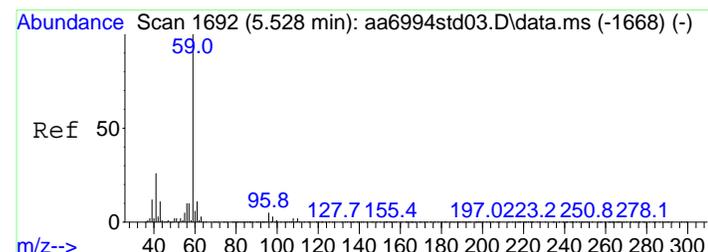
#19
 Methylene chloride
 Concen: 10.05 ppbV
 RT: 5.660 min Scan# 1733
 Delta R.T. 0.010 min
 Lab File: aa8229.D
 Acq: 7 Aug 2018 3:17 pm

Tgt Ion: 49 Resp: 680511
 Ion Ratio Lower Upper
 49 100
 84 47.9 41.6 62.4



#20
 Tert-butyl alcohol
 Concen: 0.30 ppbV
 RT: 5.525 min Scan# 1691
 Delta R.T. 0.023 min
 Lab File: aa8229.D
 Acq: 7 Aug 2018 3:17 pm

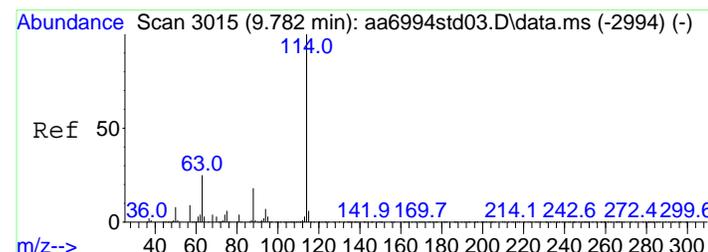
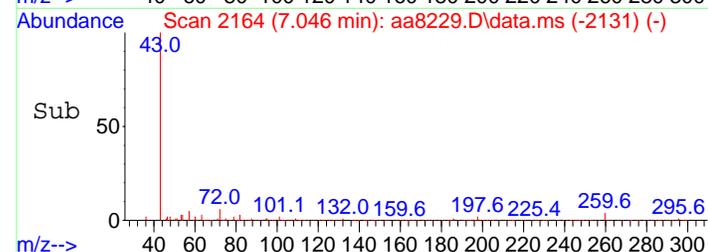
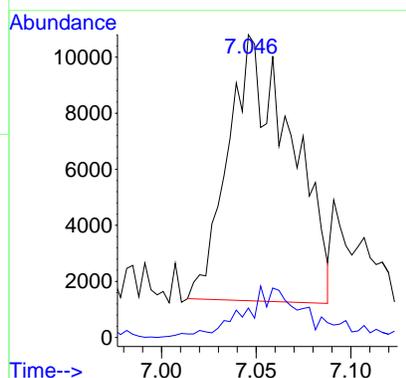
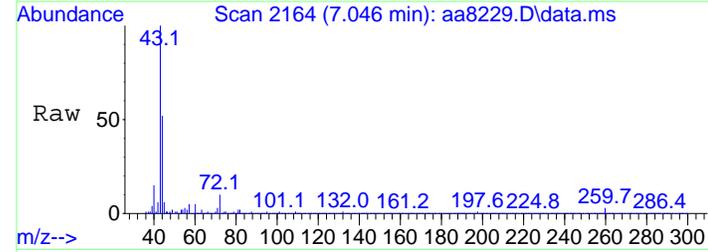
Tgt Ion: 59 Resp: 33629





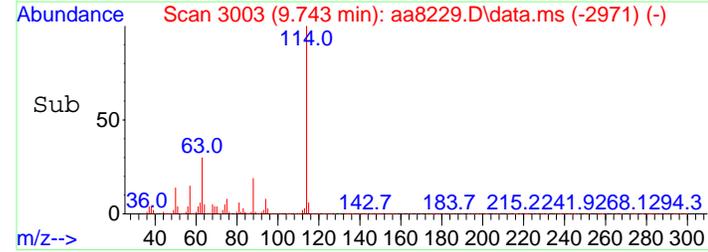
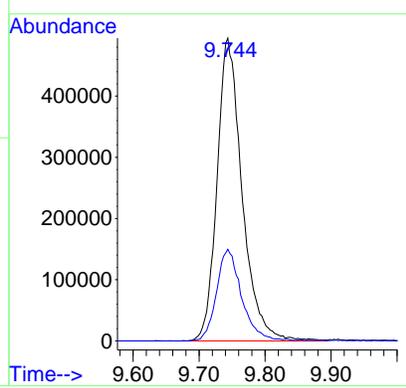
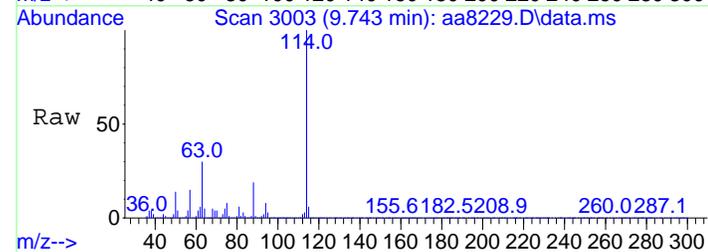
#27
 Methyl ethyl ketone
 Concen: 0.22 ppbV
 RT: 7.046 min Scan# 2164
 Delta R.T. 0.007 min
 Lab File: aa8229.D
 Acq: 7 Aug 2018 3:17 pm

