

# SOIL VAPOR INTRUSION INVESTIGATION REPORT; 690 Saint Paul Street – Off-Site NYSDEC SITE #C828159A

## Location:

691 and 705 Saint Paul Street  
Rochester, New York

## Prepared for:

Bausch & Lomb  
1400 North Goodman Street  
Rochester, New York 14609

LaBella Project No. 2170820

August 20, 2019

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## CERTIFICATIONS

*"I Daniel P. Noll certify that I am currently a NYS registered professional engineer and that this Soil Vapor Intrusion Investigation Report was prepared in accordance with all applicable statutes and regulations and in substantial conformance with the DER Technical Guidance for Site Investigation and Remediation (DER-10)."*



081996

8/20/19  
~~10/20/2012~~

NYS Professional Engineer #

Date

Signature

## 1.0 INTRODUCTION

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As required under the terms of the New State Department of Environmental Conservation (NYSDEC) Order of Consent and Administrative Settlement (Index No. R8-20161013-107) with Bausch & Lomb (B&L), this report details the findings of a Soil Vapor Intrusion (SVI) investigation that was performed by LaBella Associates, D.P.C. on behalf of B&L for the properties at 691 and 705 Saint Paul Street, City of Rochester, Monroe County, New York and identified by the NYSDEC as 690 Saint Paul Street – Off-Site, Site #C828159A. A site location map is included as Figures 1 and 2. The SVI investigation activities were conducted in accordance with a NYSDEC approved work plans entitled:

- *Soil Gas Sampling Work Plan, NYSDEC Site #828159A, 691 and 705 Saint Paul Street, Rochester, New York* dated January 30, 2017 prepared by LaBella Associates, D.P.C.
- *Soil Vapor Intrusion Work Plan, NYSDEC Site #828159A, 691 and 705 Saint Paul Street, Rochester, New York* dated March 24, 2017 prepared by LaBella Associates, D.P.C.

This SVI Report documents soil gas and SVI sampling at the site that is west of the 690 Saint Paul Street NYSDEC Brownfield Cleanup Program (BCP) Site #C828159. The SVI sampling was conducted to determine whether there is a SVI concern at the site with regard to the chlorinated solvent plumes associated with the 690 Saint Paul Street NYSDEC BCP Site #C828159. This report was also completed in accordance with the NYSDEC Division of Environmental Remediation (DER) BCP Guide dated May 2004 and the DER-10 (*Technical Guidance for Site Investigation and Remediation*) dated May 3, 2010 and the New York State Department of Health (NYSDOH) Final document “Guidance for Evaluating Soil Vapor Intrusion in the State of New York,” dated October 2006.

### 1.1 Report Organization

The remainder of this SVI Report is organized into the sections listed below.

- **Section 2** provides a general description of the site and surrounding properties and an environmental summary of the 690 Saint Paul Street BCP Site #C828159
- **Section 3** provides a description of the subsurface geologic conditions in the area of the site
- **Section 4** provides a description of the Standards, Criteria, and Guidance that are used to define the soil vapor cleanup guidelines for the site
- **Section 5** provides a description of field activities, sampling, and analysis performed during the SVI investigation
- **Section 6** provides a summary of the analytical results of samples collected and analyzed during the SVI investigation
- **Section 7** provides a conclusions of the results of the SVI work completed to date
- **Section 8** provides a recommendation with regard to potential mitigation measures at the site



## 2.0 SITE DESCRIPTION AND ENVIRONMENTAL SUMMARY

### 2.1 Site Description

The Site is located directly west of the NYSDEC BCP site #C828159 known as 690 Saint Paul Street. The 691 Saint Paul Street property is primarily used as commercial office space by: 1) Monroe County Social Services (primary tenant) that occupies the upper basement and the remaining upper floors (2) St. Michael's Woodshop occupies most of the lower basement space, and (3) Newport Gratings (photonics company) occupies the northern end of the upper and lower basement space. The 705 Saint Paul Street property is presently used as manufacturing by Richardson Gratings. Each property is improved with one building as outlined in Table 1.1 below.

TABLE 1.1 – Site Buildings Description

	691 Saint Paul Street	705 Saint Paul Street
Acreage of Site	2.94	0.82
Approximate Building Footprint (square feet)	30,630	10,627
Foundation Type	Basement	Unknown
Construction Date	1920	1930
Current Use	Commercial office space	Manufacturing

The basement of the 691 Saint Paul Street building is divided into two separate levels or elevations. The eastern half of the upper basement is approximately 8 to 10 ft higher in elevation than the lower basement that is located on the approximate western half of the site building as shown on Figures 4A and 4B.

The exterior of the 691 Saint Paul Street parcel primarily consists of an asphalt paved parking lot located south of the building and the exterior of the 705 Saint Paul Street parcel primarily consists of an asphalt paved parking lot located north of the building. Neighboring properties that border the site include:

TABLE 1.1B - Adjacent Properties

Location	Current Use
North	Asphalt paved parking lot
South	Smith Street, High Fall Brewery
East	690 Saint Paul Street, NYSDEC BCP Site# C828159
West	Suntru Street, RG&E Substation #34

## **2.2 Environmental Summary of NYSDEC 690 Saint Paul Street BCP Site #C828159**

The 690 Saint Paul Street BCP site was developed prior to 1875 and was utilized primarily for residential purposes prior to approximately the 1920s. Based on the review of historical mapping and local street directories, the BCP site was primarily utilized for industrial purposes by Bausch & Lomb, Inc., formerly known as Bausch & Lomb Optical Company (“B&L”) from sometime around 1920 until it was abandoned by the company in the late 1960s. The property was developed for industrial use by B&L to manufacture lenses and other products. From the early 1970s until 2000, the BCP site was used predominantly for light commercial and storage applications. Occupants and/or owners of the Site have included various individual residences, B&L, Thomas Edison Technical and Industrial High School, Geva Theater storage, and various manufacturing and industrial tenants.

Prior uses of the BCP site that appear to have led to site contamination include underground storage tanks (USTs) that may have leaked. These tanks appear to have contained chlorinated solvents including trichloroethene (TCE) and petroleum products including gasoline and fuel oil. In 2002, a 500-gallon UST was removed from the site and contaminated soil was encountered. In 2008, a Phase II Environmental Site Assessment (ESA) was performed to evaluate subsurface soil and groundwater. The Phase II ESA identified an area of petroleum contaminated soil. The investigation was followed by the excavation of approximately 1,650 cubic yards of petroleum impacted soil and a previously undocumented UST. TCE was identified in confirmation samples in soil and ground and thus the 690 Saint Paul Street site was entered into the BCP. A Remedial Investigation (RI) conducted through the BCP identified eight (8) Areas of Concern that related to elevated concentrations of metals, polychlorinated biphenyls (PCBs), polycyclic aromatic hydrocarbons, petroleum products, and/or chlorinated solvents. Interim Remedial Measures (IRMs) were completed after these and the Remedial Investigation (RI) work it was determined three areas required further remedial actions. One of these areas was petroleum product (light non-aqueous phase liquid) which is limited to within the 690 Saint Paul Street parcel. The other two locations contain chlorinated VOCs which were document to have off-site migration.

The areas of concern related to this report are two groundwater chlorinated solvent plumes that were identified to have migrated off the BCP site to the west and toward the Site as shown on Figure 3. It should be noted that the 690 Saint Paul Street site is currently undergoing remediation, including the chlorinated VOC impacts.

## **3.0 SUMMARY OF LOCAL SUBSURFACE GEOLOGIC CONDITIONS**

This section relies upon information obtained from the 690 Saint Paul Street NYSDEC BCP site environmental investigations. The overburden material ranges in depth up to approximately 12 feet (ft) below the ground surface (bgs) and consists primarily of glacial till. Fill material is anticipated to overly the till and may consist of sand, gravel, cinders, and ash.

The Decew Dolomite underlies the overburden material. The Decew Dolomite is the uppermost formation of the Clinton Group and consists of variably bedded, dark-gray to olive-gray, argillaceous to sandy, fine-grained dolomite that contains shaly partings and interbeds, as well as frequent pits and vugs. The thickness of this unit is generally 8 to 12 ft.

The Rochester Shale underlies the Decew Dolomite, and is a relatively uniform dark- to medium-gray, pale- and platy-weathering, highly calcareous to dolomitic mudstone. It contains abundant thin interbeds of medium gray, pale-buff weathering, laminated calcisiltites. Its thickness in Western New York is generally 58 to 65 ft.

The overburden groundwater table at the BCP site ranged from 4 to 9 ft bgs in the southern portion of the site and 8 to 10 ft bgs in the northern portion of the site. The overburden groundwater flows generally to the west-southwest.

The shallow bedrock water-bearing interval was identified as the uppermost bedrock down to depths of approximately 20 ft bgs. This interval is the uppermost water-bearing unit within the bedrock, and no low permeability horizon separates this zone from the overburden. Groundwater flow direction is generally to the west.

#### 4.0 STANDARDS CRITERIA AND GUIDELINES

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This section identifies the Standards, Criteria, and Guidance (SCGs) for the site. The SCGs identified are used in order to quantify the extent of contamination at the site that may require remedial work. The SCGs will be used to evaluate the effectiveness of any remedial measures and will be used to determine if remedial actions are warranted. The SCG's for soil vapor are provided below:

- **Soil Gas SCGs:** Currently there are no state regulatory (NYSDEC or NYSDOH) soil gas guidance values.
- **Sub-Slab Soil Vapor and Indoor Air SCGs:** The NYSDOH Final document "Guidance for Evaluating Soil Vapor Intrusion in the State of New York," dated October 2006 and updates for tetrachloroethene (September 2013), TCE (August 2015), and the soil vapor/indoor air decision matrices (May 2017). These updates are included on NYSDOH website: [https://www.health.ny.gov/environmental/indoors/vapor\\_intrusion/update.htm](https://www.health.ny.gov/environmental/indoors/vapor_intrusion/update.htm).

#### 5.0 SOIL VAPOR INTRUSION INVESTIGATION FIELD ACTIVITIES

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This section provides a description of the methodologies used during the field investigations of the site. The SVI work was completed from March to April 2017 and was conducted in accordance with the NYSDEC approved work plans. Specific tasks performed during the SVI investigation included the following:

- Soil gas and sub-slab vapor point installation, sampling, and laboratory analysis
- Quality Assurance/Quality Control

### **5.1 Soil Gas Point Installation, Sampling, and Laboratory Analysis**

The purpose of this activity was to determine if subsurface impacts from the 690 Saint Paul Street site have the potential to adversely affect indoor air quality via the vapor intrusion pathway at the site. The locations of the soil gas samples were located within the Saint Paul Street right-of-way and directly across from the chlorinated solvent plumes at the 690 Saint Paul Street site. Soil gas sample locations are shown on Figure 3, whereas soil gas sample locations SV-1 and SV-2 are located adjacent to the 705 Saint Paul Street site building and soil gas sample locations SV-3 and SV-4 are located adjacent to the southern end of the 691 Saint Paul Street site building. Sampling of soil gas was conducted on March 3, 2017.

Permanent soil gas sampling points were installed using direct push technology. A 1 foot screen section was installed to a depth of approximately 8 feet (ft) below the ground surface (bgs), which is approximately the top of bedrock and approximately just below the floor of the upper basement of the 691 Saint Paul Street site. A solid PVC riser was installed above the 1 ft screen to the ground surface. A porous inert backfill material (i.e. sand) was placed around the sampling screen to approximately 1 ft above the screened section to create a sampling zone of 2 ft in length. The soil gas sampling points were sealed above the sampling zone with approximately 7 ft of a bentonite slurry (to just below the ground surface) and finished with a protective casing that was grouted in place to minimize infiltration of water, and to prevent damage to the soil gas point. The soil gas installation log is included in Appendix 1.

Prior to sampling each soil gas point, a tracer gas (helium) was utilized at each sampling point to ensure that ambient air was not being pulled into the Summa® canister during sampling. This was accomplished by placing a clean, stainless steel enclosure over the soil vapor sampling points. Non emitting VOC modeling clay was placed on the ground surface around the edge of the enclosure where it contacted the ground to make an air-tight seal. Prior to purging and sampling activities, a helium tracer gas was released via a small diameter tube, placed through a port on the exterior side of the enclosure. The sub-slab vapor Teflon® tubing extended up through the air-tight seal to the exterior side of the enclosure and connected to a helium detector to determine the presence of helium gas and purge the sampling point. Helium was not detected on the helium detector at each sampling point indicating that each soil gas sampling point were sealed from ambient air. All sample points passed the trace gas test (i.e. less than 10% of helium was detected during each test).

Care was taken to avoid excessive purging of the samples points and the flow rate during purging did not exceed 0.2 liters per minute (L/min) to minimize outdoor air infiltration. Soil gas points were purged with a PID.

Four soil gas and one outdoor ambient air samples were collected utilizing individually certified-clean 1-liter Summa® canisters equipped with laboratory calibrated flow controllers. The samples were collected over an approximate eight (8) hour time period. The outdoor air sample was collected at a height of approximately 5 ft above the ground surface to simulate the breathing zone. Immediately after opening each Summa® canister, the initial vacuum (inches of mercury) and time was noted and recorded on the laboratory chain-of-custody. After approximately eight (8) hours, final vacuum readings (inches of mercury) were noted and the Summa® canisters were closed. A copy of the Air Sampling Field Report for each sample collected is included in Appendix 1.

The samples (including QA/QC samples) were submitted to Centek Laboratories, LLC in Syracuse, New York. Centek is a New York State Department of Health (NYSDOH) Environmental Laboratory Approval Program certified laboratory for analysis of targeted volatile organic compounds (VOCs) by United States Environmental Protection Agency (USEPA) Method TO-15.

## **5.2 Sub-Slab Sample Point Installation, Soil Vapor Intrusion Sampling, and Laboratory Analysis**

The purpose of this activity was to assess for potential SVI at the 691 Saint Paul Street building from subsurface impacts at the 690 Saint Paul Street site. LaBella conducted SVI sampling at the locations shown on Figure 4A (upper basement) and 4B (lower basement). SVI sampling was conducted on April 1, 2017. A copy of the Air Sampling Field Report for each sample collected and the NYSDOH Indoor Air Quality and Building Inventory Form are included in Appendix 1.

Sub slab vapor sample points were collected by coring a nominal 5/8 inch diameter hole through the slab to approximately 1 to 2 inches beneath the floor slab. A 5/8 inch diameter polyethylene tube and barbed fitting was inserted into the corehole. The tubing was connected to a Summa® canister with a pre-set regulator to the barbed fitting for sub-slab vapor collection.

Prior to sampling each sub-slab vapor point, a tracer gas (helium) was placed over each sampling point to ensure that ambient air was not being pulled into the Summa® canister (i.e. sampling vessel) during sampling. This was accomplished by placing a clean, stainless steel enclosure over each sub-slab sampling point. Non-emitting VOC modeling clay or equivalent was placed on the ground surface around the edge of the enclosure where it contacts the ground to make an air-tight seal. Prior to the purging and sampling activities, a helium tracer gas was released via a small diameter tube placed through a port on the exterior side of the enclosure. The sub-slab vapor Teflon tubing that extends up through the air tight seal to the exterior side of the enclosure was connected to a helium detector to determine the presence of helium gas and to purge the sampling point. All sampling points pass the trace gas test (i.e. less than 10% tracer gas in test).

Corresponding ambient air samples were collected within 5 ft and at a height of 3 to 5 ft above the ground surface at each sample locations

SVI samples were collected utilizing individually laboratory certified-clean 1-liter Summa® canisters equipped with laboratory calibrated flow controllers. The samples were collected over an approximate 8 hour time period. Immediately after opening each Summa® canister, the initial vacuum (inches of mercury) and time was noted and recorded on the laboratory chain-of-custody. After approximately 8 hours, final vacuum readings (inches of mercury) were noted and the Summa® canisters were closed.

Each sample (including QA/QC samples) was submitted to Centek Laboratories, LLC in Syracuse, New York for analysis of targeted VOCs by USEPA Method TO-15.

## **5.3 Quality Assurance/Quality Control (QA/QC)**

QA/QC procedures were implemented during the investigation in order to ensure accuracy, precision, and completeness of the chemical data collected during the investigation. Samples for QA/QC were taken during the portions of the environmental work in order to evaluate the validity of sampling and analytical methods employed throughout the work.

There were two types of field duplicates taken during sampling activities: “blind” duplicates and matrix spike/matrix spike duplicate (MS/MSD) samples. Blind duplicates were labeled in such a manner that the laboratory would not know which samples they were duplicating, nor that they were actually duplicates in some cases. This process allowed LaBella to verify laboratory reproducibility of analytical data. MS/MSD samples were also submitted and were identified as such on the chain-of-custody so the laboratory could perform internal quality checks on instrument performance.

During sampling activities, QA/QC samples were submitted to Centek Laboratories, LLC in Syracuse, New York for analysis. The samples were analyzed for targeted VOCs using USEPA Method TO-15 per the NYSDEC approved work plan. Targeted VOCs and their associated laboratory detection limits included:

**TABLE 5.3 – Targeted VOCs**

Targeted VOCs	Laboratory Detection Limit ( $\mu\text{g}/\text{m}^3$ )
TCE	0.25
cis-1,2-dichloroethene (CIS-1,2-DCE)	1.0
trans-1,2-dichloroethene	1.0
vinyl chloride	0.25
1,1-Dichloroethene (1,1-DCE)	1.0
chloroethane	1.0

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

The samples were collected prior to the NYSDOH update to the Soil Vapor/Indoor Air Decisions Matrices in May 2017.

Full data evaluation packages were submitted to LaBella for the sampling data in accordance with NYSDEC Electronic Data Deliverables (EDD) requirements. Data usability summary reports were conducted by Dataval, Inc. based on the following parameters:

- Sample Data Reporting Format
- Preservation and Holding Time Compliance
- GC/MS Instrument Performance Check
- Initial Calibration Verification (ICV)
- Continuing Calibration Verification (CCV)
- Blank Sample Analysis
- System Monitoring/Surrogate Compound Recoveries
- Laboratory Control Sample (LCS) Recoveries
- Matrix Spike/Matrix Spike Duplicate (MS/MSD) Recoveries
- Internal Standards
- Target Compound Identification
- Compound Quantitation

- Data Qualifiers
- Summary

Copies of the data usability summary reports (DUSRs) are included in Appendix 2 and the laboratory analytical reports that were provided by Centek are provided in Appendix 3 of this report. The laboratory reports are provided on a CD due to the size of the documents. Qualifiers are included in those reports as well as in the summary tables within this report.

The validation of the analytical results for samples collected from the project site indicate that the samples were processed in general compliance with applicable protocols, and the majority of results are usable as reported, or usable with minor edits or qualification as estimated or edits to non-detection. None of the results were rejected.

## 6.0 ANALYTICAL RESULTS

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This section presents and describes the analytical results for the soil gas and SVI samples collected during the investigation. Analytical result tables for the samples collected are presented in the following subsections. The analytical results were compared to applicable NYSDOH guidance values and/or standards.

Samples submitted during the SVI investigation for laboratory analysis are summarized in the following tables:

- **Soil Gas Samples** - Table 6.1 (attached)
- **Sub-Slab Vapor Samples** – Table 6.2A (attached)

### 6.1 Soil Gas Results

VOCs were detected in two (2) of the four (4) soil gas samples above the laboratory method detection (MDL) limit as summarized in Table 6.1. Sample SV-1 and SV-2 were collected adjacent to the 705 Saint Paul Street building and in the downgradient location of the northern chlorinated VOC groundwater plume at the 690 Saint Paul Street site. Soil gas samples SV-1 and SV-2 did not detect any VOCs above the laboratory method detection limit, however, one VOC (chloroethane) was identified as estimated below the detection limit. The samples with VOCs detected above the laboratory method detection limit were soil gas samples SV-3 and SV-4 that are located adjacent to the southern end of the 691 Saint Paul Street site building and are in the downgradient area of the southern chlorinated VOC plume from the 690 Saint Paul Street site. VOCs were not detected above the laboratory MDL in the outdoor ambient air/background air sample.

SCGs are not applicable to soil gas results, however, the results are used to assess the potential for a SVI concern in the buildings adjacent to these sample locations. Based on the soil gas samples results, the targeted VOCs present in soil gas samples SV-3 and SV-4 (adjacent to the 691 Saint Paul Street site building) warranted further evaluation (i.e. SVI testing of the 691 Saint Paul Street building).

## 6.2 Soil Vapor Intrusion Sample Results

The SVI (i.e. sub-slab and indoor air samples) sample results were compared to the guidance values listed in the NYSDOH SVI Guidance document. For compounds without specific indoor air guidance values, typical background levels are used for comparison purposes. The NYSDOH SVI Guidance Appendix C, includes a USEPA 2001 Building Assessment and Survey Evaluation (BASE) Database which provides background levels in indoor air of commercial and public buildings for comparison purposes. For the purposes of this evaluation, the 90<sup>th</sup> percentile values were utilized. It should be noted that this database is referenced to provide a relative benchmark for comparison to the indoor air sampling data, but does not represent regulatory standards or compliance values. The SVI sample results are summarized on Table 6.2A (attached).

Target VOCs were detected above the laboratory MDL in each of the sub-slab vapor samples and were also detected above the laboratory MDL in five (5) of the six (6) corresponding ambient air samples. The outdoor air/background ambient air sample did not detect targeted VOCs above the laboratory MDL.

A comparison of the SVI results to the NYSDOH Ambient Air Guidelines, NYSDOH Decision Matrices, and the USEPA BASE Database 90<sup>th</sup> Percentile values are summarized in Table 6.2B.

TABLE 6.2B - Summary of SVI Sample Results

Sample Location	Result above USEPA Database - 90th Percentile	Indoor Ambient Air Result above NYSDOH Ambient Air Guideline	NYSDOH Decision Matrix	
			Result Above Minimum NYSDOH Decision Matrix	NYSDOH Recommendation
Upper Basement				
691-NE	None	None	None	None
691-B15	TCE	TCE	TCE	Mitigate
691-B19	TCE	TCE	TCE	Mitigate
Lower Basement				
691-SB5B	TCE, CIS-1,2-DCE	TCE	TCE, CIS-1,2-DCE	Mitigate
691-SB5A	TCE, CIS-1,2-DCE	TCE	TCE, CIS-1,2-DCE	Mitigate
691-5A	TCE, CIS-1,2-DCE	None	CIS-1,2-DCE	Mitigate

Note: CIS-1,2-DCE denotes cis-1,2-Dichloroethene, TCE denotes Trichloroethene

As indicated in Table 6.2B, TCE was detected above the NYSDOH Ambient Air Guideline in samples 691-B15 and 691-B-19 from the upper basement and samples 691-SB5B and 691-SB5A in the lower basement. Comparison of the results to the NYSDOH Decision Matrices recommends mitigation at five (5) of the six (6) sample locations. The location that did not identify a recommended action (per NYSDOH Guidance) was location 691-NE which was in the upper basement and directly adjacent to the 705 Saint Paul Street building.