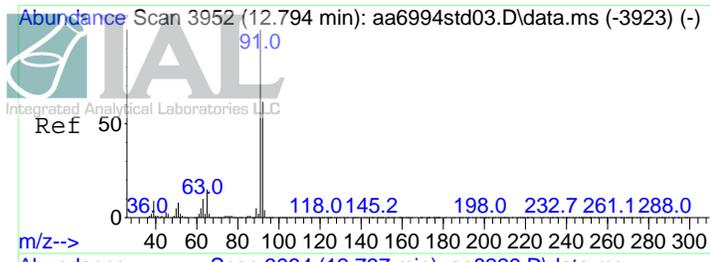
Tgt Ion	Resp	Lower	Upper
43	100		
72	18.8	14.4	21.6



#38
 1,4-Difluorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 9.743 min Scan# 3003
 Delta R.T. 0.003 min
 Lab File: aa8229.D
 Acq: 7 Aug 2018 3:17 pm

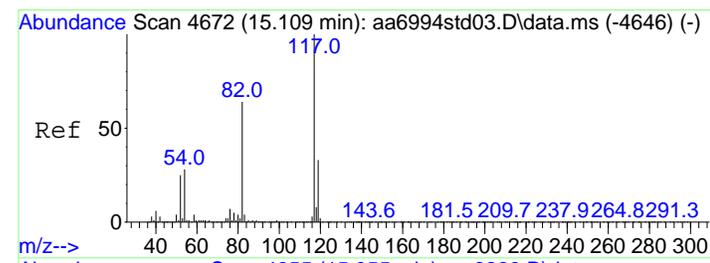
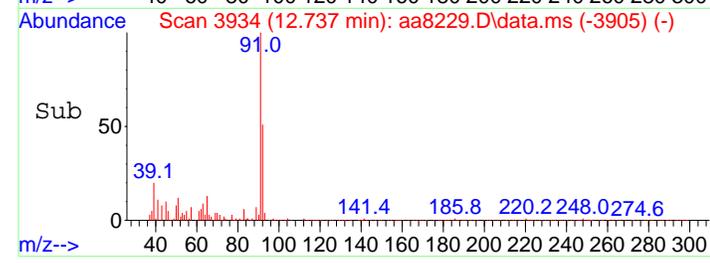
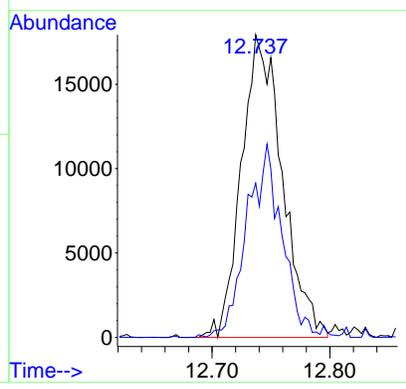
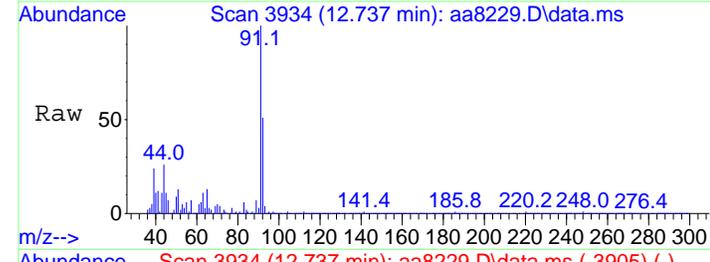
Tgt Ion	Resp	Lower	Upper
114	100		
63	30.3	20.0	30.0#





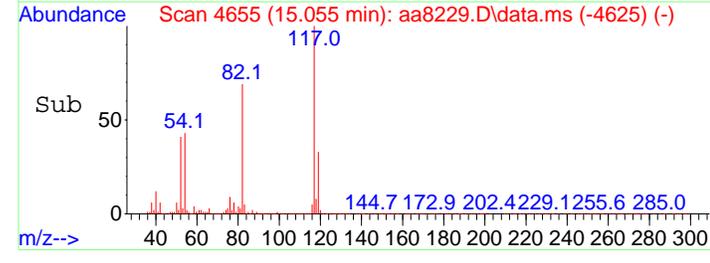
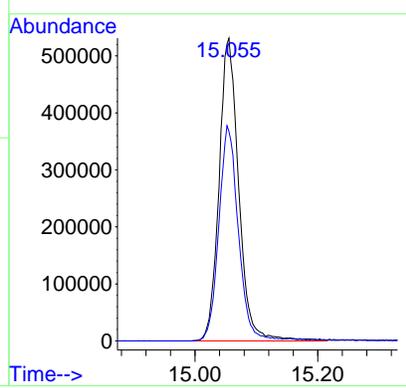
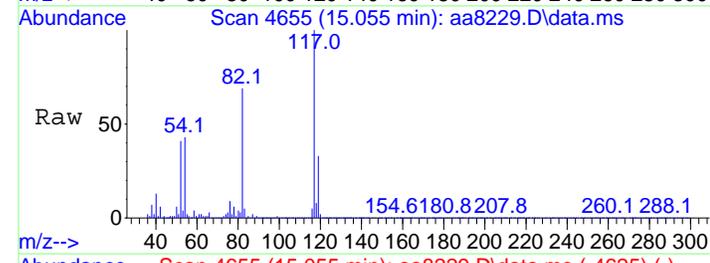
#50
 Toluene
 Concen: 0.26 ppbV
 RT: 12.737 min Scan# 3934
 Delta R.T. -0.006 min
 Lab File: aa8229.D
 Acq: 7 Aug 2018 3:17 pm

Tgt Ion	Resp	Lower	Upper
91	100		
92	54.2	50.0	75.0

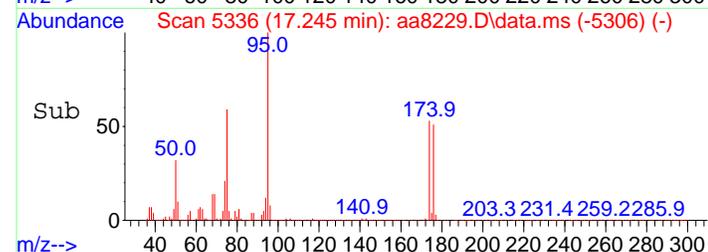
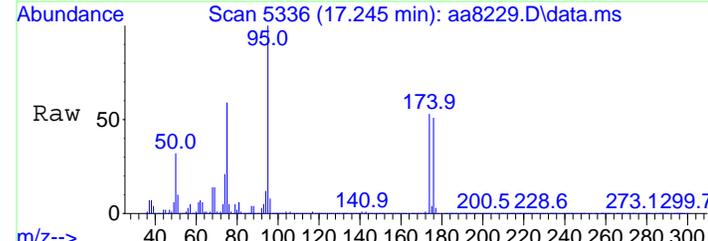
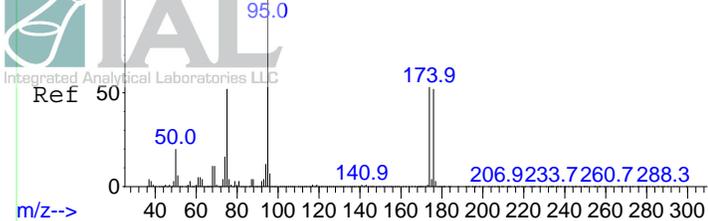


#55
 d-5 Chlorobenzene (IS)
 Concen: 10.00 ppbV
 RT: 15.055 min Scan# 4655
 Delta R.T. -0.003 min
 Lab File: aa8229.D
 Acq: 7 Aug 2018 3:17 pm

Tgt Ion	Resp	Lower	Upper
117	100		
82	70.7	56.0	84.0

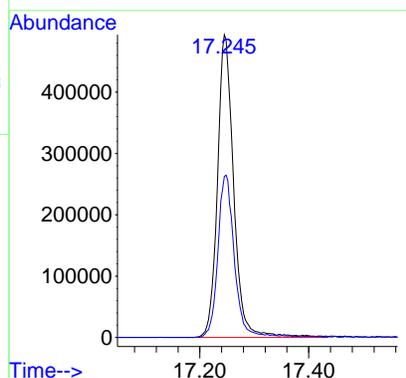


Abundance Scan 5355 (17.305 min): aa6994std03.D\data.ms (-5334) (-)



#64
Bromofluorobenzene (tune std)
Concen: 9.49 ppbV
RT: 17.245 min Scan# 5336
Delta R.T. -0.003 min
Lab File: aa8229.D
Acq: 7 Aug 2018 3:17 pm

Tgt Ion: 95 Resp: 1022916
Ion Ratio Lower Upper
95 100
174 54.5 61.5 92.3#



Integrated Analytical Laboratories
GC/MS Run Log

Instrument ID: Agilent 7890A / 5975C
Column: Restek RTX-1 SN 992567

Target Directory: D:\Agilent GCMS\

Date of Analysis: 8/6-7/2018
Date of Initial Calibration: 5/18/2018, 7/25/2018
SDG #: E18-06141

File #	Laboratory Sample ID	QC Check	Dilution Factor	Can #	Analyst	Injection Volume (cc)	Comments	Make-up Air		Acquisition		Room		TO-15 Standard	
								Added to canister (cc)	Added for dilution (cc)	Date	Time	Temp	BP "Hg	Working ID	Vendor ID (Lot#)
aa7071bfb	BFB	✓		ALM029426	JLS	50				5/18/2018	8:17	74	30.06		403-402623
aa7072std01	40 ppbv Std	✓		CC483586	JLS	200								5/18/2018	160-401140945-1
aa7073std02	20 ppbv Std	✓		CC483586	JLS	100								5/18/2018	160-401140945-1
aa7074std03	10 ppbv Std	✓		CC483586	JLS	50								5/18/2018	160-401140945-1
aa7075std04	2 ppbv Std	✓		CC483586	JLS	10								5/18/2018	160-401140945-1
aa7076std05	0.2 ppbv Std	✓		CC483586	JLS	1								5/18/2018	160-401140945-1
aa7077icvss	10 ppbv ICVSS	✓		CC483422	JLS	50									160-401186845-1

Analyst Name: Jeff Schmitt

Signature: 

IAL SDG #E18-06141

Supervisor Name: Lauren Jenkins

Signature: 

Page 244

**Integrated Analytical Laboratories
GC/MS Run Log**

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Column: Restek RTX-1 SN 992567

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Date of Initial Calibration: 5/18/2018, 7/25/2018
SDG #: E18-06141

File #	Laboratory Sample ID	QC Check	Dilution Factor	Can #	Analyst	Injection Volume (cc)	Comments	Make-up Air		Acquisition		Room		TO-15 Standard	
								Added to canister (cc)	Added for dilution (cc)	Date	Time	Temp	BP "Hg	Working ID	Vendor ID (Lot#)
aa7471bfb	BFB	✓		ALM029426	JLS	50				6/13/2018	8:30	72	30.07		403-402623
aa7472dcvs	10 ppbv DCVS	✓		AAL071156	JLS	50								5/18/2018	160-401140945-1
aa7473rlcs	0.2 ppbv RLLCS	✓		AAL071156	JLS	1								5/18/2018	160-401140945-1
aa7474blk	Method Blank	✓		1127	JLS	500									
aa7475	3059	✓			JLS	500									
aa7476	E18-04190-11 x 5	✓	5	3280	JLS	100									
aa7477	blk	✓			JLS	500									
aa7478	E18-04189-01	✓		5096	JLS	500									
aa7479	E18-04189-21	✓	dup of E18-04189-01, c		JLS	500									
aa7480	E18-04189-02	✓		5101	JLS	500									
aa7481	E18-04189-03	✓		3039A	JLS	500									
aa7482	E18-04189-04 x 100	✓	100	1771	JLS	5	not used								
aa7483	E18-04189-05 x 100	✓	100	1541	JLS	5	not used								
aa7484	blk	✓			JLS	500									
aa7485cccvs	10 ppbv CCCVS	✓		CC483586	JLS	50				6/13/2018	20:47			5/18/2018	160-401140945-1