## 7.0 SUMMARY OF FINDINGS AND CONCLUSIONS

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Based on the results of the SVI investigation, the following findings were made:

- The exterior soil gas sampling identified chlorinated VOCs above the laboratory detection limits in the southern two soil gas sample locations (i.e. SV-3 and SV-4). The northern soil gas locations (i.e. SV-1 and SV-2) did not identify VOCs above the laboratory detection limit.
- TCE was detected in four (4) of the six (6) ambient indoor air samples above the NYSDEC Ambient Air Guideline Value of  $2.0 \mu\text{g}/\text{m}^3$  at two (2) sample locations in the upper basement (i.e. 691-B15 and 691-B19) and two (2) sample locations in the lower basement (i.e. 691-SB5B and 691-SB5A). TCE was not detected in the indoor air in the sample from the northeast corner of the 691 Saint Paul Street building. (i.e. 691-NE) which is directly adjacent to the 705 Saint Paul Street building.
- Comparison of the SVI samples results to the NYSDOH Decision Matrices (updated May 2017), indicates NYSDOH recommends mitigation at five (5) of the six (6) sample locations based on the concentrations of TCE and/or CIS-1,2-DCE detected in the SVI samples. The sample location where mitigation is not recommended by the NYSDOH guidance is the location in the northeast corner adjacent to the 705 Saint Paul Street building.

Based on the above, the following conclusions are made:

- A completed SVI pathway appears present at the 691 Saint Paul Street building.
- A completed SVI pathway does not appear present for the 705 Saint Paul Street building.

## 8.0 RECOMMENDATIONS

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Based on the above conclusions, it is recommended that a sub-slab depressurization system (SSDS) be assessed to mitigate SVI vapors from entering the 691 Saint Paul Street building for both the upper and lower basement areas.

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# TABLES

## REFERENCE PAGE FOR SAMPLE RESULTS FOR TABLES 6.1 AND 6.2A

### Soil Gas and Vapor Intrusion Sampling Results

NYSDEC Site #C828159A

691 and 705 Saint Paul Street, Rochester, New York

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Concentrations displayed in micrograms per cubic meter ( $\mu\text{g}/\text{m}^3$ )

Samples analyzed by USEPA Method TO-15

NL denotes Not Listed

(1) *New York State Department of Health (NYSDOH), Guidance for Evaluating Soil Vapor Intrusion in the State of New York*, October 2006 and updates for tetrachloroethene (September 2013), TCE (August 2015), and the soil vapor/indoor air decision matrices (May 2017). These updates are included on NYSDOH website: [https://www.health.ny.gov/environmental/indoors/vapor\\_intrusion/update.htm](https://www.health.ny.gov/environmental/indoors/vapor_intrusion/update.htm). [Note: This Guidance uses a combination of indoor air and sub-slab soil vapor when comparing to the matrices. In addition, for compounds not listed in the matrices an overall site approach is employed which utilizes the USEPA BASE Database (see 2. below) as typical background for commercial buildings and also uses the outdoor air sample, refer to Guidance document for details.]

(2) USEPA 2001 Building Assessment and Survey Evaluation (BASE) Database (90th Percentile). As recommended in Section 3.2.4 of the NYSDOH Guidance (Refer to Footnote "1") this database is referenced for the indoor air sampling results. This database is also referenced to provide initial benchmarks for comparison to the air sampling data and does not represent regulatory standards or compliance values.

(3) No value was listed in NYSDOH Table C2 - USEPA Base Database. A value from Table C3 NYSDOH 1997: Control home database (90th Percentile) was used.

U indicates the value was detected below the reported laboratory method detection limit

UJ indicates the value was detected below the estimated reported laboratory method detection limit

J indicates an estimated value

**BOLD** type denotes that the compound was detected at a concentration that was found to exceed the NYSDOH Sub-Slab Vapor Concentration Decision Matrix (minimum action level).

UNDERLINED type denotes that the compound was detected at a concentration that was found to exceed the NYSDOH Indoor Air Concentration (minimum action level).

**RED** values are above Air Guideline Derived by NYSDOH in Table 3.1 of NYSDOH Guidance titled "Evaluating Soil Vapor Intrusion in the State of NY", October 2006 (and subsequent updates).

**YELLOW HIGHLIGHT** indicates the compound was detected at a concentration that was found to exceed NYSDOH Guidance 90th Percentile Database Value

\* NYSDOH Air Guideline Value included in Table 3.1 of NYSDOH Guidance Document

**TABLE 6.1****SOIL GAS SAMPLE RESULTS**

NYSDEC Site #C828159A

691 and 705 Saint Paul Street, Rochester, New York

Sample ID	Units	SV-1	SV-2	SV-3	SV-4	DUPLICATE (SV-4)	AMBIENT AIR
Sample Type		Soil Gas	Soil Gas	Soil Gas	Soil Gas	Soil Gas	Outdoor Air
Sample Date		3/3/2017	3/3/2017	3/3/2017	3/3/2017	3/3/2017	3/3/2017
1,1-Dichloroethene	ug/m <sup>3</sup>	0.59 U	0.59 U	0.59 U	0.59 U	0.59 U	0.59 U
Chloroethane	ug/m <sup>3</sup>	0.34 J	0.40 U	0.29 J	0.40 U	0.40 U	0.40 U
cis-1,2-Dichloroethene	ug/m <sup>3</sup>	0.59 U	0.59 U	0.71	4.1	3.9	0.59 U
trans-1,2-Dichloroethene	ug/m <sup>3</sup>	0.59 U	0.59 U	0.59 U	0.59 U	0.59 U	0.59 U
Trichloroethene (TCE)	ug/m <sup>3</sup>	0.81 U	0.81 U	40 J	48	47	0.21 U
Vinyl chloride	ug/m <sup>3</sup>	0.38 U	0.38 U	0.38 U	0.38 U	0.38 U	0.10 U
Total Detected VOCs	ug/m <sup>3</sup>	0	0	41.0	52.1	50.9	0

**TABLE 6.2A****SAMPLE LOCATION: 691-SB5B****Soil Vapor Intrusion Sampling Results**

NYSDEC Site #C828159A

691 and 705 Saint Paul Street, Rochester, New York

Sample ID	Units	691-SB5B-SVI	691-SB5B-IAQ	691-OUTDOOR-04012017	NYSDOH Sub-Slab Vapor Concentration Decision Matrix (minimum action level) <sup>(1)</sup>	NYSDOH Indoor Air Concentration (minimum action level) <sup>(1)</sup> / NYSDOH Air Guideline Value	NYSDOH Guidance Table C2 USEPA BASE Database 90th Percentile <sup>(2)</sup>
Sample Type		Sub-Slab	Indoor Air	Outdoor Air			
Sample Date		4/1/2017	4/1/2017	4/1/2017			
1,1-Dichloroethene	ug/m <sup>3</sup>	0.59 UJ	0.59 UJ	0.59 U	<6	<0.2	<1.4
Chloroethane	ug/m <sup>3</sup>	0.40 UJ	0.40 UJ	0.40 U	NL	NL	<1.2
cis-1,2-Dichloroethene	ug/m <sup>3</sup>	<b>23</b> J	<b>12</b> J	0.59 U	<6	<0.2	<1.8
trans-1,2-Dichloroethene	ug/m <sup>3</sup>	0.71 J	0.59 UJ	0.59 U	NL	NL	<10 <sup>(3)</sup>
Trichloroethene (TCE)	ug/m <sup>3</sup>	<b>58</b> J	<b>6.8</b> J	0.21 U	<6	<0.2 / 2 *	1.3
Vinyl chloride	ug/m <sup>3</sup>	0.28 J	0.15 J	0.10 U	<6	<0.2	<1.8

**TABLE 6.2A****SAMPLE LOCATION: 691-SB5A****Soil Vapor Intrusion Sampling Results**

NYSDEC Site #C828159A

691 and 705 Saint Paul Street, Rochester, New York

Sample ID	Units	691-SB5A-SVI	691-SB5A-IAQ	691-OUTDOOR-04012017	NYSDOH Sub-Slab Vapor Concentration Decision Matrix (minimum action level) <sup>(1)</sup>	NYSDOH Indoor Air Concentration (minimum action level) <sup>(1)</sup> / NYSDOH Air Guideline Value	NYSDOH Guidance Table C2 USEPA BASE Database 90th Percentile <sup>(2)</sup>
Sample Type		Sub-Slab	Indoor Air	Outdoor Air			
Sample Date		4/1/2017	4/1/2017	4/1/2017			
1,1-Dichloroethene	ug/m <sup>3</sup>	0.59 U	0.59 U	0.59 U	<6	<0.2	<1.4
Chloroethane	ug/m <sup>3</sup>	0.40 U	0.40 U	0.40 U	NL	NL	<1.2
cis-1,2-Dichloroethene	ug/m <sup>3</sup>	170	13	0.59 U	<6	<0.2	<1.8
trans-1,2-Dichloroethene	ug/m <sup>3</sup>	1.6	0.59 U	0.59 U	NL	NL	<10 <sup>(3)</sup>
Trichloroethene (TCE)	ug/m <sup>3</sup>	20 J	6.5	0.21 U	<6	<0.2 / 2 *	1.3
Vinyl chloride	ug/m <sup>3</sup>	0.49	0.10 U	0.10 U	<6	<0.2	<1.8

**TABLE 6.2A****SAMPLE LOCATION: 691-SB1****Soil Vapor Intrusion Sampling Results**

NYSDEC Site #C828159A

691 and 705 Saint Paul Street, Rochester, New York

Sample ID	Units	691-SB1-SVI	691-SB1-IAQ	691-OUTDOOR-04012017	NYSDOH Sub-Slab Vapor Concentration Decision Matrix (minimum action level) <sup>(1)</sup>	NYSDOH Indoor Air Concentration (minimum action level) <sup>(1)</sup> / NYSDOH Air Guideline Value	NYSDOH Guidance Table C2 USEPA BASE Database 90th Percentile <sup>(2)</sup>
Sample Type		Sub-Slab	Indoor Air	Outdoor Air			
Sample Date		4/1/2017	4/1/2017	4/1/2017			
1,1-Dichloroethene	ug/m <sup>3</sup>	0.59 U	0.59 UJ	0.59 U	<6	<0.2	<1.4
Chloroethane	ug/m <sup>3</sup>	0.40 U	0.40 UJ	0.40 U	NL	NL	<1.2
cis-1,2-Dichloroethene	ug/m <sup>3</sup>	52	2.6 J	0.59 U	<6	<0.2	<1.8
trans-1,2-Dichloroethene	ug/m <sup>3</sup>	0.59 U	0.59 UJ	0.59 U	NL	NL	<10 <sup>(3)</sup>
Trichloroethene (TCE)	ug/m <sup>3</sup>	4.0	1.3 J	0.21 U	<6	<0.2 / 2 *	1.3
Vinyl chloride	ug/m <sup>3</sup>	0.38	0.10 UJ	0.10 U	<6	<0.2	<1.8

**TABLE 6.2A****SAMPLE LOCATION: 691-NE****Soil Vapor Intrusion Sampling Results**

NYSDEC Site #C828159A

691 and 705 Saint Paul Street, Rochester, New York

Sample ID	Units	691-NE-SVI	691-NE-IAQ	691-OUTDOOR-04012017	NYSDOH Sub-Slab Vapor Concentration Decision Matrix (minimum action level) <sup>(1)</sup>	NYSDOH Indoor Air Concentration (minimum action level) <sup>(1)</sup> / NYSDOH Air Guideline Value	NYSDOH Guidance Table C2 USEPA BASE Database 90th Percentile <sup>(2)</sup>
Sample Type		Sub-Slab	Indoor Air	Outdoor Air			
Sample Date		4/1/2017	4/1/2017	4/1/2017			
1,1-Dichloroethene	ug/m <sup>3</sup>	0.59 U	0.59 U	0.59 U	<6	<0.2	<1.4
Chloroethane	ug/m <sup>3</sup>	0.40 U	0.40 U	0.40 U	NL	NL	<1.2
cis-1,2-Dichloroethene	ug/m <sup>3</sup>	0.59 U	0.59 U	0.59 U	<6	<0.2	<1.8
trans-1,2-Dichloroethene	ug/m <sup>3</sup>	0.59 U	0.59 U	0.59 U	NL	NL	<10 <sup>(3)</sup>
Trichloroethene (TCE)	ug/m <sup>3</sup>	3.9	0.21 U	0.21 U	<6	<0.2 / 2 *	1.3
Vinyl chloride	ug/m <sup>3</sup>	0.10 U	0.10 U	0.10 U	<6	<0.2	<1.8



**TABLE 6.2A****SAMPLE LOCATION: 691-B15****Soil Vapor Intrusion Sampling Results**

NYSDEC Site #C828159A

691 and 705 Saint Paul Street, Rochester, New York

Sample ID	Units	691-B15-SVI	691-B15-IAQ	691-OUTDOOR-04012017	NYSDOH Sub-Slab Vapor Concentration Decision Matrix (minimum action level) <sup>(1)</sup>	NYSDOH Indoor Air Concentration (minimum action level) <sup>(1)</sup> / NYSDOH Air Guideline Value	NYSDOH Guidance Table C2 USEPA BASE Database 90th Percentile <sup>(2)</sup>
Sample Type		Sub-Slab	Indoor Air	Outdoor Air			
Sample Date		4/1/2017	4/1/2017	4/1/2017			
1,1-Dichloroethene	ug/m <sup>3</sup>	0.59 U	0.59 U	0.59 U	<6	<0.2	<1.4
Chloroethane	ug/m <sup>3</sup>	0.37 J	0.40 U	0.40 U	NL	NL	<1.2
cis-1,2-Dichloroethene	ug/m <sup>3</sup>	0.59 U	1.6	0.59 U	<6	<0.2	<1.8
trans-1,2-Dichloroethene	ug/m <sup>3</sup>	0.59 U	0.59 U	0.59 U	NL	NL	<10 <sup>(3)</sup>
Trichloroethene (TCE)	ug/m <sup>3</sup>	220	2.2	0.21 U	<6	<0.2 / 2 *	1.3
Vinyl chloride	ug/m <sup>3</sup>	0.28	0.10 U	0.10 U	<6	<0.2	<1.8

**TABLE 6.2A****SAMPLE LOCATION: 691-B19****Soil Vapor Intrusion Sampling Results**

NYSDEC Site #C828159A

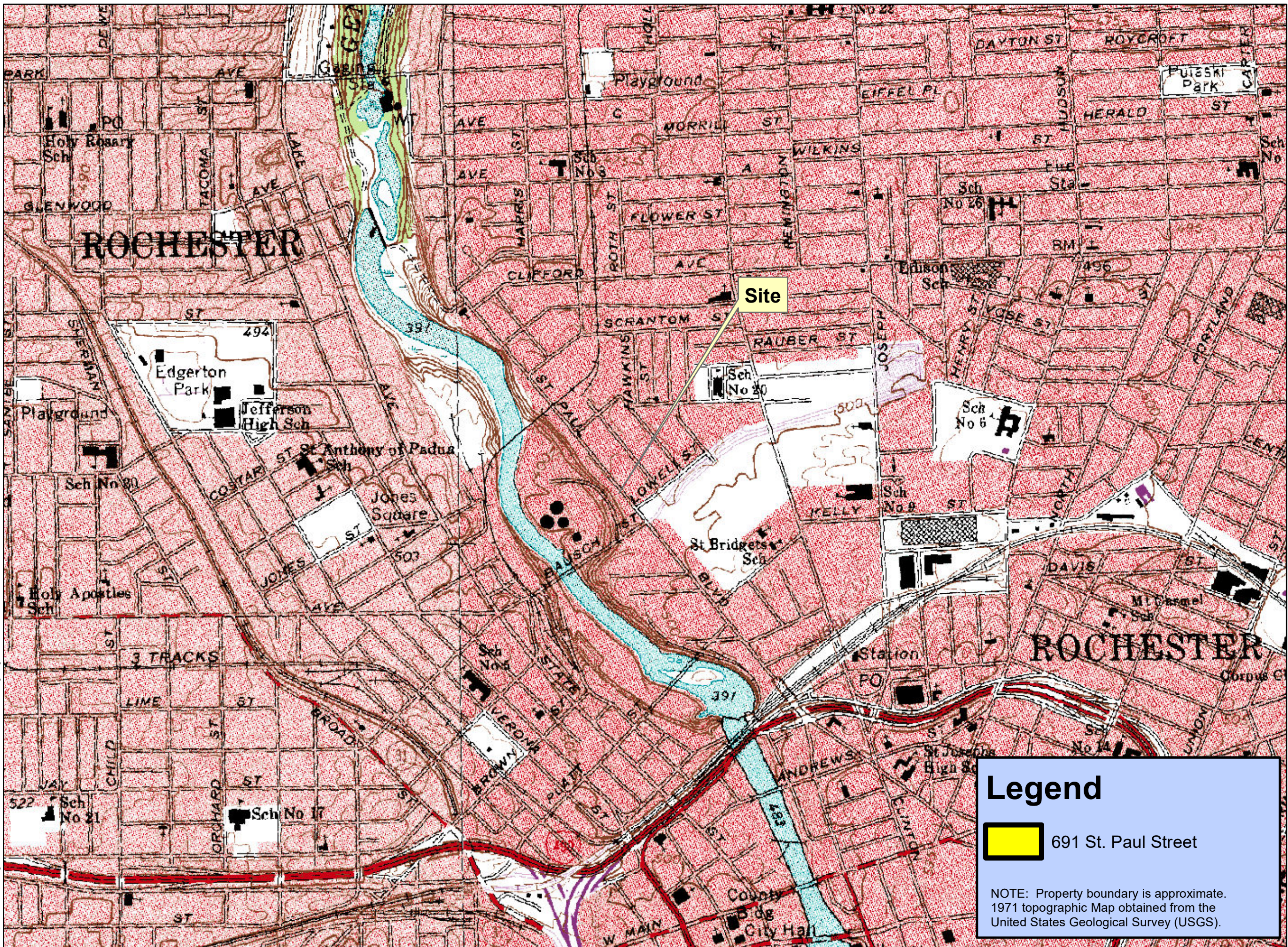
691 and 705 Saint Paul Street, Rochester, New York

Sample ID	Units	691-B19-SVI	691-B19-IAQ-1	691-B19-IAQ-2	691-OUTDOOR-04012017	NYSDOH Sub-Slab Vapor Concentration Decision Matrix (minimum action level) <sup>(1)</sup>	NYSDOH Indoor Air Concentration (minimum action level) <sup>(1)</sup> / NYSDOH Air Guideline Value	NYSDOH Guidance Table C2 USEPA BASE Database - 90th Percentile <sup>(2)</sup>
Sample Type		Sub-Slab	Indoor Air	Indoor Air	Outdoor Air			
Sample Date		4/1/2017	4/1/2017	4/1/2017	4/1/2017			
1,1-Dichloroethene	ug/m <sup>3</sup>	0.59 UJ	0.59 UJ	0.59 U	0.59 U	<6	<0.2	<1.4
Chloroethane	ug/m <sup>3</sup>	0.40 UJ	0.40 UJ	0.40 U	0.40 U	NL	NL	<1.2
cis-1,2-Dichloroethene	ug/m <sup>3</sup>	0.59 UJ	1.3 J	1.1	0.59 U	<6	<0.2	<1.8
trans-1,2-Dichloroethene	ug/m <sup>3</sup>	0.59 UJ	0.59 UJ	0.59 U	0.59 U	NL	NL	<10 <sup>(3)</sup>
Trichloroethene (TCE)	ug/m <sup>3</sup>	26 J	2.1 J	2.1	0.21 U	<6	<0.2 / 2 *	1.3
Vinyl chloride	ug/m <sup>3</sup>	0.20 J	0.10 UJ	0.10 U	0.10 U	<6	<0.2	<1.8

## FIGURES



I:\Barozzi Woodworking, Inc\213211 - 1445 Jefferson Rd, Phase II, ESADrawings\Figure2.mxd



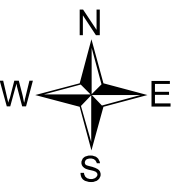
**SITE LOCATION MAP**

**SOIL VAPOR INTRUSION  
INVESTIGATION REPORT**

**690 ST. PAUL ST OFF-SITE  
NYSDEC SITE #C828159A**

**691 AND 705 SAINT PAUL ST  
ROCHESTER, NEW YORK**

**CLIENT:  
BAUSCH & LOMB**



250 0 1,000  
1 inch = 1,000 feet

DATE: 8/17/2017

INTENDED TO PRINT 11"X17"

[ 2170820 ]  
[ **FIGURE 1** ]



Path: I:\Bausch & Lomb\2170820 - 691 St Paul St SVIA BCP Site\Drawings\SVI Report\Figure 2 Site Layout Map.mxd



**Legend**

- NYSDEC BCP Site #C828159A Boundary
- BCP Site #C828159 BOUNDARY
- Monroe County Tax Parcel

NOTE:

(1) Property boundary is approximate. 2012 Aerial photograph obtained from GIS Clearinghouse. Tax parcel data obtained from Monroe County Real Property.

**TITLE**  
**SITE LAYOUT MAP**

**PROJECT**  
**SOIL VAPOR INTRUSION**  
**INVESTIGATION REPORT**

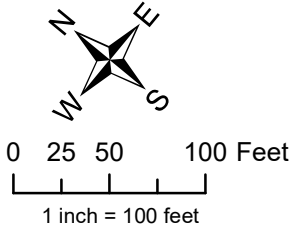
**690 ST. PAUL ST OFF-SITE**  
**NYSDEC SITE #C828159A**

**691 AND 705 SAINT PAUL ST**  
**ROCHESTER, NEW YORK**

**CLIENT:**  
**BAUSCH AND LOMB**



It is a violation of New York Education Law Article 145 Sec.7209, for any person, unless acting under the direction of a licensed architect, professional engineer, or land surveyor, to alter an item in any way. If an item bearing the seal of an architect, engineer, or land surveyor is altered; the altering architect, engineer, or land surveyor shall affix to the item their seal and notation "altered by" followed by their signature and date of such alteration, and a specific description of the alteration.



DATE: 8/17/2017

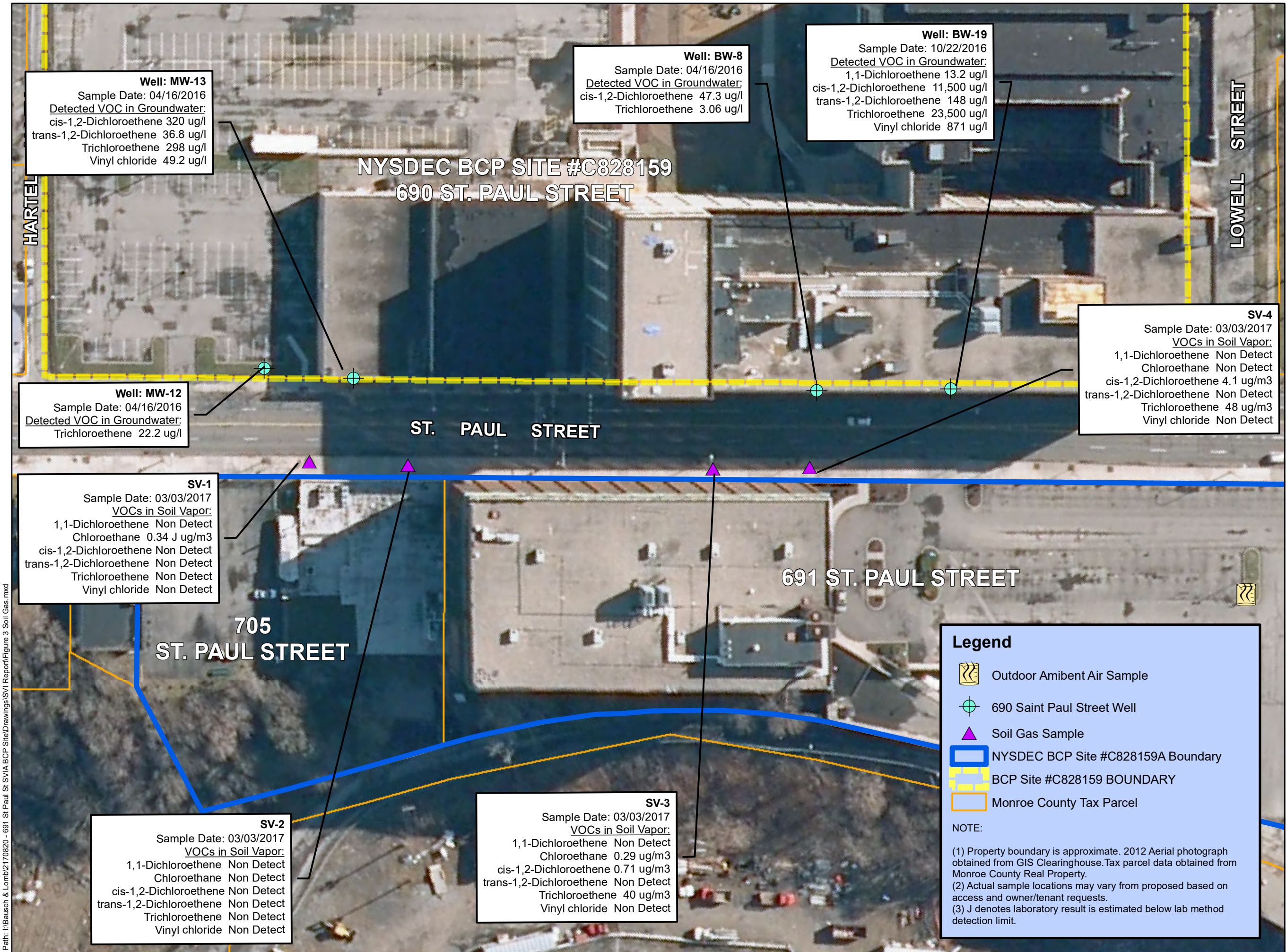
INTENDED TO PRINT 11"X17"

**2170820**

**FIGURE 2**



Path: I:\Bausch & Lomb\2170820 - 691 St Paul St SVIA BCP Site\Drawings\SVI Report\Figure 3 Soil Gas.mxd



**TITLE**  
**SOIL GAS SAMPLE RESULTS AND  
SUMMARY OF 690 SAINT PAUL ST  
GROUNDWATER IMPACTS**

**PROJECT**  
**SOIL VAPOR INTRUSION  
INVESTIGATION REPORT**

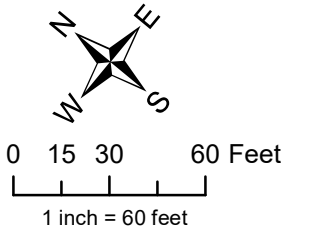
**690 ST. PAUL ST OFF-SITE  
NYSDEC SITE #C828159A**

**691 AND 705 SAINT PAUL ST  
ROCHESTER, NEW YORK**

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DATE: 8/17/2017

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**2170820**

**FIGURE 3**



ST. PAUL STREET



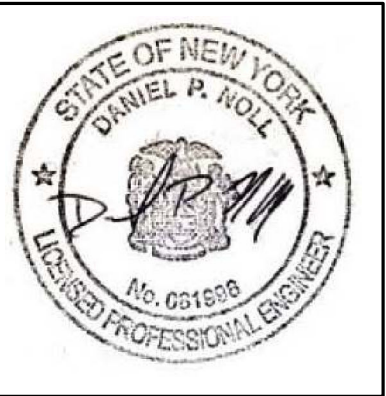
**TITLE**  
**SOIL VAPOR INTRUSION**  
**SAMPLING - UPPER BASEMENT**

**PROJECT**  
**SOIL VAPOR INTRUSION**  
**INVESTIGATION REPORT**

**690 ST. PAUL ST OFF-SITE**  
**NYSDEC SITE #C828159A**

**691 AND 705 SAINT PAUL ST**  
**ROCHESTER, NEW YORK**

**CLIENT**  
**BAUSCH AND LOMB**



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0 12.5 25 Feet  
1 inch = 25 feet

DATE: 10/05/2017

INTENDED TO PRINT 11"x17"

2170820

FIGURE 4A

**INDOOR AIR SAMPLE:**  
**691-NE-IAQ [UG/M3]**  
1,1-DCE ND  
CHLOROMETHANE ND  
CIS-1,2-DCE ND  
TRANS-1,2-DCE ND  
TCE ND  
VINYL CHLORIDE ND  
**SUB-SLAB VAPOR SAMPLE:**  
**691-NE-SVI [UG/M3]**  
1,1-DCE ND  
CHLOROETHANE ND  
CIS-1,2-DCE ND  
TRANS-1,2-DCE ND  
TCE 3.9  
VINYL CHLORIDE ND

Approximate area of  
crawl space beneath  
floor (access through  
lower basement)

**Legend**

- Outdoor Ambient Air Sample
- Sub Slab Vapor Sample - Lower Basement
- Sub Slab Vapor Sample - Upper Basement
- NYSDEC BCP Site #C828159A Boundary
- Approximate Area of Lower Basement

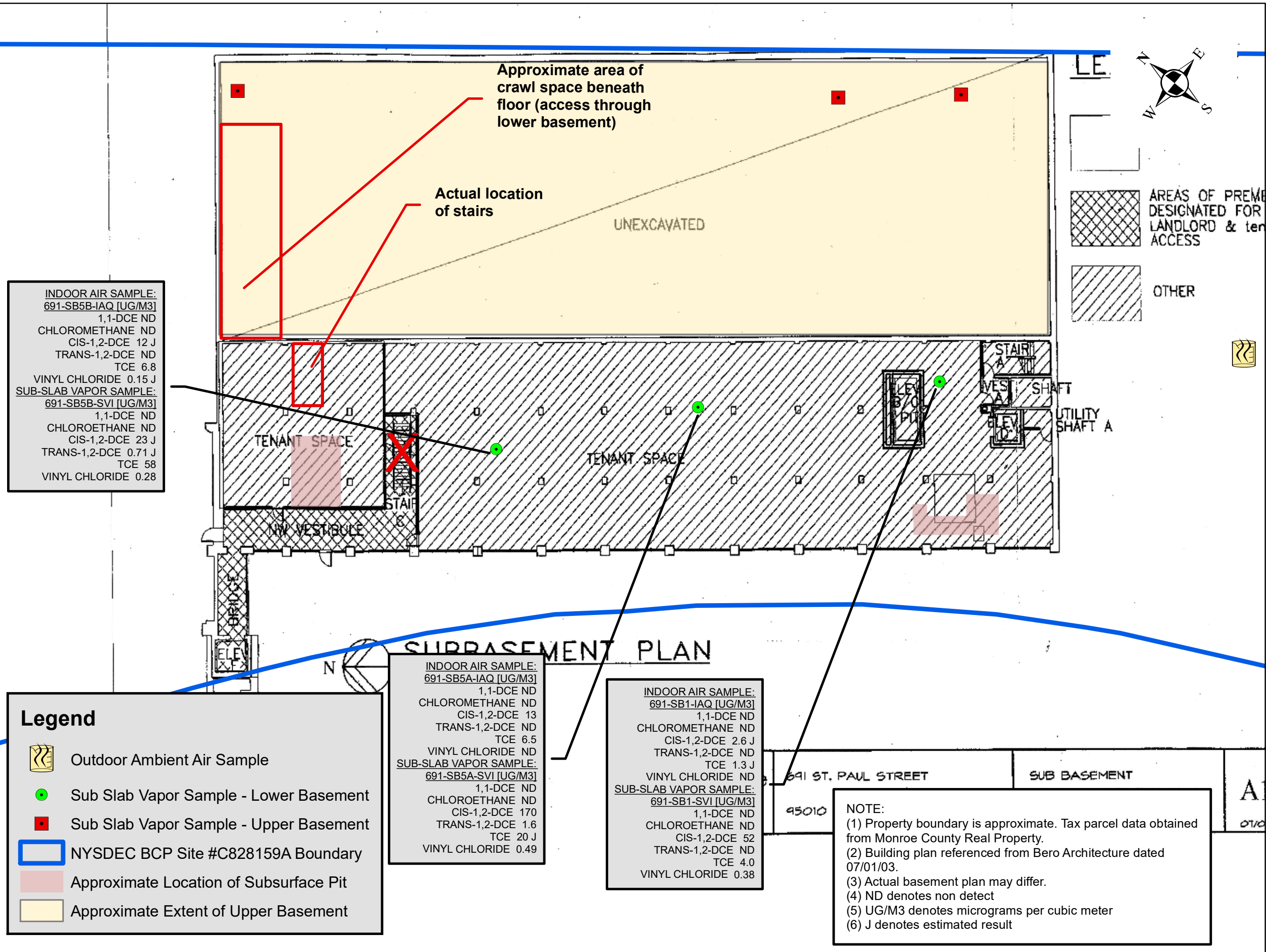
**INDOOR AIR SAMPLE:**  
**691-B15-IAQ [UG/M3]**  
1,1-DCE ND  
CHLOROMETHANE ND  
CIS-1,2-DCE 1.6  
TRANS-1,2-DCE ND  
TCE 2.2  
VINYL CHLORIDE ND  
**SUB-SLAB VAPOR SAMPLE:**  
**691-B15-SVI [UG/M3]**  
1,1-DCE ND  
CHLOROETHANE 0.37 J  
CIS-1,2-DCE ND  
TRANS-1,2-DCE ND  
TCE 220  
VINYL CHLORIDE 0.28


**INDOOR AIR SAMPLE:**  
**691-B19-IAQ-1 [UG/M3]**  
1,1-DCE ND  
CHLOROMETHANE ND  
CIS-1,2-DCE 1.3 J  
TRANS-1,2-DCE ND  
TCE 2.1 J  
VINYL CHLORIDE ND  
**SUB-SLAB VAPOR SAMPLE:**  
**691-B19-SVI [UG/M3]**  
1,1-DCE ND  
CHLOROETHANE ND  
CIS-1,2-DCE ND  
TRANS-1,2-DCE ND  
TCE 26 J  
VINYL CHLORIDE 0.20 J

**NOTE:**  
(1) Property boundary is approximate. Tax parcel data obtained from Monroe County Real Property.  
(2) Building plan referenced from Bero Architecture dated 07/01/03.  
(3) Actual basement plan may differ.  
(4) ND denotes Non Detect  
(5) UG/M3 denotes micrograms per cubic meter  
(6) J denotes estimated result



Path: I:\Bausch & Lomb\2170820 - 691 St Paul St VIA BCP Site\Drawings\SVI Report\Figure 4B Sample Map Lower Basement.mxd






**TITLE**  
SOIL VAPOR INTRUSION  
SAMPLING AT  
UPPER BASEMENT

**PROJECT**  
SOIL VAPOR INTRUSION  
INVESTIGATION REPORT

**690 ST. PAUL ST OFF-SITE  
NYSDEC SITE #C828159A**

**691 AND 705 SAINT PAUL ST  
ROCHESTER, NEW YORK**

**CLIENT**  
BAUSCH AND LOMB



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0 12.5 25 Feet

1 inch = 25 feet

DATE: 10/05/2017

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2170820

FIGURE 4B

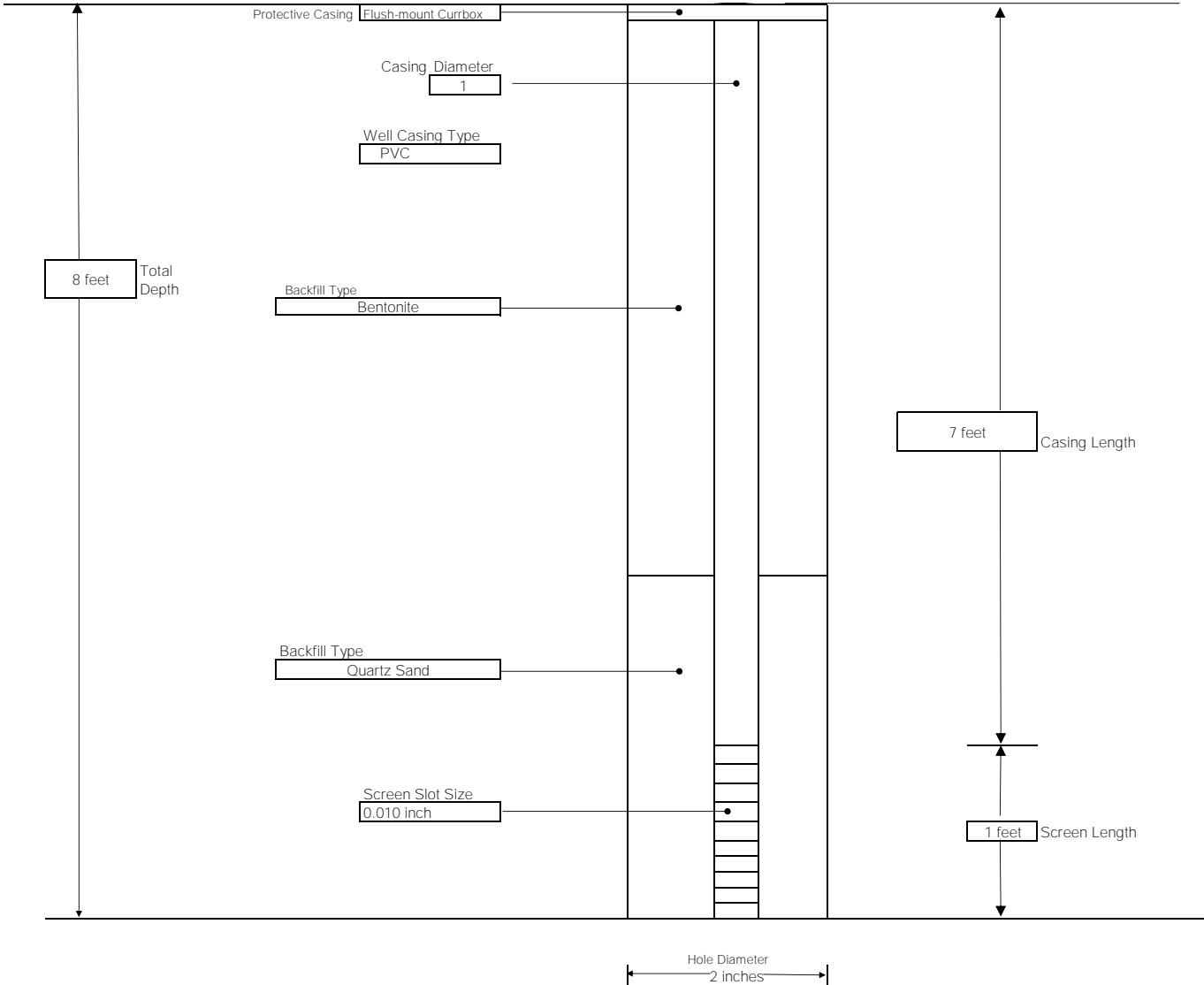


# APPENDIX 1

Field Logs and NYSDOH Building Inventory Form

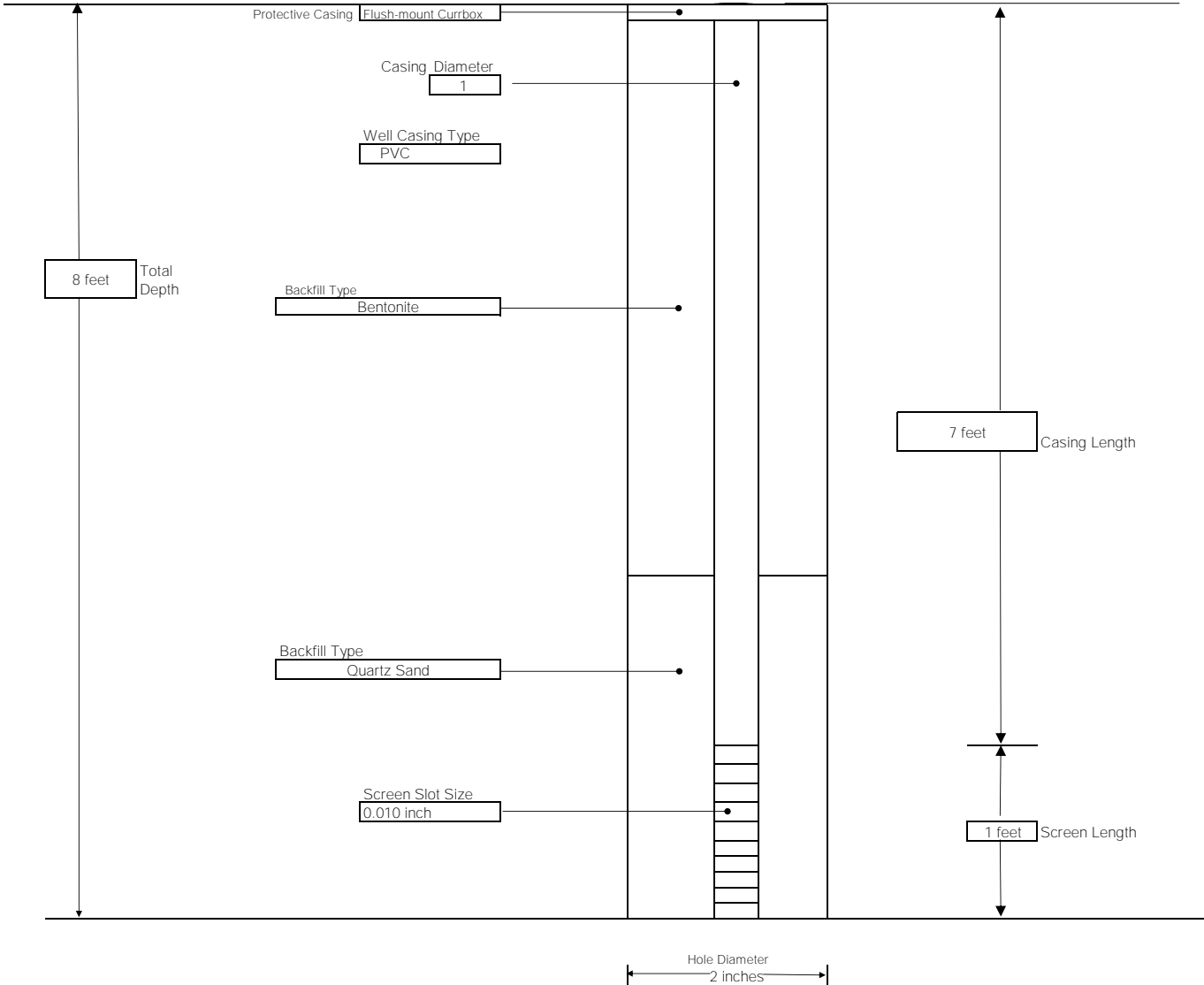
# SOIL GAS POINT INSTALLATION LOG

<b>LABELLA</b> Associates, P.C. 300 STATE STREET, ROCHESTER, NEW YORK ENVIRONMENTAL ENGINEERING CONSULTANTS	<b>PROJECT</b> 691 and 705 Saint Paul Street Rochester, New York Phase II Environmental Site Assessment		<b>Soil Gas Sampling Point: SV-1</b> <b>SHEET</b> 1 OF 1 <b>Project #</b> 2170436 <b>Client</b> Bausch & Lomb
	<b>CONTRACTOR:</b> LaBella Env. LLC <b>DRILLER:</b> M Pepe and M. Winderl <b>LABELLA REPRESENTATIVE:</b> Alex Brett		
<b>BORING LOCATION:</b> SV-1 <b>GROUND SURFACE ELEVATION:</b> NA <b>START DATE:</b> 03/01/2017		<b>DATUM:</b> NA <b>END DATE:</b> 03/01/2017	
<b>TYPE OF DRILL RIG:</b> Geoprobe 6620 DT <b>AUGER SIZE AND TYPE:</b> NA <b>OVERBURDEN SAMPLING METHOD:</b> NA <b>ROCK DRILLING METHOD:</b> NA			



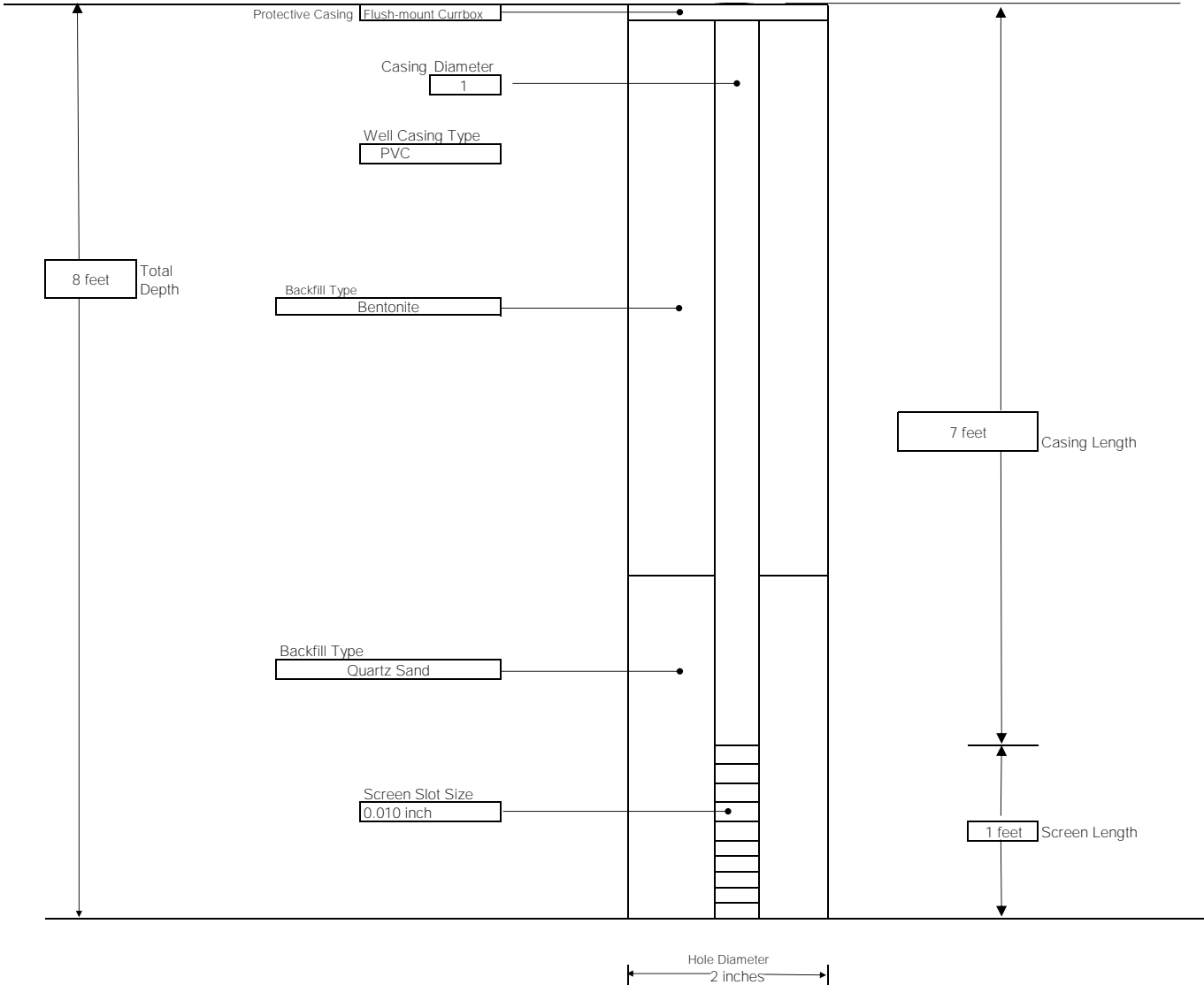
NOTE: NOT TO SCALE

<b>LABELLA</b> Associates, P.C. 300 STATE STREET, ROCHESTER, NEW YORK ENVIRONMENTAL ENGINEERING CONSULTANTS	<b>PROJECT</b> 691 and 705 Saint Paul Street Rochester, New York Phase II Environmental Site Assessment		<b>Soil Gas Sampling Point: SV-2</b> <b>SHEET</b> 1 OF 1 <b>Project #</b> 2170436 <b>Client</b> Bausch & Lomb
	CONTRACTOR: LaBella Env. LLC DRILLER: M Pepe and M. Winderl LABELLA REPRESENTATIVE: Alex Brett		
BORING LOCATION: SV-1 GROUND SURFACE ELEVATION: NA START DATE: 03/01/2017		DATUM: NA END DATE: 03/01/2017	
TYPE OF DRILL RIG: Geoprobe 6620 DT AUGER SIZE AND TYPE: NA OVERBURDEN SAMPLING METHOD: NA ROCK DRILLING METHOD: NA			