Analyst Name: Jeff Schmitt

Signature: *Jeff Schmitt*

IAL SDG #E18-06141

Supervisor Name: Lauren Jenkins

Signature: *Lauren Jenkins*

Page 245

**Integrated Analytical Laboratories
GC/MS Run Log**

Instrument ID: Agilent 7890A / 5975C
Column: Restek RTX-1 SN 992567

Target Directory: D:\Agilent GCMS\

Date of Analysis: 8/6-7/2018
Date of Initial Calibration: 5/18/2018, 7/25/2018
SDG #: E18-06141

File #	Laboratory Sample ID	QC Check	Dilution Factor	Can #	Analyst	Injection Volume (cc)	Comments	Make-up Air		Acquisition		Room		TO-15 Standard	
								Added to canister (cc)	Added for dilution (cc)	Date	Time	Temp	BP "Hg	Working ID	Vendor ID (Lot#)
aa7971bfb	BFB	✓		ALM029426	JLS	50				7/25/2018	8:35	75	30.12		403-402623
aa7972std01	40 ppbv Std	✓		CC483586	JLS	200								7/25/2018	160-401140945-1
aa7973std02	20 ppbv Std	✓		CC483586	JLS	100								7/25/2018	160-401140945-1
aa7974std03	10 ppbv Std	✓		CC483586	JLS	50								7/25/2018	160-401140945-1
aa7975std04	2 ppbv Std	✓		CC483586	JLS	10								7/25/2018	160-401140945-1
aa7976std05	0.2 ppbv Std	✓		CC483586	JLS	1								7/25/2018	160-401140945-1
aa7977icvss	10 ppbv ICVSS	✓		CC483422	JLS	50									160-401186845-1
aa7978	E18-05039-01 x 100	✓	100	5093	JLS	5	not used								
aa7979	E18-05039-02 x 100	✓	100	3033	JLS	5	not used								
aa7980	E18-05040-01 x 100	✓	100	1769	JLS	5	not used								
aa7981	E18-05041-01 x 100	✓	100	1523	JLS	5	not used								
aa7982	E18-05041-02 x 100	✓	100	1599	JLS	5	not used								

Analyst Name: Jeff Schmitt

Signature: *Jeff Schmitt*

IAL SDG #E18-06141

Supervisor Name: Lauren Jenkins

Signature: *Lauren Jenkins*

**Integrated Analytical Laboratories
GC/MS Run Log**

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Date of Analysis: 8/6-7/2018
Date of Initial Calibration: 5/18/2018, 7/25/2018
SDG #: E18-06141

File #	Laboratory Sample ID	QC Check	Dilution Factor	Can #	Analyst	Injection Volume (cc)	Comments	Make-up Air		Acquisition		Room		TO-15 Standard	
								Added to canister (cc)	Added for dilution (cc)	Date	Time	Temp	BP "Hg	Working ID	Vendor ID (Lot#)
aa8191bfb	BFB	✓		ALM029426	JLS	50				8/6/2018	11:26	75	30.20		403-402623
aa8192dcvs	10 ppbv DCVS	✓		AAL071156	JLS	50								7/25/2018	160-401140945-1
aa8193rlcs	0.2 ppbv RLLCS	✓		AAL071156	JLS	1								7/25/2018	160-401140945-1
aa8194blk	Method Blank	✓		1127	JLS	500									
aa8195	E18-06204-01	✓		3280	JLS	500									
aa8196	E18-06204-21	✓	dup of E18-06204-01, c		JLS	500									
aa8197	E18-06204-02 x 10	✓	10	1548	JLS	50									
aa8198	blk	✓			JLS	500									
aa8199	E18-05043-04 x 10	✓	10	2885	JLS	50									
aa8200	E18-05043-04	✓		2885	JLS	500									
aa8201	E18-05043-05 x 10	✓	10	2065	JLS	50									
aa8202	E18-05043-05	✓		2065	JLS	500									
aa8203	E18-05043-06 x 10	✓	10	4866	JLS	50	not used								
aa8204	E18-05043-06	✓		4866	JLS	500									
aa8205	blk	✓			JLS	500									
aa8206	E18-06141-07	✓		3014A	JLS	500									
aa8207	E18-06141-08	✓		2155	JLS	500									
aa8208	E18-06141-09	✓		5089	JLS	500									
aa8209	blk	✓			JLS	500									
aa8210	E18-06173-01 x 100	✓	100	1070	JLS	5	not used								
aa8211	E18-06173-02 x 100	✓	100	1764	JLS	5	not used								
aa8212	blk	✓			JLS	500									
aa8213ccvcs	10 ppbv CCCVS	✓		CC483586	JLS	50				8/7/2018	2:22			7/25/2018	160-401140945-1

Analyst Name: Jeff Schmitt

Signature: *Jeff Schmitt*

IAL SDG #E18-06141

Supervisor Name: Lauren Jenkins

Signature: *Lauren Jenkins*

**Integrated Analytical Laboratories
GC/MS Run Log**

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Column: Restek RTX-1 SN 992567

Target Directory: D:\Agilent GCMS\

Date of Analysis: 8/6-7/2018
Date of Initial Calibration: 5/18/2018, 7/25/2018
SDG #: E18-06141