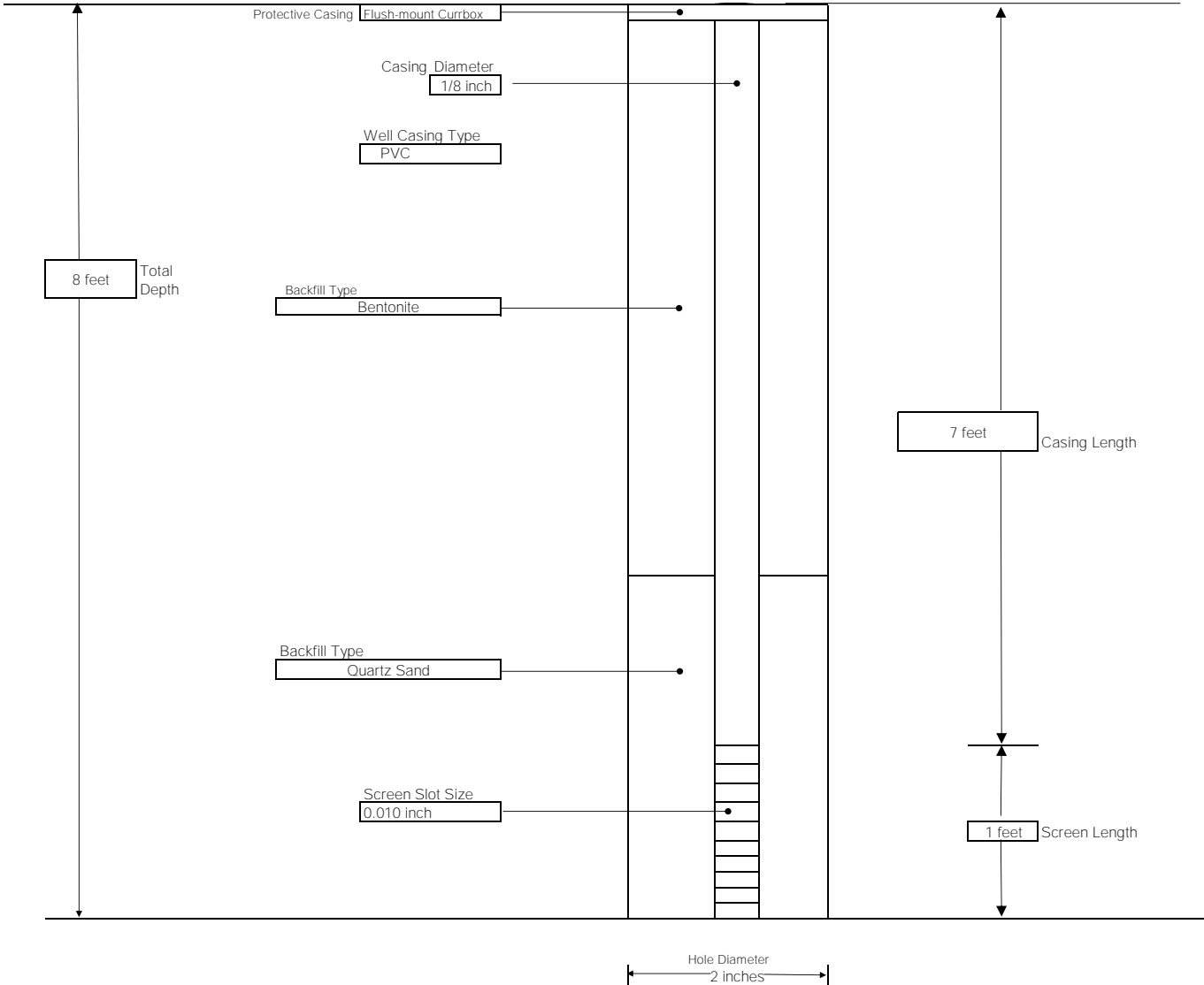
NOTE: NOT TO SCALE

<b>LABELLA</b> Associates, P.C. 300 STATE STREET, ROCHESTER, NEW YORK ENVIRONMENTAL ENGINEERING CONSULTANTS	<b>PROJECT</b> 691 and 705 Saint Paul Street Rochester, New York Phase II Environmental Site Assessment		<b>Soil Gas Sampling Point: SV-3</b> <b>SHEET</b> 1 OF 1 <b>Project #</b> 2170436 <b>Client</b> Bausch & Lomb
	CONTRACTOR: LaBella Env. LLC DRILLER: M Pepe and M. Winderl LABELLA REPRESENTATIVE: Alex Brett		
BORING LOCATION: SV-1 GROUND SURFACE ELEVATION: NA START DATE: 03/01/2017		DATUM: NA END DATE: 03/01/2017	
TYPE OF DRILL RIG: Geoprobe 6620 DT AUGER SIZE AND TYPE: NA OVERBURDEN SAMPLING METHOD: NA ROCK DRILLING METHOD: NA			



NOTE: NOT TO SCALE

<b>LABELLA</b> Associates, P.C. 300 STATE STREET, ROCHESTER, NEW YORK ENVIRONMENTAL ENGINEERING CONSULTANTS	<b>PROJECT</b> 691 and 705 Saint Paul Street Rochester, New York Phase II Environmental Site Assessment		<b>Soil Gas Sampling Point: SV-4</b> <b>SHEET</b> 1 OF 1 <b>Project #</b> 2170436 <b>Client</b> Bausch & Lomb
	<b>CONTRACTOR:</b> LaBella Env. LLC <b>DRILLER:</b> M Pepe and M. Winderl <b>LABELLA REPRESENTATIVE:</b> Alex Brett		
<b>BORING LOCATION:</b> SV-1 <b>GROUND SURFACE ELEVATION:</b> NA <b>START DATE:</b> 03/01/2017		<b>DATUM:</b> NA <b>END DATE:</b> 03/01/2017	
<b>TYPE OF DRILL RIG:</b> Geoprobe 6620 DT <b>AUGER SIZE AND TYPE:</b> NA <b>OVERBURDEN SAMPLING METHOD:</b> NA <b>ROCK DRILLING METHOD:</b> NA			



NOTE: NOT TO SCALE

# SOIL GAS SAMPLE LOG



## AIR SAMPLING FIELD REPORT

AIR  
SAMPLING  
POINT

SV-1

Project: Soil Gas Sampling  
Site Location: 691 and 705 Saint Paul Street  
Client: Bausch & Lomb  
LaBella Project No.: 2170436  
LaBella Representative: AGB  
Weather: Cloudy, moderate winds out of NW, temp in 20's

### General Information

Sample Canister Location: Sidewalk near northeast corner of 705 St Paul Street building along St Paul Street

Sample Source: ☐ Indoor Air ☐ Sub-Slab ☐ Interior Ambient Air ☐ Exterior Ambient Air  
☒ Exterior Soil Gas ☐ Other

Shipping Date: \_\_\_\_\_ Laboratory: Centek

Canister Type: ☒ 1.0 L Summa Canister ☐ 6.0 L Summa Canister Other (specify): \_\_\_\_\_

Canister Serial No.: 542 Flow Controller Serial No.: 256

### Purge Information

Leak Detection Test Date: 03/03/2017 Leak Detection: ☒ Pass ☐ Fail

Purging Method: Photoionization Detector

PID Reading Start: 0 ppm PID Reading End: 0.5 ppm

Volume of Gas Extracted: ~10 Liters

### Sampling Information

Sample Date: 03/03/2017 Sampler: AGB

Sample Depth: 7-8 ft

	Start	Stop
Canister Pressure Gauge Reading:	-30	-5
Sample Time:	1145	1925

### Comments:

Time and pressure Readings.	1145	- 30
	1310	- 25
Pressures shown are	1410	- 22
negative to show vacuum	1510	- 19
	1610	- 15
	1710	- 12
	1810	- 8.5
	1925	- 5





## AIR SAMPLING FIELD REPORT

AIR  
SAMPLING  
POINT

SV-2

Project: Soil Gas Sampling  
Site Location: 691 and 705 Saint Paul Street  
Client: Bausch & Lomb  
LaBella Project No.: 2170436  
LaBella Representative: AGB  
Weather: Cloudy, moderate winds out of NW, temp in 20's

### General Information

Sample Canister Location: Sidewalk near southeast corner of 705 St Paul Street building along St Paul Street

Sample Source: ☐ Indoor Air ☐ Sub-Slab ☐ Interior Ambient Air ☐ Exterior Ambient Air  
☒ Exterior Soil Gas ☐ Other

Shipping Date: \_\_\_\_\_ Laboratory: Centek

Canister Type: ☒ 1.0 L Summa Canister ☐ 6.0 L Summa Canister Other (specify): \_\_\_\_\_

Canister Serial No.: 237 Flow Controller Serial No.: 402

### Purge Information

Leak Detection Test Date: 03/03/2017 Leak Detection: ☒ Pass ☐ Fail

Purging Method: Photoionization Detector

PID Reading Start: 0 ppm PID Reading End: 0.7 ppm

Volume of Gas Extracted: ~10 Liters

### Sampling Information

Sample Date: 03/03/2017 Sampler: AGB

Sample Depth: 7-8 ft

	Start	Stop
Canister Pressure Gauge Reading:	-30	-5
Sample Time:	1155	1930

### Comments:

Time and pressure Readings.	1155	- 30
	1310	- 24
Pressures shown are	1410	- 22
negative to show vacuum	1510	- 19
	1610	- 15
	1710	- 12
	1810	- 9
	1930	- 5

Time and pressure Readings.	1205	- 30+	"Plus" symbol after pressure to show vacuum is greater than -30 upon start
	1310	- 27	
Pressures shown are	1410	- 24	
negative to show vacuum	1510	- 21	
	1610	- 18	
MS/MSD Collected	1710	- 14	
with Sample	1810	- 11	
	2000	- 5	

2040 - 5

# SOIL VAPOR INTRUSION SAMPLE LOG

Comments:

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Comments:

CONCRETE FLOOR 5-6" THICK'

Comments:

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Comments:

CONCRETE FLOOR 5-6" THICK'





## AIR SAMPLING FIELD REPORT

AIR  
SAMPLING  
POINT

691-SB5A-IAQ

Project: NYSDEC BCP SITE #C828159A  
Site Location: 691 SAINT PAUL ST, ROCHESTER, NY  
Client: BAUSCH & LOMB

LaBella Project No.: 2170820  
LaBella Representative: K. MILLER  
Weather: OVERCAST, RAIN, 40s F

### General Information

Sample Canister Location: LOWER BASEMENT, ROOM SB5, WOODWORKING SHOP

Sample Source: X Indoor Air        Sub-Slab        Interior Ambient Air        Exterior Ambient Air  
       Exterior Soil Gas        Other

Shipping Date: 4/1/2017 Laboratory: CEN TEK LABORATORIES

Canister Type: X 1.0 L Summa Canister        6.0 L Summa Canister Other (specify):       

Canister Serial No.: 367 Flow Controller Serial No.: 251

### NOT APPLICABLE AMBIENT AIR SAMPLE

### Purge Information

Leak Detection Test Date:        Leak Detection:        Pass        Fail

Purging Method:       

PID Reading Start:        ppm PID Reading End:        ppm

Volume of Gas Extracted:       

### Sampling Information

Sample Date: 4/1/2017 Sampler: K. MILLER

Sample Depth: NOT APPLICABLE

Start Stop

Canister Pressure Gauge Reading [IN. HG]: 30 2

Sample Time: 1112 1740

Comments:





## AIR SAMPLING FIELD REPORT

AIR  
SAMPLING  
POINT

691-SB1-IAQ

Project: NYSDEC BCP SITE #C828159A  
Site Location: 691 SAINT PAUL ST, ROCHESTER, NY  
Client: BAUSCH & LOMB

LaBella Project No.: 2170820  
LaBella Representative: K. MILLER  
Weather: OVERCAST, RAIN, 40s F

### General Information

Sample Canister Location: LOWER BASEMENT, ROOM SB1

Sample Source: X Indoor Air        Sub-Slab        Interior Ambient Air        Exterior Ambient Air  
       Exterior Soil Gas        Other

Shipping Date: 4/1/2017 Laboratory: CEN TEK LABORATORIES

Canister Type: X 1.0 L Summa Canister        6.0 L Summa Canister Other (specify):       

Canister Serial No.: 170 Flow Controller Serial No.: 387

### NOT APPLICABLE AMBIENT AIR SAMPLE

### Purge Information

Leak Detection Test Date:        Leak Detection:        Pass        Fail

Purging Method:       

PID Reading Start:        ppm PID Reading End:        ppm

Volume of Gas Extracted:       

### Sampling Information

Sample Date: 4/1/2017 Sampler: K. MILLER

Sample Depth: NOT APPLICABLE

Start Stop

Canister Pressure Gauge Reading [IN. HG]: NA 5

Sample Time: 1101 1911

Comments:



## AIR SAMPLING FIELD REPORT

AIR  
SAMPLING  
POINT

691-NE-SVI

Project: NYSDEC BCP SITE #C828159A  
Site Location: 691 SAINT PAUL ST, ROCHESTER, NY  
Client: BAUSCH & LOMB

LaBella Project No.: 2170820  
LaBella Representative: K. MILLER  
Weather: OVERCAST, RAIN, 40s F

### General Information

Sample Canister Location: STAIRWELL, NORTHEAST CORNER OF BLDG [ROOM B12]

Sample Source:            Indoor Air   X   Sub-Slab            Interior Ambient Air            Exterior Ambient Air  
           Exterior Soil Gas            Other

Shipping Date: 4/1/2017 Laboratory: CEN TEK LABORATORIES

Canister Type:   X   1.0 L Summa Canister            6.0 L Summa Canister Other (specify):           

Canister Serial No.: 1186 Flow Controller Serial No.: 145

### Purge Information

Leak Detection Test Date: 4/1/2017 Leak Detection:   X   Pass            Fail

Purging Method: MGD 2002 HELIUM LEAK DETECTOR

Volume of Gas Extracted: 3 PROBE VOLUMES (~0.02 L)

### Sampling Information

Sample Date: 4/1/2017 Sampler: K. MILLER

Sample Depth: NOT APPLICABLE

           Start            Stop

Canister Pressure Gauge Reading [IN. HG]: 30 3

Sample Time: 1045 1855

Comments: CONCRETE FLOOR 4-5" THICK'



## AIR SAMPLING FIELD REPORT

AIR  
SAMPLING  
POINT

691-NE-IAQ

Project: NYSDEC BCP SITE #C828159A  
Site Location: 691 SAINT PAUL ST, ROCHESTER, NY  
Client: BAUSCH & LOMB

LaBella Project No.: 2170820  
LaBella Representative: K. MILLER  
Weather: OVERCAST, RAIN, 40s F

### General Information

Sample Canister Location: STAIRWELL, NORTHEAST CORNER OF BLDG [ROOM B12]

Sample Source: ☒ Indoor Air ☐ Sub-Slab ☐ Interior Ambient Air ☐ Exterior Ambient Air  
☐ Exterior Soil Gas ☐ Other

Shipping Date: 4/1/2017 Laboratory: CEN TEK LABORATORIES

Canister Type: ☒ 1.0 L Summa Canister ☐ 6.0 L Summa Canister Other (specify):

Canister Serial No.: 94 Flow Controller Serial No.: 379

### NOT APPLICABLE AMBIENT AIR SAMPLE

### Purge Information

Leak Detection Test Date:  Leak Detection: ☐ Pass ☐ Fail

Purging Method:

PID Reading Start:  ppm PID Reading End:  ppm

Volume of Gas Extracted:

### Sampling Information

Sample Date: 4/1/2017 Sampler: K. MILLER

Sample Depth: NOT APPLICABLE

Start

Stop

Canister Pressure Gauge Reading [IN. HG]: 28

1

Sample Time: 1042

1705

Comments:

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## AIR SAMPLING FIELD REPORT

AIR  
SAMPLING  
POINT

691-B15-IAQ

Project: NYSDEC BCP SITE #C828159A  
Site Location: 691 SAINT PAUL ST, ROCHESTER, NY  
Client: BAUSCH & LOMB

LaBella Project No.: 2170820  
LaBella Representative: K. MILLER  
Weather: OVERCAST, RAIN, 40s F

### General Information

Sample Canister Location: ROOM B15 CUSTODIAN STORAGE OFFICE

Sample Source: X Indoor Air        Sub-Slab        Interior Ambient Air        Exterior Ambient Air  
       Exterior Soil Gas        Other

Shipping Date: 4/1/2017 Laboratory: CEN TEK LABORATORIES

Canister Type: X 1.0 L Summa Canister        6.0 L Summa Canister Other (specify):       

Canister Serial No.: 101 Flow Controller Serial No.: 299

### NOT APPLICABLE AMBIENT AIR SAMPLE

### Purge Information

Leak Detection Test Date:        Leak Detection:        Pass        Fail

Purging Method:       

PID Reading Start:        ppm PID Reading End:        ppm

Volume of Gas Extracted:       

### Sampling Information

Sample Date: 4/1/2017 Sampler: K. MILLER

Sample Depth: NOT APPLICABLE

Start Stop

Canister Pressure Gauge Reading [IN. HG]: 27.5 3

Sample Time: 1024 1825

Comments: DUPLICATE SAMPLE ALSO COLLECTED (CAN #87) USING T CONNECTION







## AIR SAMPLING FIELD REPORT

AIR  
SAMPLING  
POINT

691-B19-IAQ-2

Project: NYSDEC BCP SITE #C828159A  
Site Location: 691 SAINT PAUL ST, ROCHESTER, NY  
Client: BAUSCH & LOMB

LaBella Project No.: 2170820  
LaBella Representative: K. MILLER  
Weather: OVERCAST, RAIN, 40s F

### General Information

Sample Canister Location: UPPER BASEMENT UTILITY ROOM B19

Sample Source: X Indoor Air        Sub-Slab        Interior Ambient Air        Exterior Ambient Air  
       Exterior Soil Gas        Other

Shipping Date: 4/1/2017 Laboratory: CEN TEK LABORATORIES

Canister Type: X 1.0 L Summa Canister        6.0 L Summa Canister Other (specify):       

Canister Serial No.: 1188 Flow Controller Serial No.: 306

### NOT APPLICABLE AMBIENT AIR SAMPLE

### Purge Information

Leak Detection Test Date:        Leak Detection:        Pass        Fail

Purging Method:       

PID Reading Start:        ppm PID Reading End:        ppm

Volume of Gas Extracted:       

### Sampling Information

Sample Date: 4/1/2017 Sampler: K. MILLER

Sample Depth: NOT APPLICABLE

Start Stop

Canister Pressure Gauge Reading [IN. HG]: 30 9

Sample Time: 1135 1940

Comments:

Comments:

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FLOW REGULATOR APPEARS TO HAVE MALFUNCTIONED AFTER CONNECTION TO CANISETER, PRESSURE DROP FROM 30" TO 7" AT 1130. THEN HOLDS STEADY AT 4" FROM 1140. DEPLOYED 1 L REPLACEMENT SAMPLE 691-B19-IAQ-2

# NYSDOH BUILDING INVENTORY FORM

**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 Kyle R. Miller Date/Time Prepared 4/1/17 p.m.  
Preparer's Affiliation LaBella Assoc. Phone No. \_\_\_\_\_  
Purpose of Investigation DPC SVI Assessment #828159A

**1. OCCUPANT:**Interviewed: Y ☒ N

Last Name: \_\_\_\_\_ First Name: \_\_\_\_\_

Address: \_\_\_\_\_

County: \_\_\_\_\_

Home Phone: \_\_\_\_\_ Office Phone: \_\_\_\_\_

Number of Occupants/persons at this location \_\_\_\_\_ Age of Occupants \_\_\_\_\_

**2. OWNER OR LANDLORD:** (Check if same as occupant \_\_\_\_)Interviewed: Y ☒ N

Last Name: \_\_\_\_\_ First Name: \_\_\_\_\_

Address: \_\_\_\_\_

County: \_\_\_\_\_

Home Phone: \_\_\_\_\_ Office Phone: \_\_\_\_\_

**3. BUILDING CHARACTERISTICS****Type of Building:** (Circle appropriate response)Residential  
IndustrialSchool  
ChurchCommercial/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: _____

N/A

If multiple units, how many? \_\_\_\_\_

If the property is commercial, type?

Business Type(s) Office space, woodworking shop

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

Other characteristics:

Number of floors multiple Building age ?

Is the building insulated? Y / N unknown How air tight? Tight / Average / Not Tight unknown

#### 4. AIRFLOW

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

Airflow between floors

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Airflow near source

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Outdoor air infiltration

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Infiltration into air ducts

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### 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 \_\_\_\_\_
- 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 unknown
- k. Water in sump? Y / N / not applicable

Basement/Lowest level depth below grade: \_\_\_\_\_ (feet)

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

### 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:

(Nat. gas? and) electric

Boiler/furnace located in: Basement Outdoors Main Floor Other \_\_\_\_\_

Air conditioning: Central Air Window units Open Windows None

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.

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## 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	<hr/>
1 <sup>st</sup> Floor	<hr/>
2 <sup>nd</sup> Floor	<hr/>
3 <sup>rd</sup> Floor	<hr/>
4 <sup>th</sup> Floor	<hr/>

## 8. FACTORS THAT MAY INFLUENCE INDOOR AIR QUALITY

- |  |                                    |
|--|------------------------------------|
| a. Is there an attached garage?  | Y <u>N</u>                         |
| 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<br>Please specify <hr/> |
| d. Has the building ever had a fire?   | Y / N When? <hr/>                  |
| e. Is a kerosene or unvented gas space heater present?   | Y / N Where? <hr/>                 |
| f. Is there a workshop or hobby/craft area?  | Y / N Where & Type? <hr/>          |
| g. Is there smoking in the building?   | Y / N How frequently? <hr/>        |
| h. Have cleaning products been used recently?  | Y / N When & Type? <hr/>           |
| i. Have cosmetic products been used recently?  | Y / N When & Type? <hr/>           |

- 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? \_\_\_\_\_
- 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: \_\_\_\_\_

Do any of the building occupants use solvents at work?

(e.g., chemical manufacturing or laboratory, auto mechanic or auto body shop, painting, fuel oil delivery, boiler mechanic, pesticide application, cosmetologist)

Y / N woodworking shop

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)

Yes, use dry-cleaning infrequently (monthly or less)

Yes, work at a dry-cleaning service

No

Unknown

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 work plan*

**First Floor:**

**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 work plan*

## 13. PRODUCT INVENTORY FORM

Make &amp; Model of field instrument used: \_\_\_\_\_

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** Y/N
① Sub-basement wood working shop (SB5)		various sizes	good	latex and oil based paints and stains; mineral spirits; paint thinners; spray paints; various wood working putties and supplies and sealants		Y
② custodians storage area and office area (B15)			good	several spray bottles and 1-gal. containers of <del>cleaners</del> cleaning solutions stored on janitors carts		Y

\* 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.

③ Loading dock on SW corner of bldg. connected to upper basement area by man door and o/H door has yellow fire proof cabinet containing gasoline, oil/gas mix and spray lube.



Chemical storage in wood shop



Chemical storage in wood shop



Chemical storage in wood shop



Chemical storage in wood shop



Typical chemicals in upper basement space



Typical chemicals in upper basement space

## APPENDIX 2

### Data Usability Summary Reports



DATA USABILITY SUMMARY REPORT

for

LaBella Associates, P.C.

300 State Street

Rochester, NY 14614

691 ST PAUL SITE

Project 2170820

SDG: C1704014

Sampled 4/4/2017

TO-15 AIR SAMPLES

691-B19-IAQ-01	(C1704014-01)	691-B19-IAQ-02	(C1704014-02)
691-B19-SVI	(C1704014-03)	691-B15-IAQ	(C1704014-04)
691-B15-SVI	(C1704014-05)	691-OUTDOOR-0401217	(C1704014-06)
691-DUPLICATE	(C1704014-07)	691-NE-IAQ	(C1704014-08)
691-NE-SVI	(C1704014-09)	691-SBI-IAQ	(C1704014-10)
691-SBI-SVI	(C1704014-11)	691-SB5A-IAQ	(C1704014-12)
691-SB5B-IAQ	(C1704014-13)	691-SB5A-SVI	(C1704014-14)
691-SB5B-SVI	(C1704014-15)		

## DATA ASSESSMENT

A TO-15 data package containing analytical results for fifteen air samples was received from LaBella Associates, P.C. on 08May17. The ASP deliverables package included formal reports, raw data, the necessary QC, and supporting information. The samples, taken from the 691 St. Paul Site, were identified by Chain of Custody documents and traceable through the work of Centek Laboratories, LLC, the laboratory contracted for analysis. The analyses were performed using US EPA Method TO-15 and addressed measurements of sixty-three volatile organic compounds. Laboratory data was evaluated according to the quality assurance / quality control requirements of the New York State Department of Environmental Conservation's Analytical Services Protocol (ASP), September 1989, Rev. 07/2005. When the required protocol was not followed, the current EPA Region II Functional Guidelines (SOP HW-31, Rev. #4, October 2006, Volatile Organic Analysis of Ambient Air in Canisters by Method TO-15) was used as a technical reference.

The results from 691-B19-IAQ-01, 691-B19SVI, 691-SB5B-IAQ and 691-SB5B-SVI have been qualified as estimations because the samples were not collected properly.

The results reported from 691-SBI-IAQ have been qualified as estimations because the sample collection was not properly documented.

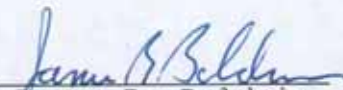
The trichloroethene results from 691-SB5A-SVI, 691-SB5B-SVI and 691-B19-SVI have been qualified as estimations due to poor internal standard performance.

## CORRECTNESS AND USABILITY

Reported data should be considered technically defensible and completely usable in its present form. Results presenting a usable estimation of the conditions at the time of sampling have been flagged "J" or "UJ". Estimated data should be used with caution. A detailed discussion of the review process follows.

Two facts should be considered by all data users. No compound concentration, even if it has passed all QC testing, can be guaranteed to be accurate. Strict QC serves to increase confidence in data, but any value potentially contains error. Secondly, DATAVAL, Inc. guarantees the quality of this data assessment. However, DATAVAL, Inc. does not warrant any interpretation or utilization of this data by a third party.

Reviewer's signature:

  
James B. Baldwin  
DATAVAL, Inc.

Date:

19 MAY 17



## SAMPLE HISTORY

Analyte concentrations can deteriorate with time due to chemical instability, bacterial degradation or volatility. Samples that are not properly preserved or are not analyzed within established holding times may no longer be considered representative. Holding times are calculated from the date of sampling. TO-15 samples must be analyzed within 14 days of collection.

This sample delivery group contained fifteen air samples that were collected in 1-liter SUMMA canisters. Sampling was completed on 01Apr17. The canisters were shipped back to the laboratory, via FedEx, on 04Apr17 and were received the following morning. Although the sample canisters were received intact, custody seals were not present on the packaging. It is noted that 691-B19-IAQ-1 was collected in a 1.4-liter canister to facilitate the preparation of MS/MSD samples.

Although the SUMMA canisters were set in the laboratory to collect 8-hour samples, most of the samples were terminated after 6.5-9.5 hours, based on the canister vacuum readings. The collection of 691-SB5B-IAQ and 691-SB5B-SVI, however, was terminated after just four hours. The results from this pair of samples have been qualified as estimations because the samples were not collected properly.

Numerous samples failed to satisfy the ASP gauge reading requirement of  $5 \pm 1$ "Hg at the completion of sampling. Due to the accuracy of the canister gauges, only results falling outside of the limits of -2"Hg to -8"Hg have been qualified as estimations. 691-B19-IAQ-1, 691-B19-SVI and 691-SB5B-SVI were affected.

SAMPLE	PRIOR TO SHIPMENT ("Hg)	PRIOR TO SAMPLING ("Hg)	POST SAMPLING ("Hg)	LAB ANALYSIS ("Hg)
691-B19-IAQ-1	-30	-30	-4	-4
691-B19-IAQ-2	-30	-30	-9	-9
691-B19-SVI	-30	-30	-9	-9
691-B15-IAQ	-30	-27.5	-3	-3
691-B15-SVI	-30	-30	-2	-3
691-OUTDOOR	-30	-28	-8	-8
691-DUPLICATE	-30	-27.5	-3	-3
691-NE-IAQ	-30	-28	-1	-2
691-NE-SVI	-30	-30	-3	-3
691-SB1-IAQ	-30		-5	-5
691-SB1-SVI	-30	-30	-6	-6
691-SB5A-IAQ	-30	-30	-2	-2
691-SB5B-IAQ	-30	-30	-3	-3
691-SB5A-SVI	-30	-28	-2	-2
691-SB5B-SVI	-30	-29.5	-0.5	-1

An examination of the vacuum readings following sampling and at the time of analysis indicates that the integrity of each sample was maintained during this period.



It is noted that a vacuum reading was not recorded prior to the sampling of 691-SB1-IAQ. The results reported from this sample have been qualified as estimations due to this omission. Although these results are assumed to be usable, based on the laboratory's measurement, they would not withstand a legal challenge.

The analysis of this group of samples was completed between 06Apr17 and 08Apr17, satisfying the ASP holding time limitation.

#### CANISTER CERTIFICATION

The canisters used for this project were pressure tested at 30 psig for 24 hours. Each canister demonstrated a change  $\leq 0.5$  psig over this period.

The canisters for this project were cleaned in four batches. A blank analysis of a clean canister from each of these batches was free of targeted analyte contamination.

#### BLANKS

Blanks are analyzed to evaluate various sources of sample contamination. Trip Blanks monitor sampling activities, sample transport, and storage. Method blanks are analyzed to verify instrument integrity. Samples are considered compromised by conditions causing contamination in any blank.

Two method blanks were analyzed with this group of samples. Both of these blanks demonstrated acceptable chromatography and were free of targeted analyte contamination.

#### MS TUNING

Mass spectrometer tuning and performance criteria are established to ensure sufficient mass resolution and sensitivity to accurately detect and identify targeted analytes. Verification is accomplished using a certified standard.

BFB ion abundance criteria was reported from standards run before the initial instrument calibration and prior to the analysis of program samples on 06Apr17 and 07Apr17. Each of these checks satisfied the ASP acceptance criteria.

#### CALIBRATION

Requirements for instrument calibration are established to ensure that laboratory equipment is capable of producing accurate, quantitative data. Initial calibrations demonstrate a range through which measurements may be made. Continuing calibration check standards verify instrument stability.

The initial instrument calibration was performed on 31Mar17. Standards of 0.04, 0.10, 0.15, 0.30, 0.50, 0.75, 1.0, 1.25, 1.50 and 2.0 ppbV were included. Each targeted analyte produced the required levels of instrument response and demonstrated an acceptable degree of linearity during this calibration.



Continuing calibration check standards were analyzed on 06Apr17 and 07Apr17, prior to the 24-hour periods of instrument operation that included samples from this program. When compared to the initial calibration, each targeted analyte produced the required levels of instrument response and demonstrated an acceptable degree of linearity during these checks.

#### SURROGATES

Each sample, blank and standard is spiked with surrogate compounds prior to analysis. The structures of surrogates are similar to analytes of interest, but they are not normally found in environmental samples. Surrogate recoveries are monitored to evaluate overall laboratory performance and the efficiency of laboratory technique.

Although surrogate summary sheets were properly prepared, an incorrect acceptance criteria was applied. However, when compared to the ASP requirements, acceptable surrogate recoveries were reported for each addition to this group of samples.

#### INTERNAL STANDARDS

Internal standards are added to each sample, blank and standard just prior to injection. Analyte concentrations are calculated relative to the response of a specific internal standard. Internal standard performance criteria ensure that GC/MS sensitivity and response are stable during the analysis of each sample. The area of internal standard peaks may not vary by more than 40%. When compared to the preceding calibration check, retention times may not vary by more than 10 seconds.

The laboratory recorded the response of each internal standard addition to this group of samples and the response obtained from the preceding CCV standard. Although the control limits based on the response of the CCV were not reported, they were calculated by this reviewer. When compared to these limits, unacceptable performance was reported for the 1,4-difluorobenzene and chlorobenzene-d5 additions to 691-SB5A-SVI, 691-SB5B-SVI and the 1:10 dilution of 691-B19-SVI. The trichloroethene (TCE) results from 691-SB5A-SVI, 691-SB5B-SVI and the 1:10 dilution of 691-B19-SVI have been qualified as estimations based on this performance. The remaining analytes reported from these samples were associated with the response of bromochloromethane and remain unqualified.

Internal standard retention times were not addressed. The ASP retention time acceptance criteria was calculated by this reviewer. The retention times produced by each program sample satisfied these requirements.

#### MATRIX SPIKES / MATRIX SPIKE DUPLICATES / MATRIX SPIKED BLANKS

Matrix spiking refers to the addition of known analyte concentrations to a sample, prior to analysis. Analyte recoveries provide an indication of laboratory accuracy. The analysis of a duplicate spiked aliquot provides a measurement of precision.

691-B19-IAQ-1 was selected for matrix spiking. The six targeted analytes were added to two volumes of this sample. The recoveries reported for these additions demonstrated acceptable levels of measurement precision and accuracy.

Two pairs of spiked blanks (LCS/LCSD) were also analyzed with this delivery group. Each of these spiked blanks produced acceptable recoveries of the six targeted analytes.

#### DUPLICATES

Two aliquots of the same sample are processed separately through all aspects of sample preparation and analysis. Results produced by the analysis of this pair of samples are compared as a measurement of precision. Poor precision may be indicative of sample non-homogeneity, method defects, or poor laboratory technique.

The blind duplicate sample that was included in this delivery group was not identified.

#### REPORTED ANALYTES

Formal reports were provided for each sample. The data package also included total ion chromatograms and raw instrument print-outs. Reference mass spectra were provided to confirm the identification of each analyte that was detected in this group of samples.



# SUMMARY OF QUALIFIED DATA

691 ST PAUL SITE

SAMPLED APRIL 2017

	SAMPLING	VACUUM	SAMPLING RECORD	INT STD TCE
691-B19-IAQ-01	(C1704014-01)	ALL J/UJ		
691-B19-IAQ-02	(C1704014-02)			
691-B19-SVI	(C1704014-03)	ALL J/UJ		26J
691-B15-IAQ	(C1704014-04)			
691-B15-SVI	(C1704014-05)			
691-OUTDOOR-0401217	(C1704014-06)			
691-DUPLICATE	(C1704014-07)			
691-NE-IAQ	(C1704014-08)			
691-NE-SVI	(C1704014-09)			
691-SBI-IAQ	(C1704014-10)		ALL J/UJ	
691-SBI-SVI	(C1704014-11)			
691-SB5A-IAQ	(C1704014-12)			
691-SB5B-IAQ	(C1704014-13)	ALL J/UJ		20J
691-SB5A-SVI	(C1704014-14)			57J
691-SB5B-SVI	(C1704014-15)	ALL J/UJ		

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-001A

Client Sample ID: 691-B19-1AQ-1  
Tag Number: 1321.1163  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC						Analyst: RJP
1,1-Dichloroethene	< 0.59 <i>UJ</i>	0.59		ug/m3	1	4/6/2017 9:55:00 PM
Chloroethane	< 0.40 <i>UJ</i>	0.40		ug/m3	1	4/6/2017 9:55:00 PM
cis-1,2-Dichloroethene -	1.3 <i>J</i>	0.59		ug/m3	1	4/6/2017 9:55:00 PM
trans-1,2-Dichloroethene	< 0.59 <i>UJ</i>	0.59		ug/m3	1	4/6/2017 9:55:00 PM
Trichloroethene -	2.1 <i>J</i>	0.21		ug/m3	1	4/6/2017 9:55:00 PM
Vinyl chloride	< 0.10 <i>UJ</i>	0.10		ug/m3	1	4/6/2017 9:55:00 PM

*7/15*

Qualifiers: \*\* Quantitation Limit  
B Analytic detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analytic. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analytic detected below quantitation limit  
ND Not Detected at the Limit of Detection

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# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-002A

Client Sample ID: 691-B19-IAQ-2  
Tag Number: 1188.306  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 12:09:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 12:09:00 AM
cis-1,2-Dichloroethene -	1.1	0.59		ug/m3	1	4/7/2017 12:09:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 12:09:00 AM
Trichloroethene -	2.1	0.21		ug/m3	1	4/7/2017 12:09:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 12:09:00 AM

*Handwritten signature/initials in red ink.*

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-003A

Client Sample ID: 691-B19-SV1  
Tag Number: 362446  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC				TO-15		Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 3:23:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 3:23:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 3:23:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 3:23:00 PM
Trichloroethene	26	2.1		ug/m3	10	4/7/2017 8:07:00 PM
Vinyl chloride	0.20	0.10		ug/m3	1	4/7/2017 3:23:00 PM

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Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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## Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-004A

Client Sample ID: 691-B15-IAQ  
Tag Number: 87.299  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15				Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 12:50:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 12:50:00 AM
cis-1,2-Dichloroethene	1.6	0.59		ug/m3	1	4/7/2017 12:50:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 12:50:00 AM
Trichloroethene	2.2	0.21		ug/m3	1	4/7/2017 12:50:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 12:50:00 AM



Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-005A

Client Sample ID: 691-B15-SV1  
Tag Number: 550.266  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:03:00 PM
Chloroethene -	0.37	0.40	J	ug/m3	1	4/7/2017 4:03:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:03:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:03:00 PM
Trichloroethene -	220	19		ug/m3	90	4/7/2017 9:20:00 PM
Vinyl chloride -	0.28	0.10		ug/m3	1	4/7/2017 4:03:00 PM

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Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-006A

Client Sample ID: 691-Outdoor-04012017  
Tag Number: 240.340  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 1:31:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 1:31:00 AM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 1:31:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 1:31:00 AM
Trichloroethene	< 0.21	0.21		ug/m3	1	4/7/2017 1:31:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 1:31:00 AM

7/14/17

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
IN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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## Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-007A

Client Sample ID: 691-Duplicate  
Tag Number: 101.299  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 2:12:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 2:12:00 AM
cis-1,2-Dichloroethene	1.7	0.59		ug/m3	1	4/7/2017 2:12:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 2:12:00 AM
Trichloroethene	2.2	0.21		ug/m3	1	4/7/2017 2:12:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 2:12:00 AM

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Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte, Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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## Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-008A

Client Sample ID: 691-NE-IAQ  
Tag Number: 94.379  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15				Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 2:52:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 2:52:00 AM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 2:52:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 2:52:00 AM
Trichloroethene	< 0.21	0.21		ug/m3	1	4/7/2017 2:52:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 2:52:00 AM

9/15

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routing analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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## Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-009A

Client Sample ID: 691-NE-SV1  
Tag Number: 1186.145  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15				Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 5:28:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 5:28:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 5:28:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 5:28:00 PM
Trichloroethene	3.9	0.21		ug/m3	1	4/7/2017 5:28:00 PM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 5:28:00 PM



Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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## Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1704014  
 Project: 691 St Paul Street  
 Lab ID: C1704014-010A

Client Sample ID: 691-SB1-1AQ  
 Tag Number: 170.387  
 Collection Date: 4/1/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59 <i>UJ</i>	0.59		ug/m3	1	4/7/2017 3:33:00 AM
Chloroethane	< 0.40 <i>UJ</i>	0.40		ug/m3	1	4/7/2017 3:33:00 AM
cis-1,2-Dichloroethene --	2.6 <i>J</i>	0.59		ug/m3	1	4/7/2017 3:33:00 AM
trans-1,2-Dichloroethene	< 0.59 <i>UJ</i>	0.59		ug/m3	1	4/7/2017 3:33:00 AM
Trichloroethene --	1.3 <i>J</i>	0.21		ug/m3	1	4/7/2017 3:33:00 AM
Vinyl chloride	< 0.10 <i>UJ</i>	0.10		ug/m3	1	4/7/2017 3:33:00 AM

*2/15*

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

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# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-011A

Client Sample ID: 691-SB1-SV1  
Tag Number: 203.372  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 6:09:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 6:09:00 PM
cis-1,2-Dichloroethene -	52	5.9		ug/m3	10	4/7/2017 11:10:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 6:09:00 PM
Trichloroethene -	4.0	0.21		ug/m3	1	4/7/2017 6:09:00 PM
Vinyl chloride -	0.38	0.10		ug/m3	1	4/7/2017 6:09:00 PM

MS

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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## Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-012A

Client Sample ID: 691-SB5A-1AQ  
Tag Number: 367.251  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:13:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 4:13:00 AM
cis-1,2-Dichloroethene	13	5.9		ug/m3	10	4/7/2017 1:50:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:13:00 AM
Trichloroethene	6.5	0.21		ug/m3	1	4/7/2017 4:13:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 4:13:00 AM

7/15

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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## Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1704014  
 Project: 691 St Paul Street  
 Lab ID: C1704014-013A

Client Sample ID: 691-SB5B-1AQ  
 Tag Number: 285.344  
 Collection Date: 4/1/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59 <i>UJ</i>	0.59		ug/m3	1	4/7/2017 4:53:00 AM
Chloroethane	< 0.40 <i>UJ</i>	0.40		ug/m3	1	4/7/2017 4:53:00 AM
cis-1,2-Dichloroethene -	12 <i>J</i>	5.9		ug/m3	10	4/7/2017 2:27:00 PM
trans-1,2-Dichloroethene	< 0.59 <i>UJ</i>	0.59		ug/m3	1	4/7/2017 4:53:00 AM
Trichloroethene -	6.8 <i>J</i>	0.21		ug/m3	1	4/7/2017 4:53:00 AM
Vinyl chloride -	0.15 <i>J</i>	0.10		ug/m3	1	4/7/2017 4:53:00 AM

*JHS*

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

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## Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-014A

Client Sample ID: 691-SB5A-SV1  
Tag Number: 475.1170  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 6:50:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 6:50:00 PM
cis-1,2-Dichloroethene -	170	55		ug/m3	90	4/8/2017 1:00:00 AM
trans-1,2-Dichloroethene -	1.6	0.59		ug/m3	1	4/7/2017 6:50:00 PM
Trichloroethene -	20 J	1.9		ug/m3	9	4/8/2017 12:23:00 AM
Vinyl chloride -	0.49	0.10		ug/m3	1	4/7/2017 6:50:00 PM

115

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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## Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1704014  
 Project: 691 St Paul Street  
 Lab ID: C1704014-015A

Client Sample ID: 691-SB5B-SV1  
 Tag Number: 467.1167  
 Collection Date: 4/1/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59 <i>J</i>	0.59		ug/m3	1	4/7/2017 7:30:00 PM
Chloroethane	< 0.40 <i>J</i>	0.40		ug/m3	1	4/7/2017 7:30:00 PM
cis-1,2-Dichloroethene —	23 <i>J</i>	5.9		ug/m3	10	4/8/2017 1:37:00 AM
trans-1,2-Dichloroethene —	0.71 <i>J</i>	0.59		ug/m3	1	4/7/2017 7:30:00 PM
Trichloroethene —	57 <i>J</i>	2.1		ug/m3	10	4/8/2017 1:37:00 AM
Vinyl chloride —	0.28 <i>J</i>	0.10		ug/m3	1	4/7/2017 7:30:00 PM

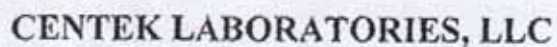
*JAS*

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

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## QC SUMMARY REPORT SURROGATE RECOVERIES

**Matrix: A**

\* Surrogate recovery outside acceptance limits

## GC/MS QA-QC Check Report

Tune File : C:\HPCHEM\1\DATA\AO040602.D

Tune Time : 6 Apr 2017 9:31 am

Daily Calibration File : C:\HPCHEM\1\DATA\AO040602.D

				37398	170379	156465
(BFB)				(IS1)	(IS2)	(IS3)
				26713	121699	111761
				16028	730194	67057
File	Sample	DL	Surrogate Recovery %	Internal Standard Responses		
AO040603.D	ALCS1UG-040617	95		26450	123600	109377
AO040604.D	AMB1UG-040617	90		26038	118374	102379
AO040620.D	ALCS1UGD-040617	98		24809	117535	100863
AO040621.D	C1704014-001A	96		25396	113281	97086
AO040622.D	C1704014-001A MS	97		26903	116940	105999
AO040623.D	C1704014-001A MSD	99		26299	120898	106126
AO040624.D	C1704014-002A	89		24868	113360	102277
AO040625.D	C1704014-004A	92		25596	114646	102450
AO040626.D	C1704014-006A	88		25816	110508	98499
AO040627.D	C1704014-007A	91		24957	110993	99559
AO040628.D	C1704014-008A	93		25088	108028	95812
AO040629.D	C1704014-010A	92		24485	107880	98198
AO040630.D	C1704014-012A	103		25278	111053	100304
AO040631.D	C1704014-013A	102		25079	113761	101873

t - fails 24hr time check \* - fails criteria

Created: Thu May 04 12:00:40 2017 MSD #1/



## GC/MS QA-QC Check Report

Tune File : C:\HPCHEM\1\DATA\AO040702.D

Tune Time : 7 Apr 2017 11:20 am

Daily Calibration File : C:\HPCHEM\1\DATA\AO040702.D

(BFB)				32494	138901	124351
				(IS1)	(IS2)	(IS3)
				23210	99215	88822
				13926	59529	53293
File	Sample	DL	Surrogate Recovery %	Internal	Standard	Responses
AO040703.D	ALCS1UG-040717	97		22420 ✓	98698	86489
AO040704.D	AMB1UG-040717	86		21100	92197	83384
AO040706.D	C1704014-012A 10X	91		20808	86837	77606
AO040707.D	C1704014-013A 10X	92		19443	85901	76902
AO040708.D	C1704014-003A	109		20468	96070	92232
AO040709.D	C1704014-005A	118		24680	109137	104661
AO040710.D	C1704014-009A	113		27394	125972	121869
AO040711.D	C1704014-011A	109		30170	132762	121408
AO040712.D	C1704014-014A	107		31361	139148	131462
AO040713.D	C1704014-015A	116		31915	147799	139968
AO040714.D	C1704014-003A 10x	95		31326	147339	127930
AO040716.D	C1704014-005A 90x	92		26112	119025	102419
AO040719.D	C1704014-011A 10x	93		21591	93800	84591
AO040721.D	C1704014-014A 9x	95		20532	89989	83171
AO040722.D	C1704014-014A 90x	89		20827	88198	80228
AO040723.D	C1704014-015A 10x	94		20780	88679	79938
AO040725.D	ALCS1UGD-040717	97		19313	89890	78098

c - fails 24hr time check \* - fails criteria

Created: Thu May 04 12:02:38 2017 MSD #1/

## CENTEK LABORATORIES, LLC

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.