File #	Laboratory Sample ID	QC Check	Dilution Factor	Can #	Analyst	Injection Volume (cc)	Comments	Make-up Air		Acquisition		Room		TO-15 Standard	
								Added to canister (cc)	Added for dilution (cc)	Date	Time	Temp	BP "Hg	Working ID	Vendor ID (Lot#)
aa8221bfb	BFB	✓		ALM029426	JJW	50				8/7/2018	9:33	74	30.11		403-402623
aa8222dcvs	10 ppbv DCVS	✓		AAL071156	JJW	50								7/25/2018	160-401140945-1
aa8223rlcs	0.2 ppbv RLLCS	✓		AAL071156	JJW	1								7/25/2018	160-401140945-1
aa8224blk	Method Blank	✓		1127	JJW	500									
aa8225	E18-06204-02 x 100	✓	100	1548	JJW	5	not used								
aa8226	E18-06204-02 x 10	✓	10	1548	JJW	50									
aa8227	E18-06204-02 x 20	✓	20	1548	JJW	25									
aa8228	E18-06173-01 x 10	✓	10	1070	JJW	50									
aa8229	E18-06173-21 x 10	✓	10	06173-01 x 10	JJW	50									
aa8230	E18-06173-02 x 10	✓	10	1764	JJW	50									
aa8231	blk	✓			JJW	500									
aa8232	E18-06141-08 x 100	✓	100	2155	JJW	5									
aa8233	E18-06141-09 x 10	✓	10	5089	JJW	50									
aa8234	blk	✓			JJW	500									
aa8235	E18-06174-01 x 100	✓	100	1603	JJW	5	not used								
aa8236	E18-06174-02 x 100	✓	100	1779	JJW	5									
aa8237	E18-06174-03 x 100	✓	100	1600	JJW	5									
aa8238	E18-06174-04 x 100	✓	100	1570	JJW	5	not used								
aa8239	E18-06174-05 x 100	✓	100	1400	JJW	5	not used								
aa8240	E18-06174-06 x 100	✓	100	1521	JJW	5									
aa8241	blk	✓			JJW	500									
aa8242cccvs	10 ppbv CCCVS	✓		CC483586	JJW	50				8/7/2018	2:22			7/25/2018	160-401140945-1

Analyst Name: Jeff Schmitt

Signature: *Jeff Schmitt*

IAL SDG #E18-06141

Supervisor Name: Lauren Jenkins

Signature: *Lauren Jenkins*

Client Contact Information			Project Information					Carrier (check one): <input checked="" type="checkbox"/> IAL Courier <input type="checkbox"/> Client Courier <input type="checkbox"/> FedEx/UPS				pg <u>1</u> of <u>2</u>															
Company: <u>BEI</u>			Project Name: <u>Congers C.P. / 060141</u>					Invoice Information				Analysis		Report		Matrix											
Address: <u>19 Chatham Rd. Summit, NJ 07901</u>			Project Location (State): <u>NY</u>					Attn:				EPA TO-15 NJDEP LLTO-15 (includes 30 TICs) Library Search (10, 20, or 30 TICs) Other (Explain in Comments) Regulatory/ NY Cat B / Full (NJ Required) Reduced / NY Cat A Data Package Results Only Indoor Air Ambient / Outdoor Air Sub Slab / Soil Gas / Near Slab (Circle One) Stack Emission / SVE System High Concentrations Expected															
Phone: <u>908-918-1702</u>			Project Manager: <u>Jeff McCurdy</u>					Address:																			
Fax: <u>908-918-1707</u>			PM Signature: <u>[Signature]</u>					PO #:																			
Report to:			PM E-Mail: <u>Jmccurdy@bei-enr.com</u>					Quote #:																			
Sampler: <u>AS</u>																											
Analysis Turnaround Time - IF NO TAT IS SPECIFIED, 2 WEEK TAT IS ASSUMED										Barometric Pressure																	
IAL Standard: 2 weeks (10 business days)										Start		<u>30.10</u>		Stop		<u>30.10</u>											
Rush (**pre-approved by lab):										24hr**		48hr**		72hr**		96hr**		1wk**									
Sample Identification	Start DATE & TIME (24hr Clock)	End DATE & TIME (24hr Clock)	Starting Vacuum ("Hg)	Ending Vacuum ("Hg)	Starting Temp. ("F)	Ending Temp. ("F)	Outgoing Vacuum - Lab ("Hg)	Incoming Vacuum - Lab ("Hg)	Flow Regulator ID	Canister ID	Canister Size (1L or 6L)	Flow Controller Readout (cc/min)	EPA TO-15	NJDEP LLTO-15 (includes 30 TICs)	Library Search (10, 20, or 30 TICs)	Other (Explain in Comments)	Regulatory/ NY Cat B / Full (NJ Required)	Reduced / NY Cat A Data Package	Results Only	Indoor Air	Ambient / Outdoor Air	Sub Slab / Soil Gas / Near Slab (Circle One)	Stack Emission / SVE System	High Concentrations Expected			
1) AA-101	7/31/18 935	7/31/18 408	-30	-2.0	72	92	-29.0	-2.0	7342673	5076	6L	12.4	H				X					X					
2) IA-FF-101	837	411	-29.5	-2.5	69	69	-29.0	-2.5A	0702495-3	3013	6L	12.3	H				X				X						
3) IA-FF-102	839	355	-28.5	-2.0	68	69	-29.0	-2.0	7342708	3814	6L	12.4	H				X				X						
4) IA-FF-103	844	412	-30	-2.5	68	69	-29.0	-2.5A	0302495-5	2753	6L	12.5	H				X				X						
5) IA-B-101	840	348	-29	-2.5	69	68	-29.0	-2.5A	0121687-1	3277	6L	12.5	H				X				X						
6) IA-B-102	842	410	-29.5	-2.5	68	68	-29.0	-2.5	7340286	3026A	6L	12.6	H				X				X						
Comments/ Special Analysis Instructions / QC Requirements:													Note: Hold or contingent samples may be designated by writing an "H" or "C" in the appropriate analysis box. ALL FIELDS IN RED ARE REQUIRED														
3013, 2753, 3026A 06081801 06131822 3059 H = hold X = run Report VOCs only 5076, 3814, 3277 05221801 05291822 5078																											
Shipping Information / Canister Preparation (for laboratory use only)										Laboratory Canister Certification																	
Individual Preparing Canisters / Title: <u>P. Jenkins, J. Walukiewicz / Air Department Sample Custodians</u>										GC/MS Analyst Signature																	
Lab Affixed Seal Number(s): <u>IAL-20180200, -0201, -0202</u>										[Signature] IAL SDG#: <u>06141</u>																	
Date/Time Shipping Container Sealed: <u>7-26-18 1130</u>																											
External Chain of Custody																											
Relinquished			Received			Date / Time			Reason for Change of External Custody																		
[Signature]			[Signature]			7-26-18 1130			shipment from laboratory to client Rec'd @ air lab																		
[Signature]			[Signature]			8/21/18 1606 8/2/18 1659																					
Name/Title Resealing Shipping Container Name: _____										NJDEP Affixed Seal Number: _____																	
Date/Time Sample Shipping Container Resealed: _____										Individual Opening Sample Shipping Container: <u>Padraic Jenkins / Joseph Walukiewicz</u>																	
Date/Time Sample Shipping Container Opened: _____										Date/Time Internal Chain of Custody Initiated: _____																	