Work Order: C1704014

Project: 691 St Paul Street

TestCode: 0.25CT-TCE-VC

Sample ID	ALCS1UG-040617	SampType: LCS	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12114					
Client ID:	ZZZZZ	Batch ID: R12114	TestNo: TO-15		Analysis Date: 4/6/2017	SeqNo: 141660					
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	0.9700	0.15	1	0	97.0	70	130				
Chloroethane	1.090	0.15	1	0	109	70	130				
cis-1,2-Dichloroethene	0.9700	0.15	1	0	97.0	70	130				
trans-1,2-Dichloroethene	0.9500	0.15	1	0	95.0	70	130				
Trichloroethene	0.9700	0.040	1	0	97.0	70	130				
Vinyl chloride	0.9700	0.040	1	0	97.0	70	130				

Sample ID	ALCS1UG-040717	SampType: LCS	TestCode: 0.25CT-TCE-	Units: ppbv	Prep Date:	RunNo: 12118					
Client ID: ZZZZZ		Batch ID: R12118	TestNo: TO-15		Analysis Date: 4/7/2017	SeqNo: 141731					
Analyte	PQL	Result	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	0.15	0.9900	1	0	99.0	70	130				
Chloroethane	0.15	1.180	1	0	118	70	130				
cis-1,2-Dichloroethene	0.15	1.000	1	0	100	70	130				
trans-1,2-Dichloroethene	0.15	0.9800	1	0	98.0	70	130				
Trichloroethene	0.040	1.060	1	0	106	70	130				
Vinyl chloride	0.040	1.150	1	0	115	70	130				

Sample ID	ALCS1UGD-040617	SampType: LCSD	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12114					
Client ID: ZZZZZ	Batch ID: R12114	TestNo: TO-15	Analysis Date: 4/6/2017	SeqNo: 141661							
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	0.9900	0.15	1	0	99.0	70	130	0.97	2.04	30	
Chloroethane	1.000	0.15	1	0	100	70	130	1.09	8.61	30	

Qualifiers: J Results reported are not blank corrected  
 S Analyte detected below quantitation limit  
 ND Spike Recovery outside accepted recovery limits  
 E Estimated Value above quantitation range  
 ND Not Detected at the Limit of Detection  
 H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits



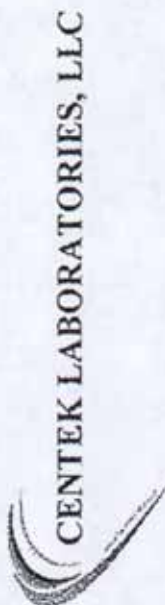
CLIENT: LaBella Associates, P.C.  
 Work Order: C1704014  
 Project: 691 St Paul Street

TestCode: 0.25CT-TCE-VC

Sample ID	ALCS1UGD-040617	Sample Type	LCSD	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12114				
Client ID:	ZZZZZ	Batch ID:	R12114	TestNo: TO-15		Analysis Date: 4/8/2017	SeqNo: 141661				
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPOLimit	Qual
cis-1,2-Dichloroethene	0.9600	0.15	1	0	96.0	70	130	0.97	1.04	30	
trans-1,2-Dichloroethene	0.9800	0.15	1	0	98.0	70	130	0.95	3.11	30	
Trichloroethene	1.000	0.040	1	0	100	70	130	0.97	3.05	30	
Vinyl chloride	1.060	0.040	1	0	106	70	130	0.97	8.87	30	

Sample ID	ALCS1UGD-040717	SampType: LCSD	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12118					
Client ID: ZZZZZ	Batch ID: R12118	TestNo: TO-15	Analysis Date: 4/8/2017	SeqNo: 141732							
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	1.080	0.15	1	0	108	70	130	0.99	8.70	30	
Chloroethane	1.180	0.15	1	0	118	70	130	1.18	0	30	
cis-1,2-Dichloroethene	1.120	0.15	1	0	112	70	130	1	11.3	30	
trans-1,2-Dichloroethene	1.070	0.15	1	0	107	70	130	0.98	8.78	30	
Trichloroethene	1.100	0.040	1	0	110	70	130	1.06	3.70	30	
Vinyl chloride	1.240	0.040	1	0	124	70	130	1.15	7.53	30	

Qualifiers: Results reported are not blank corrected  
 J Analyte detected below quantitation limit  
 S Spike Recovery outside accepted recovery limits  
 E Estimated Value above quantitation range  
 ND Not Detected at the Limit of Detection  
 H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits



Date: 04-May-17

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.

Work Order: C1704014

Project: 691 St Paul Street

TestCode: 0.25CT-TCE-VC

Sample ID	C1704014-001A MS	SampType: MS	TestCode: 0.25CT-TCE-	Units: ppbv	Prep Date:	RunNo: 12114					
Client ID:	691-B19-IAQ-1	Batch ID: R12114	TestNo: TO-15	Analysis Date: 4/6/2017		SeqNo: 141726					
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,1-Dichloroethene

Chloroethane

cis-1,2-Dichloroethene

trans-1,2-Dichloroethene

Trichloroethene

Vinyl chloride

Sample ID	C1704014-001A MS	SampType: MSD	TestCode: 0.25CT-TCE-	Units: ppbv	Prep Date:	RunNo: 12114					
Client ID:	691-B19-IAQ-1	Batch ID: R12114	TestNo: TO-15	Analysis Date: 4/6/2017	SeqNo: 141727						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,1-Dichloroethene

Chloroethane

cis-1,2-Dichloroethene

trans-1,2-Dichloroethene

Trichloroethene

Vinyl chloride

Qualifiers: J Results reported are not blank corrected  
S Analyte detected below quantitation limit  
E Estimated Value above quantitation range  
ND Not Detected at the Limit of Detection  
H Holding times for preparation or analysis exceeded  
R RPD outside accepted recovery limits

Page 1 of 1



Date: 04-May-17



# ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.  
 Work Order: C1704014  
 Project: 691 St Paul Street

TestCode: 0.25CT-TCE-VC

Sample ID	AMB1UG-040617	SampType: MBLK	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12114						
Client ID:	ZZZZZ	Batch ID: R12114	TestNo: TO-15		Analysis Date: 4/8/2017	SeqNo: 141659						
Analyte		Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,1-Dichloroethene	< 0.15	0.15
Chloroethane	< 0.15	0.15
cis-1,2-Dichloroethene	< 0.15	0.15
trans-1,2-Dichloroethene	< 0.15	0.15
Trichloroethene	< 0.040	0.040
Vinyl chloride	< 0.040	0.040

Sample ID	AMB1UG-040717	SampType: MBLK	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12118						
Client ID:	ZZZZZ	Batch ID: R12118	TestNo: TO-15		Analysis Date: 4/7/2017	SeqNo: 141730						
Analyte		Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,1-Dichloroethene	< 0.15	0.15
Chloroethane	< 0.15	0.15
cis-1,2-Dichloroethene	< 0.15	0.15
trans-1,2-Dichloroethene	< 0.15	0.15
Trichloroethene	< 0.040	0.040
Vinyl chloride	< 0.040	0.040

Qualifiers: J Results reported are not blank corrected  
 S Analyte detected below quantitation limit  
 E Estimated Value above quantitation range  
 ND Not Detected at the Limit of Detection  
 H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits

DATA USABILITY SUMMARY REPORT

for

LaBella Associates, P.C.

300 State Street

Rochester, NY 14614

691 ST PAUL SITE

Project 2170436

SDG: C1703015

Sampled 3/3/2017

TO-15 AIR SAMPLES

SV-1	(C1703015-01)
SV-2	(C1703015-02)
SV-3	(C1703015-03)
SV-2	(C1703015-04)
DUPLICATE	(C1703015-05)
AMBIENT AIR	(C1703015-06)



## DATA ASSESSMENT

A TO-15 data package containing analytical results for six air samples was received from LaBella Associates, P.C. on 22May17. The ASP deliverables package included formal reports, raw data, the necessary QC, and supporting information. The samples, taken from the 691 St. Paul Site, were identified by Chain of Custody documents and traceable through the work of Centek Laboratories, LLC, the laboratory contracted for analysis. The analyses were performed using US EPA Method TO-15 and addressed measurements of sixty-three volatile organic compounds. Laboratory data was evaluated according to the quality assurance / quality control requirements of the New York State Department of Environmental Conservation's Analytical Services Protocol (ASP), September 1989, Rev. 07/2005. When the required protocol was not followed, the current EPA Region II Functional Guidelines (SOP HW-31, Rev. #4, October 2006, Volatile Organic Analysis of Ambient Air in Canisters by Method TO-15) was used as a technical reference.

The chloroethane concentration found in SV-1 has been qualified as an estimation due to a high surrogate standard recovery.

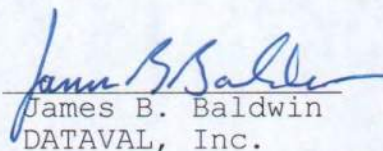
The trichloroethene concentration found in SV-3 has been qualified as an estimation due to a high spiked sample recovery.

## CORRECTNESS AND USABILITY

Reported data should be considered technically defensible and completely usable in its present form. Results presenting a usable estimation of the conditions at the time of sampling have been flagged "J". Estimated data should be used with caution. A detailed discussion of the review process follows.

Two facts should be considered by all data users. No compound concentration, even if it has passed all QC testing, can be guaranteed to be accurate. Strict QC serves to increase confidence in data, but any value potentially contains error. Secondly, DATAVAL, Inc. guarantees the quality of this data assessment. However, DATAVAL, Inc. does not warrant any interpretation or utilization of this data by a third party.

Reviewer's signature:

 Date: 25 May 17  
James B. Baldwin  
DATAVAL, Inc.



### SAMPLE HISTORY

Analyte concentrations can deteriorate with time due to chemical instability, bacterial degradation or volatility. Samples that are not properly preserved or are not analyzed within established holding times may no longer be considered representative. Holding times are calculated from the date of sampling. TO-15 samples must be analyzed within 14 days of collection.

This sample delivery group contained six air samples that were collected in 1-liter SUMMA canisters. Sampling was completed on 03Mar17. The canisters were shipped back to the laboratory, via FedEx, on 04Mar17 and were received on 07Mar17. Although the sample canisters were received intact, custody seals were not present on the packaging. It is noted that SV-3 was collected in a 1.4-liter canister to facilitate the preparation of MS/MSD samples.

Although the SUMMA canisters were set in the laboratory to collect 8-hour samples, sampling was terminated at a gauge reading of -5"Hg to comply with the ASP requirement of  $5 \pm 1$ "Hg. The discrepancies between vacuum readings recorded following sampling and prior to analysis are assumed to reflect the quality of the cylinder vacuum gauges.

SAMPLE	PRIOR TO SHIPMENT ("Hg)	PRIOR TO SAMPLING ("Hg)	POST SAMPLING ("Hg)	LAB ANALYSIS ("Hg)
SV-1	-30	-30	-5	-2
SV-2	-30	-30	-5	-1
SV-3	-30	-30	-5	-1
SV-4	-30	-28	-5	-3
DUPLICATE	-30	-28	-5	-3
AMBIENT AIR	-30	-30	-5	-1

The analysis of this group of samples was completed between 08Mar17 and 09Mar17, satisfying the ASP holding time limitation.

It is noted that sampling start and stop times were not provided. This made it impossible to determine if the proper sampling rate was maintained during sample collection.

### CANISTER CERTIFICATION

The canisters used for this project were pressure tested at 30 psig for 24 hours. Each canister demonstrated a change  $\leq 0.5$  psig over this period.

The canisters for this project were cleaned in four batches. A blank analysis of a clean canister from each of these batches was free of targeted analyte contamination.

### BLANKS

Blanks are analyzed to evaluate various sources of sample contam-



ination. Trip Blanks monitor sampling activities, sample transport, and storage. Method blanks are analyzed to verify instrument integrity. Samples are considered compromised by conditions causing contamination in any blank.

One method blank was analyzed with this group of samples. This blank demonstrated acceptable chromatography and were free of targeted analyte contamination.

#### MS TUNING

Mass spectrometer tuning and performance criteria are established to ensure sufficient mass resolution and sensitivity to accurately detect and identify targeted analytes. Verification is accomplished using a certified standard.

BFB ion abundance criteria was reported from standards run before the initial instrument calibration and prior to the analysis of program samples on 08Mar17. Both of these checks satisfied the ASP acceptance criteria.

#### CALIBRATION

Requirements for instrument calibration are established to ensure that laboratory equipment is capable of producing accurate, quantitative data. Initial calibrations demonstrate a range through which measurements may be made. Continuing calibration check standards verify instrument stability.

The initial instrument calibration was performed on 27Feb17. Standards of 0.04, 0.10, 0.15, 0.30, 0.50, 0.75, 1.0, 1.25, 1.50 and 2.0 ppbV were included. Each targeted analyte produced the required levels of instrument response and demonstrated an acceptable degree of linearity during this calibration.

A continuing calibration check standard was analyzed on 08Mar17, prior to the 24-hour period of instrument operation that included samples from this program. When compared to the initial calibration, each targeted analyte produced the required levels of instrument response and demonstrated an acceptable degree of linearity during this check.

#### SURROGATES

Each sample, blank and standard is spiked with surrogate compounds prior to analysis. The structures of surrogates are similar to analytes of interest, but they are not normally found in environmental samples. Surrogate recoveries are monitored to evaluate overall laboratory performance and the efficiency of laboratory technique.

Although surrogate summary sheets were properly prepared, an incorrect acceptance criteria was applied. When compared to the ASP requirements, unacceptably high recoveries were reported for the bromofluorobenzene additions to SV-1 and SV-2. The chloroethane (CLEANE) concentration found in SV-1 has been qualified as an estimation based on these indications of positive bias. The remaining results from SV-1 and SV-2 were negative and remain unqualified.



INTERNAL STANDARDS

Internal standards are added to each sample, blank and standard just prior to injection. Analyte concentrations are calculated relative to the response of a specific internal standard. Internal standard performance criteria ensure that GC/MS sensitivity and response are stable during the analysis of each sample. The area of internal standard peaks may not vary by more than 40%. When compared to the preceding calibration check, retention times may not vary by more than 10 seconds.

The laboratory recorded the response of each internal standard addition to this group of samples and the response obtained from the preceding CCV standard. Although the control limits based on the response of the CCV were not reported, they were calculated by this reviewer. When compared to these limits, unacceptable performance was reported for the chlorobenzene-d5 additions to SV-1, SV-2, SV-3 and the DUPLICATE. This performance had no impact on reported data, however, because none of the targeted analytes were associated with this internal standard.

Internal standard retention times were not addressed. The ASP retention time acceptance criteria was calculated by this reviewer. The retention times produced by each program sample satisfied these requirements.

MATRIX SPIKES / MATRIX SPIKE DUPLICATES / MATRIX SPIKED BLANKS

Matrix spiking refers to the addition of known analyte concentrations to a sample, prior to analysis. Analyte recoveries provide an indication of laboratory accuracy. The analysis of a duplicate spiked aliquot provides a measurement of precision.

SV-3 was selected for matrix spiking. The six targeted analytes were added to two volumes of this sample. The recoveries reported for these additions included a high trichloroethene (155%) result. The trichloroethene (TCE) concentration found in SV-3 has been qualified as an estimation based on this indication of positive bias.

A pair of spiked blanks (LCS/LCSD) was also analyzed with this delivery group. Both of these spiked blanks produced acceptable recoveries of the six targeted analytes.

DUPLICATES

Two aliquots of the same sample are processed separately through all aspects of sample preparation and analysis. Results produced by the analysis of this pair of samples are compared as a measurement of precision. Poor precision may be indicative of sample non-homogeneity, method defects, or poor laboratory technique.

The blind duplicate sample that was included in this delivery group was not identified.

REPORTED ANALYTES

Formal reports were provided for each sample. The data package also included total ion chromatograms and raw instrument print-



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outs. Reference mass spectra were provided to confirm the identification of each analyte that was detected in this group of samples.

# SUMMARY OF QUALIFIED DATA

691 ST PAUL SITE

SAMPLED 3/3/2017

	SURROGATE CLEANE	SPIKE TCE
SV-1	(C1703015-01)	
SV-2	(C1703015-02)	
SV-3	(C1703015-03)	
SV-2	(C1703015-04)	40J
DUPLICATE	(C1703015-05)	
AMBIENT AIR	(C1703015-06)	



## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-001A

Client Sample ID: SV-1  
 Tag Number: 542.256  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY METHOD TO15		TO-15				Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 1:28:00 PM
Chloroethane	0.34 J	0.40 J		ug/m3	1	3/8/2017 1:28:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 1:28:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 1:28:00 PM
Trichloroethene	< 0.81	0.81		ug/m3	1	3/8/2017 1:28:00 PM
Vinyl chloride	< 0.38	0.38		ug/m3	1	3/8/2017 1:28:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

7/15

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

Page 1 of 6



## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-002A

Client Sample ID: SV-2  
 Tag Number: 237.402  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY METHOD TO15		TO-15				Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 2:09:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	3/8/2017 2:09:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 2:09:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 2:09:00 PM
Trichloroethene	< 0.81	0.81		ug/m3	1	3/8/2017 2:09:00 PM
Vinyl chloride	< 0.38	0.38		ug/m3	1	3/8/2017 2:09:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

JHS

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

Page 2 of 6



## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-003A

Client Sample ID: SV-3  
 Tag Number: 1206.249  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY METHOD TO15		TO-15				Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 2:48:00 PM
Chloroethane -	0.29	0.40	J	ug/m3	1	3/8/2017 2:48:00 PM
cis-1,2-Dichloroethene -	0.71	0.59		ug/m3	1	3/8/2017 2:48:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 2:48:00 PM
Trichloroethene -	40 J	8.1		ug/m3	10	3/8/2017 11:34:00 PM
Vinyl chloride	< 0.38	0.38		ug/m3	1	3/8/2017 2:48:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

11AS

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

Page 3 of 6



## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-004A

Client Sample ID: SV-4  
 Tag Number: 290.299  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY METHOD TO15		TO-15			Analyst: RJP	
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 4:59:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	3/8/2017 4:59:00 PM
cis-1,2-Dichloroethene -	4.1	0.59		ug/m3	1	3/8/2017 4:59:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 4:59:00 PM
Trichloroethene -	48	8.1		ug/m3	10	3/9/2017 12:47:00 AM
Vinyl chloride	< 0.38	0.38		ug/m3	1	3/8/2017 4:59:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

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## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-005A

Client Sample ID: Duplicate  
 Tag Number: 1184.299  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY METHOD TO15		TO-15				Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 5:39:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	3/8/2017 5:39:00 PM
cis-1,2-Dichloroethene -	3.9	0.59		ug/m3	1	3/8/2017 5:39:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 5:39:00 PM
Trichloroethene -	47	8.1		ug/m3	10	3/9/2017 2:01:00 AM
Vinyl chloride	< 0.38	0.38		ug/m3	1	3/8/2017 5:39:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

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## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-006A

Client Sample ID: Ambient Air  
 Tag Number: 556.267  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC				TO-15		Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 12:48:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	3/8/2017 12:48:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 12:48:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 12:48:00 PM
Trichloroethene	< 0.21	0.21		ug/m3	1	3/8/2017 12:48:00 PM
Vinyl chloride	< 0.10	0.10		ug/m3	1	3/8/2017 12:48:00 PM

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

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Date: 27-Mar-17



CENTEK LABORATORIES, LLC

# QC SUMMARY REPORT SURROGATE RECOVERIES

CLIENT: LaBella Associates, P.C.

Work Order: C1703015

Project: 691 and 705 St Paul St

Test No: TO-15

Matrix: A

Sample ID	BR4FBZ								
ALCSIUG-030817	105								
ALCSIUGD-030817	104								
AMBIUG-030817	85.0								
C1703015-001A	239 *								
C1703015-002A	235 *								
C1703015-003A	97.0								
C1703015-003A MS	170 *								
C1703015-003A MSD	173 *								
C1703015-004A	100								
C1703015-005A	98.0								
C1703015-006A	87.0								

Acronym	Surrogate	QC Limits
BR4FBZ	= Bromofluorobenzene	<del>70-130</del> 80-120

\* Surrogate recovery outside acceptance limits



## Centek Laboratories, LLC

## GC/MS QA-QC Check Report

June File : C:\HPCHEM\1\DATA\AO030803.D

June Time : 8 Mar 2017 10:33 am

Daily Calibration File : C:\HPCHEM\1\DATA\AO030803.D

File	Sample	DL	Surrogate Recovery %	(BFB)		
				(IS1)	(IS2)	(IS3)
				51607	232876	196939
				30964	139726	118163
Internal Standard Responses						
0030804.D	ALCS1UG-030817	105		54344	232438	196865
0030805.D	AMB1UG-030817	85		51456	199472	169731
0030806.D	C1703015-006A	87		47960	186316	173836
0030807.D	C1703015-001A	239*		59338	270574	278336
0030808.D	C1703015-002A	235*		72867	317032	316562*
0030809.D	C1703015-003A	166*		80896	366728	336237*
0030810.D	C1703015-003A MS	170*		80202	356830	344072*
0030811.D	C1703015-003A MSD	173*		79391	370639	337498*
0030812.D	C1703015-004A	190*		76299	343779	323940*
0030813.D	C1703015-005A	187*		77699	348681	330439*
0030822.D	C1703015-003A 10X	97		53827	227536	201273
0030824.D	C1703015-004A 10X	100		46536	210199	197259
0030826.D	C1703015-005A 10X	98		46525	198463	190799
0030828.D	ALCS1UGD-030817	104		45106	199550	175594

t - fails 24hr time check \* - fails criteria

Created: Mon Mar 27 11:04:40 2017 MSD #1/

Retention Times			
CCV	9.64	12.03	16.90
C1703015-1	9.65	12.04	16.90
-2	9.65	12.03	16.90
-3	9.65	12.03	16.90
-3 10X	9.65	12.04	16.90
-4	9.65	12.03	16.90
-4 10X	9.65	12.03	16.90
-5	9.65	12.03	16.90
-5 10X	9.65	12.03	16.90
-6	9.65	12.03	16.90



Date: 27-Mar-17

## CENTEK LABORATORIES, LLC

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.