Client Contact Information				Project Information				Carrier (check one): <input checked="" type="checkbox"/> IAL Courier <input type="checkbox"/> Client Courier <input type="checkbox"/> FedEx/UPS				pg <u>2</u> of <u>2</u>													
Company: <u>BEI</u>				Project Name: <u>Congers C.P./066141</u>				Invoice Information				Analysis		Report		Matrix									
Address: <u>19 Chatham Rd. Summit, NJ 07901</u>				Project Location (State): <u>NY</u>				Attn:				EPA TO - 15 NJDEP LLTO-15 (Includes 30 TICs) Library Search (10, 20, or 30 TICs) Other (Explain in Comments) Regulatory/ NY Cat B / Full (NJ Required) Reduced / NY Cat A Data Package Results Only Indoor Air Ambient / Outdoor Air Sub Slab / Soil Gas / Near Slab (Circle One) Stack Emission / SVE System High Concentrations Expected													
Phone: <u>908-918-1702</u>				Project Manager: <u>Jeff McCurdy</u>				Address:																	
Fax: <u>908-918-1707</u>				PM Signature: <u>[Signature]</u>				PO #:																	
Report to:				PM E-Mail: <u>jmccurdy@be-enr.com</u>				Quote #:																	
Sampler: <u>AS</u>																									
Analysis Turnaround Time - IF NO TAT IS SPECIFIED, 2 WEEK TAT IS ASSUMED										Barometric Pressure															
IAL Standard: 2 weeks (10 business days)										Start		30.10		Stop		30.10									
Rush (**pre-approved by lab):										24hr**		48hr**		72hr**		96hr**		1wk**							
Sample Identification	Start DATE & TIME (24hr Clock)	End DATE & TIME (24hr Clock)	Starting Vacuum ("Hg)	Ending Vacuum ("Hg)	Starting Temp. (°F)	Ending Temp. (°F)	Outgoing Vacuum - Lab ("Hg)	Incoming Vacuum - Lab ("Hg)	Flow Regulator ID	Canister ID	Canister Size (1L or 6L)	Flow Controller Readout (cc/min)													
7) 55-101	8/1/18 917	8/1/18 415	-29.0	-2.5	68	68	-29.0	-2.5	A00988641-6	3014A	6L	12.7	X				X								
8) 55-102	↓ 833	↓ 419	-30.0	-2.5	68	68	-29.0	-2.5	A0302498-2	2155	6L	12.6	X				X								
9) 55-103	↓ 839	↓ 411	-30.0	-2.5	68	69	-29.0	-2.5	A0302497	5089	6L	12.5	X				X								
Comments/ Special Analysis Instructions / QC Requirements:												Note: Hold or contingent samples may be designated by writing an "H" or "C" in the appropriate analysis box. ALL FIELDS IN RED ARE REQUIRED													
3014A, 2155, 5089 06091801 061318aa 3059 H=hold X=run Report CVOCS only																									
Shipping Information / Canister Preparation (for laboratory use only)								Laboratory Canister Certification																	
Individual Preparing Canisters / Title: P. Jenkins, J. Walukiewicz / Air Department Sample Custodians								GC/MS Analyst Signature																	
Lab Affixed Seal Number(s): <u>IAL-20180200, -0201, -0202</u>								[Signature] IAL SDG#: <u>06141</u>																	
Date/Time Shipping Container Sealed: <u>7-26-18 1130</u>																									
External Chain of Custody																									
Relinquished				Received				Date / Time				Reason for Change of External Custody													
[Signature]				[Signature]				7-26-18 1130				shipment from laboratory to client													
								8/2/18 1606				Rec'd @ air lab													
								8/2/18 1659																	
Name/Title Resealing Shipping Container Name:								NJDEP Affixed Seal Number:																	
Date/Time Sample Shipping Container Resealed:								Individual Opening Sample Shipping Container: Padraic Jenkins / Joseph Walukiewicz																	
Date/Time Sample Shipping Container Opened:								Date/Time Internal Chain of Custody Initiated:																	