Work Order: C1703015

Project: 691 and 705 St Paul St

TestCode: 0.25CT-TCE-VC

Sample ID	ALCS1UG-030817	SampType: LCS	TestCode: 0.25CT-TCE-	Units: ppbv	Prep Date:	RunNo: 12014					
Client ID: ZZZZZ		Batch ID: R12014	TestNo: TO-15		Analysis Date: 3/8/2017	SeqNo: 140535					
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	1.060	0.15	1	0	106	70	130				
Chloroethane	0.9000	0.15	1	0	90.0	70	130				
cis-1,2-Dichloroethene	1.100	0.15	1	0	110	70	130				
trans-1,2-Dichloroethene	1.070	0.15	1	0	107	70	130				
Trichloroethene	1.060	0.040	1	0	106	70	130				
Vinyl chloride	0.8700	0.040	1	0	87.0	70	130				

Sample ID	ALCS1UGD-030817	SampType: LCSD	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12014					
Client ID: ZZZZZ		Batch ID: R12014	TestNo: TO-15		Analysis Date: 3/9/2017	SeqNo: 140536					
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	1.100	0.15	1	0	110	70	130	1.06	3.70	30	
Chloroethane	1.000	0.15	1	0	100	70	130	0.9	10.5	30	
cis-1,2-Dichloroethene	1.130	0.15	1	0	113	70	130	1.1	2.69	30	
trans-1,2-Dichloroethene	1.100	0.15	1	0	110	70	130	1.07	2.76	30	
Trichloroethene	1.100	0.040	1	0	110	70	130	1.06	3.70	30	
Vinyl chloride	0.9100	0.040	1	0	91.0	70	130	0.87	4.49	30	

Qualifiers: J Results reported are not blank corrected  
 S Analyte detected below quantitation limit  
 E Estimated Value above quantitation range  
 ND Not Detected at the Limit of Detection  
 H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits

Page 1 of 1



Date: 27-Mar-17

## CENTEK LABORATORIES, LLC

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.

Work Order: C1703015

Project: 691 and 705 St Paul St

TestCode: 0.25CT-TCE-VC

Sample ID	AMB1UG-030817	SampType:	MBLK	TestCode: 0.25CT-TCE-	Units: ppbv	Prep Date:	RunNo: 12014				
Client ID:	ZZZZ	Batch ID:	R12014	TestNo: TO-15		Analysis Date: 3/8/2017	SeqNo: 140534				
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	< 0.15	✓	0.15								
Chloroethane	< 0.15		0.15								
cis-1,2-Dichloroethene	< 0.15		0.15								
trans-1,2-Dichloroethene	< 0.15		0.15								
Trichloroethene	< 0.040		0.040								
Vinyl chloride	< 0.040		0.040								

## Qualifiers:

J Results reported are not blank corrected  
 S Analyte detected below quantization limit  
 S Spike Recovery outside accepted recovery limits

E Estimated Value above quantitation range  
 ND Not Detected at the Limit of Detection

H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits

Page 1 of 1



Date: 27-Mar-17

## CENTEK LABORATORIES, LLC

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.

Work Order: C1703015

Project: 691 and 705 St Paul St

TestCode: 1ugM3\_TO15

Sample ID	C1703015-003A MS	SampType: MS	TestCode: 1ugM3_TO15	Units: ppbV	Prep Date:	RunNo: 12014					
Client ID:	SV-3	Batch ID: R12014	TestNo: TO-15		Analysis Date: 3/8/2017	SeqNo: 140546					
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	1.090	0.15	1	0	109	70	130				
Chloroethane	0.8100	0.15	1	0.11	70.0	70	130				
cis-1,2-Dichloroethene	1.400	0.15	1	0.18	122	70	130				
trans-1,2-Dichloroethene	1.140	0.15	1	0	114	70	130				
Trichloroethene	9.890	0.15	1	8.34	155	70	130				S
Vinyl chloride	0.7200	0.15	1	0	72.0	70	130				

Sample ID C1703015-003A MS		SampType: MSD	TestCode: 1ugM3_TO15		Units: ppbV	Prep Date:		RunNo: 12014			
Client ID: SV-3		Batch ID: R12014	TestNo: TO-15			Analysis Date: 3/8/2017		SeqNo: 140547			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	1.160	0.15	1	0	116	70	130	1.09	6.22	30	
Chloroethane	0.8600	0.15	1	0.11	75.0	70	130	0.81	5.99	30	
cis-1,2-Dichloroethene	1.440	0.15	1	0.18	126	70	130	1.4	2.82	30	
trans-1,2-Dichloroethene	1.180	0.15	1	0	118	70	130	1.14	3.45	30	
Trichloroethene	9.550	0.15	1	8.34	121	70	130	9.89	3.50	30	
Vinyl chloride	0.7500	0.15	1	0	75.0	70	130	0.72	4.08	30	

Qualifiers: J Results reported are not blank corrected  
 S Analyte detected below quantitation limit  
 E Estimated Value above quantitation range  
 ND Not Detected at the Limit of Detection  
 H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits

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## APPENDIX 3

### Laboratory Analytical Reports



## TO-15 Package Review Checklist

Client: La Bella Project: 691 St Paul St. SDG: C1704014

		YES	NO	NA
Analytical Results	Present and Complete	✓		
TIC's present	Present and Complete	✓		
	Holding Times Met	✓		

Comments: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

Chain-of-Custody	Present and Complete	✓		
Surrogate Recovery	Present and Complete	✓		
	Recoveries within limits	✓		
	Sample(s) reanalyzed			✓
Internal Standards Recovery	Present and Complete	✓		
	Recoveries within limits	✓		
	Sample(s) reanalyzed			✓

Comments: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

Lab Control Sample (LCS)	Present and Complete	✓		
	Recoveries within limits	✓		
Lab Control Sample Dupe (LCSD)	Present and Complete	✓		
	Recoveries within limits	✓		
MS/MSD	Present and Complete	✓		
	Recoveries within limits	✓		

Comments: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

Sample Raw Data	Present and Complete	✓		
	Spectra present for all samples	✓		

Comments: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

## TO-15 Package Review Checklist

Client: LaBella Project: 691 St. Paul St SDG: C1704014

		<u>YES</u>	<u>NO</u>	<u>NA</u>
<b>Standards Data</b>				
Initial Calibration Summary	Present and Complete	✓	—	—
	Calibration(s) met criteria	✓	—	—
Continuing Calibration Summary	Present and Complete	✓	—	—
	Calibration(s) met criteria	✓	—	—
Standards Raw Data	Present and Complete	✓	—	—
Comments: _____				

<b>Raw Quality Control Data</b>				
Tune Criteria Report	Present and Complete	✓	—	—
Method Blank Data	MB Results <PQL	✓	—	—
	Associated results flagged "B"	—	—	✓
LCS sample data	Present and Complete	✓	—	—
LCSD sample data	Present and Complete	✓	—	—
MS/MSD sample data	Present and Complete	✓	—	—
Comments: _____				

<b>Logbooks</b>				
Injection Log	Present and Complete	✓	—	—
Standards Log	Present and Complete	✓	—	—
Can Cleaning Log	Present and Complete	✓	—	—
	Raw Data Present	✓	—	—
Calculation sheet	Present and Complete	✓	—	—
IDL's	Present and Complete	✓	—	—
Bottle Order Form	Present and Complete	✓	—	—
Sample Tracking Form	Present and Complete	✓	—	—

Additional Comments: \_\_\_\_\_

Section Supervisor: Walt Dahl Date: 5/8/17

QC Supervisor: Michael Date: 5/8/17



## CENTEK LABORATORIES, LLC

143 Midler Park Drive • Syracuse, NY 13206

Phone (315) 431-9730 • Emergency 24/7 (315) 416-2752

NYSDOH ELAP

Certificate No. 11830

### **Analytical Report**

Daniel Noll  
LaBella Associates, P.C.  
300 State Street, Suite 201  
Rochester, NY 14614

Wednesday, April 12, 2017

Order No.: C1704014

TEL: (585) 454-6110

FAX (585) 454-3066

RE: 691 St Paul Street

Dear Daniel Noll:

Centek Laboratories, LLC received 15 sample(s) on 4/5/2017 for the analyses presented in the following report.

I certify that this data package is in compliance with the terms and conditions of the Contract, both technically and for completeness. Release of the data contained in this hardcopy data package and/or in the computer readable data submitted has been authorized by the Laboratory Manager or his designee, as verified by the following signature.

All method blanks, laboratory spikes, and/or matrix spikes met quality assurance objective except as indicated in the case narrative. All samples were received and analyzed within the EPA recommended holding times. Test results are not Method Blank (MB) corrected for contamination.

Centek Laboratories is distinctively qualified to meet your needs for precise and timely volatile organic compound analysis. We perform all analyses according to EPA, NIOSH or OSHA-approved analytical methods. Centek Laboratories is dedicated to providing quality analyses and exceptional customer service. Samples were analyzed using the methods outlined in the following references:

Compendium of Methods for the Determination of Toxic Organic Compounds, Compendium Method TO-15, January 1999.

Centek Laboratories SOP TS-80

Analytical results relate to samples as received at laboratory. We do our best to make our reporting format clear and understandable and hope you are thoroughly satisfied with our services.

Please contact your client service representative at (315) 431-9730 or myself, if you would like any additional information regarding this report.



This report cannot be reproduced except in its entirety, without prior written authorization.

Sincerely,



William Dobbin  
Lead Technical Director

Disclaimer: The test results and procedures utilized, and laboratory interpretations of the data obtained by Centek as contained in this report are believed by Centek to be accurate and reliable for sample(s) tested. In accepting this report, the customer agrees that the full extent of any and all liability for actual and consequential damages of Centek for the services performed shall be equal to the fee charged to the customer for the services as liquidated damages. ELAP does not offer certification for the following parameters by this method at present time, they are: 4-ethyltoluene, ethyl acetate, propylene, tetrahydrofuran, 4-PCH, sulfur derived and silicon series compounds.

#### Centek Laboratories, LLC Terms and Conditions

##### Sample Submission

All samples sent to Centek Laboratories should be accompanied by our Request for Analysis Form or Chain of Custody Form. A Chain of Custody will be provided with each order shipped for all sampling events, or if needed, one is available at our website [www.CentekLabs.com](http://www.CentekLabs.com). Samples received after 3:00pm are considered to be a part of the next day's business.

##### Sample Media

Samples can be collected in an canister or a Tedlar bag. Depending on your analytical needs, Centek Laboratories may receive a bulk, liquid, soil or other matrix sample for headspace analysis.

##### Blanks

Every sample is run with a surrogate or tracer compound at a pre-established concentration. The surrogate compound run with each sample is used as a standard to measure the performance of each run of the instrument. If required, a Minican can be provided containing nitrogen to be run as a trip blank with your samples.

##### Sampling Equipment

Centek Laboratories will be happy to provide the canisters to carry-out your sampling event at no charge. The necessary accessories, such as regulators, tubing or personal sampling belts, are also provided to meet your sampling needs. The customer is responsible for all shipping charges to the client's destination and return shipping to the laboratory. Client assumes all responsibility for lost, stolen and any damages of equipment.

##### Turn Around time (TAT)

Centek Laboratories will provide results to its clients in one business-week by 6:00pm EST after receipt of samples. For example, if samples are received on a Monday they are due on the following Monday by 6:00pm EST. Results are faxed or emailed to the requested location indicated on the Chain of Custody. Non-routine analysis may require more than the one business-week turnaround time. Please confirm non-routine sample turnaround times.

### Reporting

Results are emailed or faxed at no additional charge. A hard copy of the result report is mailed within 24 hours of the faxing or emailing of your results. Cat "B" like packages are within 3-4 weeks from time of analysis. Standard Electronic Disk Deliverables (EDD) is also available at no additional charge.

### Payment Terms

Payment for all purchases shall be due within 30 days from date of invoice. The client agrees to pay a finance charge of 1.5% per month on the overdue balance and cost of collection, including attorney fees, if collection proceedings are necessary. You must have a completed credit application on file to extend credit. Purchase orders or checks information must be submitted for us to release results

### Rush Turnaround Samples

Expedited turn around times is available. Please confirm rush turnaround times with Client Services before submitting samples.

Applicable Surcharges for Rush Turnaround Samples:

Same day TAT = 200%

Next business day TAT by Noon = 150%

Next business day TAT by 6:00pm = 100%

Second business day TAT by 6:00pm = 75%

Third business day TAT by 6:00pm = 50%

Fourth business day TAT by 6:00pm = 35%

Fifth business day = Standard

### Statement of Confidentiality

Centek Laboratories, LLC is aware of the importance of the confidentiality of results to many of our clients. Your name and data will be held in the strictest of confidence. We will not accept business that may constitute a conflict of interest. We commonly sign Confidential Nondisclosure Agreements with clients prior to beginning work. All research, results and reports will be kept strictly confidential. Secrecy Agreements and Disclosure Statements will be signed for the client if so specified. Results will be provided only to the addressee specified on the Chain of Custody Form submitted with the samples unless law requires release. Written permission is required from the addressee to release results to any other party.

### Limitation on Liability

Centek Laboratories, LLC warrants the test results to be accurate to the methodology and sample type for each sample submitted to Centek Laboratories, LLC. In no event shall Centek Laboratories, LLC be liable for direct, indirect, special, punitive, incidental, exemplary or consequential damages, or any damages whatsoever, even if Centek Laboratories, LLC has been previously advised of the possibility of such damages whether in an action under contract, negligence, or any other theory, arising out of or in connection with the use, inability to use or performance of the information, services, products and materials available from the laboratory or this site. These limitations shall apply notwithstanding any failure of essential purpose of any limited remedy. Because some jurisdictions do not allow limitations on how long an implied warranty lasts, or the exclusion or limitation of liability for consequential or incidental damages, the above limitations may not apply to you. This is a comprehensive limitation of

liability that applies to all damages of any kind, including (without limitation) compensatory, direct, indirect or consequential damages, loss of data, income or profit and or loss of or damage to property and claims of third parties.



# **ASP CAT B DELIVERABLE PACKAGE**

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## CENTEK LABORATORIES, LLC

Date: 08-May-17

CLIENT: LaBella Associates, P.C.

Project: 691 St Paul Street

Lab Order: C1704014

## CASE NARRATIVE

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Samples were analyzed using the methods outlined in the following references:

Centek Laboratories, LLC SOP TS-80

Compendium of Methods for the Determination of Toxic Organic Compounds, Compendium Method TO-15, January 1999

All method blanks, laboratory spikes, and/or matrix spikes met quality assurance objective except as indicated in the corrective action report(s). All samples were received and analyzed within the EPA recommended holding times. Test results are not Method Blank (MB) corrected for contamination.

### NYSDEC ASP samples:

Canisters should be evacuated to a reading of less than or equal to 50 millitorr prior to shipment to sampling personnel. The vacuum in the canister will be field checked prior to sampling, and must read 28" of Hg ( $\pm 2$ ", vacuum, absolute) before a sample can be collected. After the sample has been collected, the pressure of the canister will be read and recorded again, and must be 5" of Hg ( $\pm 1$ ", vacuum, absolute) for the sample to be valid. Once received at the laboratory, the canister vacuum should be confirmed to be 5" of Hg,  $\pm 1$ ". Please record and report the pressure/vacuum of received canisters on the sample receipt paperwork. A pressure/vacuum reading should also be taken just prior to the withdrawal of sample from the canister, and recorded on the sample preparation log sheet. All regulators are calibrated to meet these requirements before they leave the laboratory. However, due to environmental conditions and use of the equipment Centek can not guarantee that this criteria can always be achieved.







# CEN TEK LABORATORIES, LLC

Date: 08-May-17

CLIENT: LaBella Associates, P.C.  
Project: 691 St Paul Street  
Lab Order: C1704014

## Work Order Sample Summary

Lab Sample ID	Client Sample ID	Tag Number	Collection Date	Date Received
C1704014-001A	691-B19-IAQ-1	1321.1163	4/1/2017	4/5/2017
C1704014-002A	691-B19-IAQ-2	1188.306	4/1/2017	4/5/2017
C1704014-003A	691-B19-SVI	362.446	4/1/2017	4/5/2017
C1704014-004A	691-B15-IAQ	87.299	4/1/2017	4/5/2017
C1704014-005A	691-B15-SVI	550.266	4/1/2017	4/5/2017
C1704014-006A	691-Outdoor-04012017	240.340	4/1/2017	4/5/2017
C1704014-007A	691-Duplicate	101.299	4/1/2017	4/5/2017

**CLIENT:** LaBella Associates, P.C.  
**Project:** 691 St Paul Street  
**Lab Order:** C1704014

## Work Order Sample Summary

Lab Sample ID	Client Sample ID	Tag Number	Collection Date	Date Received
C1704014-008A	691-NE-IAQ	94.379	4/1/2017	4/5/2017
C1704014-009A	691-NE-SVI	1186.145	4/1/2017	4/5/2017
C1704014-010A	691-SB1-IAQ	170.387	4/1/2017	4/5/2017
C1704014-011A	691-SB1-SVI	203.372	4/1/2017	4/5/2017
C1704014-012A	691-SB5A-IAQ	367.251	4/1/2017	4/5/2017
C1704014-013A	691-SB5B-IAQ	285.344	4/1/2017	4/5/2017
C1704014-014A	691-SB5A-SVI	475.1170	4/1/2017	4/5/2017
C1704014-015A	691-SB5B-SVI	467.1167	4/1/2017	4/5/2017



# CEN TEK LABORATORIES, LLC

## Sample Receipt Checklist

Client Name LABELLA - ROCHESTER

Date and Time Receive

4/5/2017

Work Order Number C1704914

Received by NM

Checklist completed by

Signature

Date

Reviewed by

Initials

Date

Matrix:

Carrier name: FedEx Ground

Shipping container/cooler in good condition?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	Not Present <input type="checkbox"/>
Custody seals intact on shipping container/cooler?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	Not Present <input checked="" type="checkbox"/>
Custody seals intact on sample bottles?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	Not Present <input checked="" type="checkbox"/>
Chain of custody present?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Chain of custody signed when relinquished and received?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Chain of custody agrees with sample labels?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Samples in proper container/bottle?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Sample containers intact?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Sufficient sample volume for indicated test?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
All samples received within holding time?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Container/Temp Blank temperature in compliance?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Water - VOA vials have zero headspace?	No VOA vials submitted <input checked="" type="checkbox"/>	Yes <input type="checkbox"/>	No <input type="checkbox"/>
Water - pH acceptable upon receipt?	Yes <input type="checkbox"/>	No <input checked="" type="checkbox"/>	

Adjusted? \_\_\_\_\_ Checked by \_\_\_\_\_

Any No and/or NA (not applicable) response must be detailed in the comments section below

Client contacted \_\_\_\_\_ Date contacted: \_\_\_\_\_ Person contacted \_\_\_\_\_

Contacted by: \_\_\_\_\_ Regarding: \_\_\_\_\_

Comments: \_\_\_\_\_

Corrective Action \_\_\_\_\_



Lab Order: C1704014  
 Client: LaBella Associates, P.C.  
 Project: 691 St Paul Street

## DATES REPORT

Sample ID	Client Sample ID	Collection Date	Matrix	Test Name	TCLP Date	Prep Date	Analysis Date
C1704014-001A	691-B19-IAQ-1	4/1/2017	Air	1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/6/2017
C1704014-002A	691-B19-IAQ-2			1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
C1704014-003A	691-B19-SVI			1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
				1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
C1704014-004A	691-B15-IAQ			1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
C1704014-005A	691-B15-SVI			1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
				1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
C1704014-006A	691-Outdoor-04012017			1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
C1704014-007A	691-Duplicate			1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
C1704014-008A	691-NE-IAQ			1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
C1704014-009A	691-NE-SVI			1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
C1704014-010A	691-SB1-IAQ			1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
C1704014-011A	691-SB1-SVI			1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
				1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
C1704014-012A	691-SB5A-IAQ			1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
				1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
C1704014-013A	691-SB5B-IAQ			1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
				1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
C1704014-014A	691-SB5A-SVI			1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
				1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
C1704014-015A	691-SB5B-SVI			1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017
				1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			4/7/2017



# CEN TEK LABORATORIES, LLC

*Air Quality Testing...It's a Gas*

143 Midler Park Drive \* Syracuse, NY 13206  
TEL: 315-431-9730 \* FAX: 315-431-9731

## CANISTER ORDER

**6417**

08-May-17

### SHIPPED TO:

Company: LaBella Associates, P.C.  
Contact: Benjamin Stracuzzi  
Address: 300 State Street, Suite 201  
Rochester, NY 14614  
Phone: (585) 454-6110  
Quote ID: 0  
Project:  
PO: 2170820

### Submitted By:

MadeBy: jan

Ship Date: 3/28/2017  
VIA: FedEx Ground  
Due Date: 3/29/2017

Bottle Code	Bottle Type	TEST(s)	QTY
MC1400CC	1.4L Mini-Can	1ug/m3 w/ 0.25ug/M3 CT-TCE-VC	1
MC1000CC	1L Mini-Can	1ug/m3 w/ 0.25ug/M3 CT-TCE-VC	20

Can / Reg ID	Description
87	1L Mini-Can - 1104 VI
94	1L Mini-Can - 1086 VI
101	1L Mini-Can - 1101 VI
106	1L Mini-Can - 1056 VI
130	1L Mini-Can - 1078 VI
139	1L Mini-Can - 1113 VI
145	Time-Set Reg - 640 VI
170	1L Mini-Can - 1141 VI
203	1L Mini-Can - 1158 VI
240	1L Mini-Can - 1172 VI
251	Time-Set Reg - 689 VI
258	Time-Set Reg - 696 VI
266	Time-Set Reg - 704 VI
268	Time-Set Reg - 706 VI
285	1L Mini-Can - 1061 VI
286	1L Mini-Can - 1262 VI
299	Time-Set Reg - 722 VI
306	Time-Set Reg - 729 VI
324	1L Mini-Can - 1287 VI
328	1L Mini-Can - 1291 VI
339	Time-Set Reg - 736 VI
340	Time-Set Reg - 737 VI
344	Time-Set Reg - 741 VI
362	1L Mini-Can - 1311 VI
367	1L Mini-Can - 1316 VI
372	Time-Set Reg - 746 VI
379	Time-Set Reg - 753 VI
387	Time-Set Reg - 761 VI
446	Time-Set Reg - 825 VI
467	1L Mini-Can - 1371 VI
475	1L Mini-Can - 1377 VI
550	1L Mini-Can - 118 VI