Example Calculation (EPA TO-15)

$$\frac{\text{Area of Sample}}{\text{Area of Internal Standard}} \times \frac{\text{Concentration of Internal Standard (10 ppbv)}}{\text{Response Factor}} = \text{Concentration of Sample (ppbv)}$$

Conversion from ppbv to $\mu\text{g}/\text{m}^3$

$$\frac{\text{Concentration of Compound (ppbv)}}{24.45} \times \text{Molecular Weight of Compound} = \text{Concentration of Compound } (\mu\text{g}/\text{m}^3)$$

Clean Canister Certification Report

Lab Sample Name: Clean Canister, Batch Master 3059
Field Sample Name: Canister 3059
Sample Volume: 500ml

Data File: AA7475
Date Analyzed: 6/13/2018
Matrix: Air

Canisters associated with this run: 3059,2155(used for E18-06141-08),2164,2753,3013,3014A(used for E18-06141-07),3026A,5089(used for E18-06141-09)

Runs with this Clean Canister Certification:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7471BFB]	06/13/2018 08:30
10 PPBV DCVS [AA7472DCVS]	06/13/2018 09:18
METHOD BLANK [AA7473BLK]	06/13/2018 10:25
02 PPBV RLLCS [AA7474RLLCS]	06/13/2018 11:14
CLEAN CAN CERTIFICATION, BATCH MASTER 3059 [AA7475]	06/13/2018 12:20
10 PPBV CCCVS [AA7485CCCVS]	06/13/2018 20:47

This canister has been certified clean, all compounds are below 0.2 ppbv.

Compound	CAS #	RL (ppbv)	Calculated Amount (ppbv)
Acetone	67-64-1	0.20	ND
Benzene	71-43-2	0.20	ND
Bromodichloromethane	75-27-4	0.20	ND
Bromoform	75-25-2	0.20	ND
Bromomethane	74-83-9	0.20	ND
1,3-Butadiene	106-99-0	0.20	ND
Chlorobenzene	108-90-7	0.20	ND
Chloroethane	75-00-3	0.20	ND
Chloroform	67-66-3	0.20	ND
Chloromethane	74-87-3	0.20	ND
Carbon disulfide	75-15-0	0.20	ND
Carbon tetrachloride	56-23-5	0.20	ND
Cyclohexane	110-82-7	0.20	ND
Dibromochloromethane	124-48-1	0.20	ND
1,2-Dibromoethane	106-93-4	0.20	ND
1,2-Dichlorobenzene	95-50-1	0.20	ND
1,3-Dichlorobenzene	541-73-1	0.20	ND
1,4-Dichlorobenzene	106-46-7	0.20	ND
Dichlorodifluoromethane	75-71-8	0.20	ND
1,1-Dichloroethane	75-34-3	0.20	ND
1,2-Dichloroethane	107-06-2	0.20	ND
1,1-Dichloroethene	75-35-4	0.20	ND
1,2-Dichloroethene (cis)	156-59-2	0.20	ND
1,2-Dichloroethene (trans)	156-60-5	0.20	ND
1,2-Dichloropropane	78-87-5	0.20	ND
1,3-Dichloropropene (cis)	10061-01-5	0.20	ND
1,3-Dichloropropene (trans)	10061-02-6	0.20	ND
1,2-Dichlorotetrafluoroethane	76-14-2	0.20	ND
1,4-Dioxane	123-91-1	0.20	ND
Ethylbenzene	100-41-4	0.20	ND
n-Heptane	142-82-5	0.20	ND
1,3-Hexachlorobutadiene	87-68-3	0.20	ND
n-Hexane	110-54-3	0.20	ND
Methylene chloride	75-09-2	0.20	ND
Methyl ethyl ketone	78-93-3	0.20	ND

Clean Canister Certification Report

Lab Sample Name: Clean Canister, Batch Master 3059
Field Sample Name: Canister 3059
Sample Volume: 500ml

Data File: AA7475
Date Analyzed: 6/13/2018
Matrix: Air

Canisters associated with this run: 3059,2155(used for E18-06141-08),2164,2753,3013,3014A(used for E18-06141-07),3026A,5089(used for E18-06141-09)

Runs with this Clean Canister Certification:

Standard/Sample Run	Date/Time of Sample/Standard Injection
BFB [AA7471BFB]	06/13/2018 08:30
10 PPBV DCVS [AA7472DCVS]	06/13/2018 09:18
METHOD BLANK [AA7473BLK]	06/13/2018 10:25
02 PPBV RLLCS [AA7474RLLCS]	06/13/2018 11:14
CLEAN CAN CERTIFICATION, BATCH MASTER 3059 [AA7475]	06/13/2018 12:20
10 PPBV CCCVS [AA7485CCCVS]	06/13/2018 20:47

This canister has been certified clean, all compounds are below 0.2 ppbv.

Compound	CAS #	RL (ppbv)	Calculated Amount (ppbv)
Methyl isobutyl ketone	108-10-1	0.20	ND
Methyl tert-butyl ether	1634-04-4	0.20	ND
Styrene	100-42-5	0.20	ND
Tert-butyl alcohol	75-65-0	0.20	ND
1,1,2,2-Tetrachloroethane	79-34-5	0.20	ND
Tetrachloroethene	127-18-4	0.20	ND
Toluene	108-88-3	0.20	ND
1,2,4-Trichlorobenzene	120-82-1	0.20	ND
1,1,1-Trichloroethane	71-55-6	0.20	ND
1,1,2-Trichloroethane	79-00-5	0.20	ND
Trichloroethene	79-01-6	0.20	ND
Trichlorofluoromethane	75-69-4	0.20	ND
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	0.20	ND
1,2,4-Trimethylbenzene	95-63-6	0.20	ND
1,3,5-Trimethylbenzene	108-67-8	0.20	ND
2,2,4-Trimethylpentane	540-84-1	0.20	ND
Vinyl bromide	593-60-2	0.20	ND
Vinyl chloride	75-01-4	0.20	ND
Xylenes (m&p)	179601-23-1	0.40	ND
Xylenes (o)	95-47-6	0.20	ND

Data Path : C:\DATA\06-13-18\
 Data File : aa7475.D
 Acq On : 13 Jun 2018 12:20 pm
 Operator : jls
 Sample : 3059
 Misc : 2155, 2164, 2753, 3013, 3014A, 3026A, 5089
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Jun 18 08:48:05 2018
 Quant Method : C:\msdchem\1\METHODS\0518.M
 Quant Title : TO-15 on the Agilent 7890A / 5975C
 QLast Update : Fri May 18 13:51:08 2018
 Response via : Initial Calibration

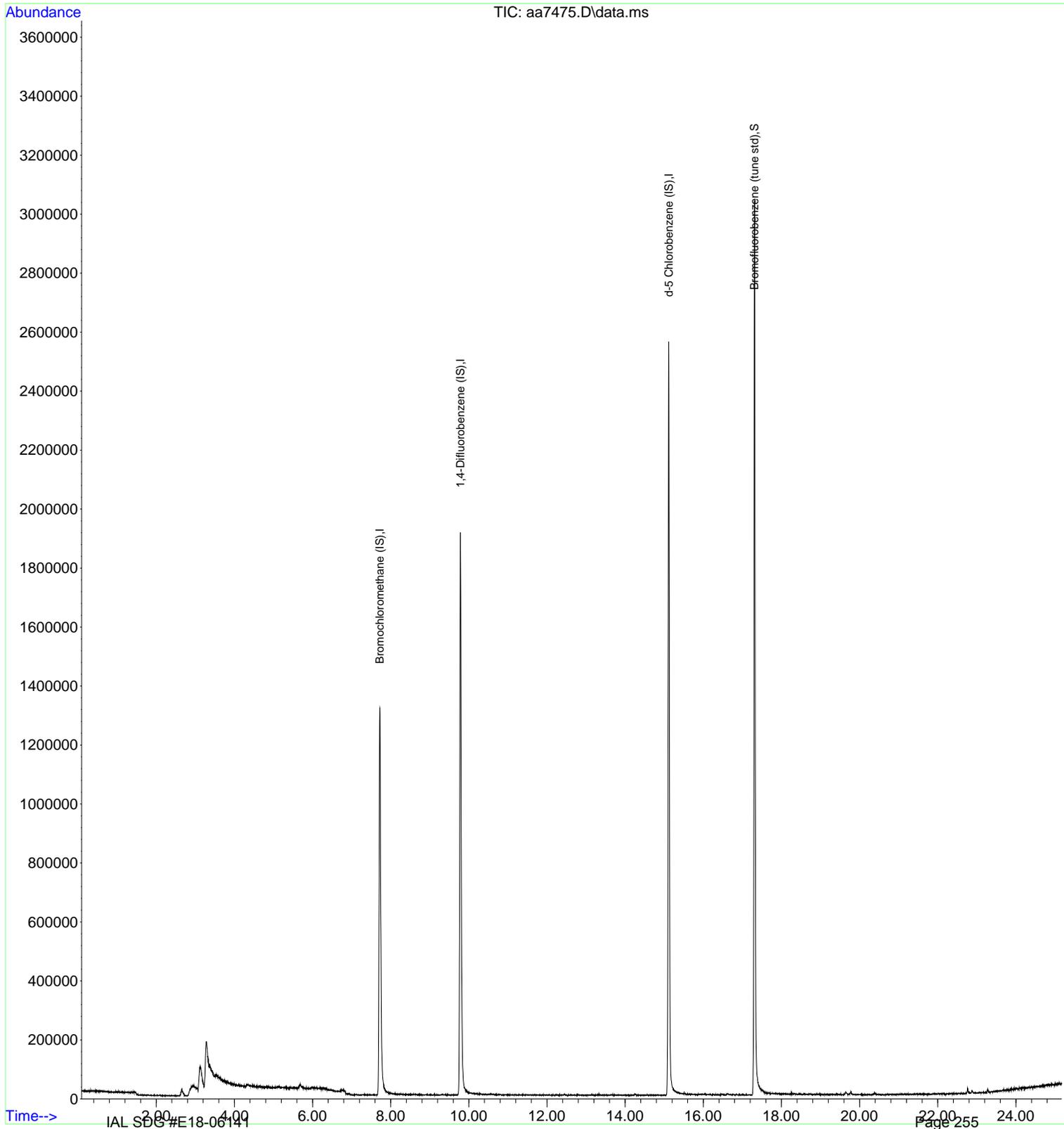
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Bromochloromethane (IS)	7.724	130	486608	10.00	ppbV	0.00
38) 1,4-Difluorobenzene (IS)	9.785	114	1675394	10.00	ppbV #	0.00
55) d-5 Chlorobenzene (IS)	15.116	117	1426411	10.00	ppbV	0.00
System Monitoring Compounds						
64) Bromofluorobenzene (tu...	17.312	95	1270032	9.67	ppbV	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\DATA\06-13-18\
Data File : aa7475.D
Acq On : 13 Jun 2018 12:20 pm
Operator : jls
Sample : 3059
Misc : 2155, 2164, 2753, 3013, 3014A, 3026A, 5089
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Jun 18 08:48:05 2018
Quant Method : C:\msdchem\1\METHODS\0518.M
Quant Title : TO-15 on the Agilent 7890A / 5975C
QLast Update : Fri May 18 13:51:08 2018
Response via : Initial Calibration



LAST PAGE OF DOCUMENT