## SHIPPED TO:

Company: LaBella Associates, P.C.  
Contact: Benjamin Stracuzzi  
Address: 300 State Street, Suite 201  
Rochester, NY 14614  
Phone: (585) 454-6110  
Quote ID: 0  
Project:  
PO: 2170820

## Submitted By:

MadeBy: jan

Ship Date: 3/28/2017  
VIA: FedEx Ground  
Due Date: 3/29/2017

Bottle Code	Bottle Type	TEST(s)	QTY
1163	Time-Set Reg-0676 VI		
1166	Time-Set Reg-0791 VI		
1167	Time-Set Reg-0792 VI		
1168	Time-Set Reg-0793 VI		
1170	Time-Set Reg-0795 VI		
1171	Time-Set Reg-0796 VI		
1186	1L Mini-Can - 1235 VI		
1188	1L Mini-Can - 1256 VI		
1321	1.4L Mini-Can - 0252 VI		

Comments: 20 (1L) @ 8hr + Ms/MSD +helium shroud and detector +clay +tubing WAC 031317A-C, 032017 D-H



**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**ANALYTICAL RESULTS**

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-001A

Client Sample ID: 691-B19-1AQ-1  
Tag Number: 1321.1163  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>		<b>FLD</b>		<b>Analyst:</b>		
Lab Vacuum In	-4			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>		<b>TO-15</b>		<b>Analyst: RJP</b>		
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/6/2017 9:55:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	4/6/2017 9:55:00 PM
cis-1,2-Dichloroethene	0.33	0.15		ppbV	1	4/6/2017 9:55:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/6/2017 9:55:00 PM
Trichloroethene	0.40	0.040		ppbV	1	4/6/2017 9:55:00 PM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/6/2017 9:55:00 PM
Surr: Bromofluorobenzene	96.0	70-130		%REC	1	4/6/2017 9:55:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-001A

Client Sample ID: 691-B19-IAQ-1  
Tag Number: 1321.1163  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/6/2017 9:55:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/6/2017 9:55:00 PM
cis-1,2-Dichloroethene	1.3	0.59		ug/m3	1	4/6/2017 9:55:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/6/2017 9:55:00 PM
Trichloroethene	2.1	0.21		ug/m3	1	4/6/2017 9:55:00 PM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/6/2017 9:55:00 PM

Qualifiers: \*\* Quantitation Limit . Results reported are not blank corrected  
B Analyte detected in the associated Method Blank E Estimated Value above quantitation range  
H Holding times for preparation or analysis exceeded J Analyte detected below quantitation limit  
JN Non-routine analyte. Quantitation estimated. ND Not Detected at the Limit of Detection  
S Spike Recovery outside accepted recovery limits



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-002A

Client Sample ID: 691-B19-IAQ-2  
Tag Number: 1188.306  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>		<b>FLD</b>		<b>Analyst:</b>		
Lab Vacuum In	-9			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>		<b>TO-15</b>		<b>Analyst: RJP</b>		
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 12:09:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 12:09:00 AM
cis-1,2-Dichloroethene	0.29	0.15		ppbV	1	4/7/2017 12:09:00 AM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 12:09:00 AM
Trichloroethene	0.39	0.040		ppbV	1	4/7/2017 12:09:00 AM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 12:09:00 AM
Surr: Bromofluorobenzene	89.0	70-130		%REC	1	4/7/2017 12:09:00 AM

<b>Qualifiers:</b>	<b>**</b>	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

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# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-002A

Client Sample ID: 691-B19-1AQ-2  
Tag Number: 1188.306  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 12:09:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 12:09:00 AM
cis-1,2-Dichloroethene	1.1	0.59		ug/m3	1	4/7/2017 12:09:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 12:09:00 AM
Trichloroethene	2.1	0.21		ug/m3	1	4/7/2017 12:09:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 12:09:00 AM

Qualifiers: \*\* Quantitation Limit . Results reported are not blank corrected  
B Analyte detected in the associated Method Blank E Estimated Value above quantitation range  
H Holding times for preparation or analysis exceeded J Analyte detected below quantitation limit  
JN Non-routine analyte. Quantitation estimated. ND Not Detected at the Limit of Detection  
S Spike Recovery outside accepted recovery limits

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-003A

Client Sample ID: 691-B19-SVI  
Tag Number: 362.446  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						Analyst:
Lab Vacuum In	-9			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 3:23:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 3:23:00 PM
cis-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 3:23:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 3:23:00 PM
Trichloroethene	4.8	0.40		ppbV	10	4/7/2017 8:07:00 PM
Vinyl chloride	0.080	0.040		ppbV	1	4/7/2017 3:23:00 PM
Surr: Bromofluorobenzene	109	70-130		%REC	1	4/7/2017 3:23:00 PM

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-003A

Client Sample ID: 691-B19-SVI  
Tag Number: 362.446  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15				Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 3:23:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 3:23:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 3:23:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 3:23:00 PM
Trichloroethene	26	2.1		ug/m3	10	4/7/2017 8:07:00 PM
Vinyl chloride	0.20	0.10		ug/m3	1	4/7/2017 3:23:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-004A

Client Sample ID: 691-B15-1AQ  
Tag Number: 87.299  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
FIELD PARAMETERS		FLD		Analyst:		
Lab Vacuum In	-3			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 12:50:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 12:50:00 AM
cis-1,2-Dichloroethene	0.40	0.15		ppbV	1	4/7/2017 12:50:00 AM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 12:50:00 AM
Trichloroethene	0.41	0.040		ppbV	1	4/7/2017 12:50:00 AM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 12:50:00 AM
Surr: Bromofluorobenzene	92.0	70-130		%REC	1	4/7/2017 12:50:00 AM

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-004A

Client Sample ID: 691-B15-IAQ  
Tag Number: 87.299  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 12:50:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 12:50:00 AM
cis-1,2-Dichloroethene	1.6	0.59		ug/m3	1	4/7/2017 12:50:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 12:50:00 AM
Trichloroethene	2.2	0.21		ug/m3	1	4/7/2017 12:50:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 12:50:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-005A

Client Sample ID: 691-B15-SVI  
Tag Number: 550.266  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-2			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
		TO-15				Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 4:03:00 PM
Chloroethane	0.14	0.15	J	ppbV	1	4/7/2017 4:03:00 PM
cis-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 4:03:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 4:03:00 PM
Trichloroethene	40	3.6		ppbV	90	4/7/2017 9:20:00 PM
Vinyl chloride	0.11	0.040		ppbV	1	4/7/2017 4:03:00 PM
Surr: Bromofluorobenzene	118	70-130		%REC	1	4/7/2017 4:03:00 PM

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte, Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-005A

Client Sample ID: 691-B15-SVI  
Tag Number: 550.266  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:03:00 PM
Chloroethane	0.37	0.40	J	ug/m3	1	4/7/2017 4:03:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:03:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:03:00 PM
Trichloroethene	220	19		ug/m3	90	4/7/2017 9:20:00 PM
Vinyl chloride	0.28	0.10		ug/m3	1	4/7/2017 4:03:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.

Client Sample ID: 691-Outdoor-04012017

Lab Order: C1704014

Tag Number: 240.340

Project: 691 St Paul Street

Collection Date: 4/1/2017

Lab ID: C1704014-006A

Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-8			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 1:31:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 1:31:00 AM
cis-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 1:31:00 AM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 1:31:00 AM
Trichloroethene	< 0.040	0.040		ppbV	1	4/7/2017 1:31:00 AM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 1:31:00 AM
Surr: Bromofluorobenzene	88.0	70-130		%REC	1	4/7/2017 1:31:00 AM

Qualifiers:	**	Quantitation Limit	,	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte, Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-006A

Client Sample ID: 691-Outdoor-04012017  
Tag Number: 240.340  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 1:31:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 1:31:00 AM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 1:31:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 1:31:00 AM
Trichloroethene	< 0.21	0.21		ug/m3	1	4/7/2017 1:31:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 1:31:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-007A

Client Sample ID: 691-Duplicate  
Tag Number: 101.299  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>		<b>FLD</b>		<b>Analyst:</b>		
Lab Vacuum In	-3			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>		<b>TO-15</b>		<b>Analyst: RJP</b>		
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 2:12:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 2:12:00 AM
cis-1,2-Dichloroethene	0.42	0.15		ppbV	1	4/7/2017 2:12:00 AM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 2:12:00 AM
Trichloroethene	0.41	0.040		ppbV	1	4/7/2017 2:12:00 AM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 2:12:00 AM
Surr: Bromofluorobenzene	91.0	70-130		%REC	1	4/7/2017 2:12:00 AM

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-007A

Client Sample ID: 691-Duplicate  
Tag Number: 101.299  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 2:12:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 2:12:00 AM
cis-1,2-Dichloroethene	1.7	0.59		ug/m3	1	4/7/2017 2:12:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 2:12:00 AM
Trichloroethene	2.2	0.21		ug/m3	1	4/7/2017 2:12:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 2:12:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.

Client Sample ID: 691-NE-IAQ

Lab Order: C1704014

Tag Number: 94.379

Project: 691 St Paul Street

Collection Date: 4/1/2017

Lab ID: C1704014-008A

Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-1			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 2:52:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 2:52:00 AM
cis-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 2:52:00 AM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 2:52:00 AM
Trichloroethene	< 0.040	0.040		ppbV	1	4/7/2017 2:52:00 AM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 2:52:00 AM
Surr: Bromofluorobenzene	93.0	70-130		%REC	1	4/7/2017 2:52:00 AM

## Qualifiers:

\*\* Quantitation Limit

B Analyte detected in the associated Method Blank

H Holding times for preparation or analysis exceeded

JN Non-routine analyte. Quantitation estimated.

S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected

E Estimated Value above quantitation range

J Analyte detected below quantitation limit

ND Not Detected at the Limit of Detection

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# Centek Laboratories, LLC

Date: 04-May-17

**CLIENT:** LaBella Associates, P.C.  
**Lab Order:** C1704014  
**Project:** 691 St Paul Street  
**Lab ID:** C1704014-008A

**Client Sample ID:** 691-NE-IAQ  
**Tag Number:** 94.379  
**Collection Date:** 4/1/2017  
**Matrix:** AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 2:52:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 2:52:00 AM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 2:52:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 2:52:00 AM
Trichloroethene	< 0.21	0.21		ug/m3	1	4/7/2017 2:52:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 2:52:00 AM

**Qualifiers:** \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits  
 . Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-009A

Client Sample ID: 691-NE-SVI  
Tag Number: 1186.145  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>		<b>FLD</b>		<b>Analyst:</b>		
Lab Vacuum In	-3			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>		<b>TO-15</b>		<b>Analyst: RJP</b>		
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 5:28:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 5:28:00 PM
cis-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 5:28:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 5:28:00 PM
Trichloroethene	0.72	0.040		ppbV	1	4/7/2017 5:28:00 PM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 5:28:00 PM
Surr: Bromofluorobenzene	113	70-130		%REC	1	4/7/2017 5:28:00 PM

Qualifiers: \*\* Quantitation Limit , Results reported are not blank corrected  
B Analyte detected in the associated Method Blank E Estimated Value above quantitation range  
H Holding times for preparation or analysis exceeded J Analyte detected below quantitation limit  
JN Non-routine analyte. Quantitation estimated. ND Not Detected at the Limit of Detection  
S Spike Recovery outside accepted recovery limits

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-009A

Client Sample ID: 691-NE-SVI  
Tag Number: 1186.145  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 5:28:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 5:28:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 5:28:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 5:28:00 PM
Trichloroethene	3.9	0.21		ug/m3	1	4/7/2017 5:28:00 PM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 5:28:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.

Client Sample ID: 691-SB1-IAQ

Lab Order: C1704014

Tag Number: 170.387

Project: 691 St Paul Street

Collection Date: 4/1/2017

Lab ID: C1704014-010A

Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-5			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 3:33:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 3:33:00 AM
cis-1,2-Dichloroethene	0.66	0.15		ppbV	1	4/7/2017 3:33:00 AM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 3:33:00 AM
Trichloroethene	0.24	0.040		ppbV	1	4/7/2017 3:33:00 AM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 3:33:00 AM
Surr: Bromofluorobenzene	92.0	70-130		%REC	1	4/7/2017 3:33:00 AM

Qualifiers:	** Quantitation Limit	.	Results reported are not blank corrected
	B Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S Spike Recovery outside accepted recovery limits		

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# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-010A

Client Sample ID: 691-SB1-1AQ  
Tag Number: 170.387  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 3:33:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 3:33:00 AM
cis-1,2-Dichloroethene	2.6	0.59		ug/m3	1	4/7/2017 3:33:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 3:33:00 AM
Trichloroethene	1.3	0.21		ug/m3	1	4/7/2017 3:33:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 3:33:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte, Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-011A

Client Sample ID: 691-SB1-SVI  
Tag Number: 203.372  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-6			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 6:09:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 6:09:00 PM
cis-1,2-Dichloroethene	13	1.5		ppbV	10	4/7/2017 11:10:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 6:09:00 PM
Trichloroethene	0.74	0.040		ppbV	1	4/7/2017 6:09:00 PM
Vinyl chloride	0.15	0.040		ppbV	1	4/7/2017 6:09:00 PM
Surr: Bromofluorobenzene	109	70-130		%REC	1	4/7/2017 6:09:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte, Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

, Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-011A

Client Sample ID: 691-SB1-SVI  
Tag Number: 203.372  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 6:09:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 6:09:00 PM
cis-1,2-Dichloroethene	52	5.9		ug/m3	10	4/7/2017 11:10:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 6:09:00 PM
Trichloroethene	4.0	0.21		ug/m3	1	4/7/2017 6:09:00 PM
Vinyl chloride	0.38	0.10		ug/m3	1	4/7/2017 6:09:00 PM

Qualifiers:	** Quantitation Limit	Results reported are not blank corrected
	B Analyte detected in the associated Method Blank	E Estimated Value above quantitation range
	H Holding times for preparation or analysis exceeded	J Analyte detected below quantitation limit
	JN Non-routine analyte. Quantitation estimated.	ND Not Detected at the Limit of Detection
	S Spike Recovery outside accepted recovery limits	

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# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-012A

Client Sample ID: 691-SB5A-JAQ  
Tag Number: 367.251  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-2			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 4:13:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 4:13:00 AM
cis-1,2-Dichloroethene	3.3	1.5		ppbV	10	4/7/2017 1:50:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 4:13:00 AM
Trichloroethene	1.2	0.040		ppbV	1	4/7/2017 4:13:00 AM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 4:13:00 AM
Surr: Bromofluorobenzene	103	70-130		%REC	1	4/7/2017 4:13:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte, Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
/ Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-012A

Client Sample ID: 691-SB5A-IAQ  
Tag Number: 367.251  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:13:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 4:13:00 AM
cis-1,2-Dichloroethene	13	5.9		ug/m3	10	4/7/2017 1:50:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:13:00 AM
Trichloroethene	6.5	0.21		ug/m3	1	4/7/2017 4:13:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 4:13:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-013A

Client Sample ID: 691-SB5B-1AQ  
Tag Number: 285.344  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-3			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 4:53:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 4:53:00 AM
cis-1,2-Dichloroethene	3.1	1.5		ppbV	10	4/7/2017 2:27:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 4:53:00 AM
Trichloroethene	1.3	0.040		ppbV	1	4/7/2017 4:53:00 AM
Vinyl chloride	0.060	0.040		ppbV	1	4/7/2017 4:53:00 AM
Surr: Bromofluorobenzene	102	70-130		%REC	1	4/7/2017 4:53:00 AM

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1704014  
 Project: 691 St Paul Street  
 Lab ID: C1704014-013A

Client Sample ID: 691-SB5B-IAQ  
 Tag Number: 285.344  
 Collection Date: 4/1/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:53:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 4:53:00 AM
cis-1,2-Dichloroethene	12	5.9		ug/m3	10	4/7/2017 2:27:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:53:00 AM
Trichloroethene	6.8	0.21		ug/m3	1	4/7/2017 4:53:00 AM
Vinyl chloride	0.15	0.10		ug/m3	1	4/7/2017 4:53:00 AM

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits  
 . Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-014A

Client Sample ID: 691-SB5A-SVI  
Tag Number: 475.1170  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-2			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 6:50:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 6:50:00 PM
cis-1,2-Dichloroethene	44	14		ppbV	90	4/8/2017 1:00:00 AM
trans-1,2-Dichloroethene	0.41	0.15		ppbV	1	4/7/2017 6:50:00 PM
Trichloroethene	3.7	0.36		ppbV	9	4/8/2017 12:23:00 AM
Vinyl chloride	0.19	0.040		ppbV	1	4/7/2017 6:50:00 PM
Surr: Bromofluorobenzene	107	70-130		%REC	1	4/7/2017 6:50:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-014A

Client Sample ID: 691-SB5A-SVI  
Tag Number: 475.1170  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 6:50:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 6:50:00 PM
cis-1,2-Dichloroethene	170	55		ug/m3	90	4/8/2017 1:00:00 AM
trans-1,2-Dichloroethene	1.6	0.59		ug/m3	1	4/7/2017 6:50:00 PM
Trichloroethene	20	1.9		ug/m3	9	4/8/2017 12:23:00 AM
Vinyl chloride	0.49	0.10		ug/m3	1	4/7/2017 6:50:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte, Quantitation estimated,  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-015A

Client Sample ID: 691-SB5B-SVI  
Tag Number: 467.1167  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-1			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 7:30:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 7:30:00 PM
cis-1,2-Dichloroethene	5.8	1.5		ppbV	10	4/8/2017 1:37:00 AM
trans-1,2-Dichloroethene	0.18	0.15		ppbV	1	4/7/2017 7:30:00 PM
Trichloroethene	11	0.40		ppbV	10	4/8/2017 1:37:00 AM
Vinyl chloride	0.11	0.040		ppbV	1	4/7/2017 7:30:00 PM
Surr: Bromofluorobenzene	116	70-130		%REC	1	4/7/2017 7:30:00 PM

Qualifiers: \*\* Quantitation Limit . Results reported are not blank corrected  
B Analyte detected in the associated Method Blank E Estimated Value above quantitation range  
H Holding times for preparation or analysis exceeded J Analyte detected below quantitation limit  
JN Non-routine analyte. Quantitation estimated. ND Not Detected at the Limit of Detection  
S Spike Recovery outside accepted recovery limits

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-015A

Client Sample ID: 691-SB5B-SV1  
Tag Number: 467.1167  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15				Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 7:30:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 7:30:00 PM
cis-1,2-Dichloroethene	23	5.9		ug/m3	10	4/8/2017 1:37:00 AM
trans-1,2-Dichloroethene	0.71	0.59		ug/m3	1	4/7/2017 7:30:00 PM
Trichloroethene	57	2.1		ug/m3	10	4/8/2017 1:37:00 AM
Vinyl chloride	0.28	0.10		ug/m3	1	4/7/2017 7:30:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**QUALITY CONTROL SUMMARY**



**QC SUMMARY REPORT  
SURROGATE RECOVERIES**

Test No: TO-15

[illegible]

### \* Surrogate recovery outside acceptance limits



## GC/MS QA-QC Check Report

Tune File : C:\HPCHEM\1\DATA\AO040602.D

Tune Time : 6 Apr 2017 9:31 am

Daily Calibration File : C:\HPCHEM\1\DATA\AO040602.D

(BFB)

(IS1)

(IS2)

(IS3)

26713

121699

111761

File	Sample	DL	Surrogate Recovery %	Internal Standard Responses		
AO040603.D	ALCS1UG-040617	95		26450	123600	109377
AO040604.D	AMB1UG-040617	90		26038	118374	102379
AO040620.D	ALCS1UGD-040617	98		24809	117535	100863
AO040621.D	C1704014-001A	96		25396	113281	97086
AO040622.D	C1704014-001A MS	97		26903	116940	105999
AO040623.D	C1704014-001A MSD	99		26299	120898	106126
AO040624.D	C1704014-002A	89		24868	113360	102277
AO040625.D	C1704014-004A	92		25596	114646	102450
AO040626.D	C1704014-006A	88		25816	110508	98499
AO040627.D	C1704014-007A	91		24957	110993	99559
AO040628.D	C1704014-008A	93		25088	108028	95812
AO040629.D	C1704014-010A	92		24485	107880	98198
AO040630.D	C1704014-012A	103		25278	111053	100304
AO040631.D	C1704014-013A	102		25079	113761	101873

t - fails 24hr time check \* - fails criteria

Created: Thu May 04 12:00:40 2017 MSD #1/

## GC/MS QA-QC Check Report

Tune File : C:\HPCHEM\1\DATA\AO040702.D

Tune Time : 7 Apr 2017 11:20 am

Daily Calibration File : C:\HPCHEM\1\DATA\AO040702.D

(BFB)

(IS1)

(IS2)

(IS3)

23210

99215

88822

File	Sample	DL	Surrogate Recovery %	Internal Standard Responses		
AO040703.D	ALCS1UG-040717		97	22420	98698	86489
AO040704.D	AMB1UG-040717		86	21100	92197	83384
AO040706.D	C1704014-012A 10X		91	20808	86837	77606
AO040707.D	C1704014-013A 10X		92	19443	85901	76902
AO040708.D	C1704014-003A		109	20468	96070	92232
AO040709.D	C1704014-005A		118	24680	109137	104661
AO040710.D	C1704014-009A		113	27394	125972	121869
AO040711.D	C1704014-011A		109	30170	132762	121408
AO040712.D	C1704014-014A		107	31361	139148	131462
AO040713.D	C1704014-015A		116	31915	147799	139968
AO040714.D	C1704014-003A 10x		95	31326	147339	127930
AO040716.D	C1704014-005A 90x		92	26112	119025	102419
AO040719.D	C1704014-011A 10x		93	21591	93800	84591
AO040721.D	C1704014-014A 9x		95	20532	89989	83171
AO040722.D	C1704014-014A 90x		89	20827	88198	80228
AO040723.D	C1704014-015A 10x		94	20780	88679	79938
AO040725.D	ALCS1UGD-040717		97	19313	89890	78098

t - fails 24hr time check \* - fails criteria

Created: Thu May 04 12:02:38 2017 MSD #1/



Date: 04-May-17

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.  
Work Order: C1704014  
Project: 691 St Paul Street

TestCode: 0.25CT-TCE-VC

Sample ID	ALCS1UG-040617	Sample Type: LCS	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12114					
Client ID: ZZZZZ	Batch ID: R12114	TestNo: TO-15	Analysis Date: 4/6/2017	SeqNo: 141660							
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	0.9700	0.15	1	0	97.0	70	130				
Chloroethane	1.090	0.15	1	0	109	70	130				
cis-1,2-Dichloroethene	0.9700	0.15	1	0	97.0	70	130				
trans-1,2-Dichloroethene	0.9500	0.15	1	0	95.0	70	130				
Trichloroethene	0.9700	0.040	1	0	97.0	70	130				
Vinyl chloride	0.9700	0.040	1	0	97.0	70	130				

Sample ID	ALCS1UG-040717	SampleType: LCS	Batch ID: R12118	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12118				
Client ID: ZZZZZ				TestNo: TO-15		Analysis Date: 4/7/2017	SeqNo: 141731				
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	0.9900	0.15	1	0	99.0	70	130				
Chloroethane	1.180	0.15	1	0	118	70	130				
cis-1,2-Dichloroethene	1.000	0.15	1	0	100	70	130				
trans-1,2-Dichloroethene	0.9800	0.15	1	0	98.0	70	130				
Trichloroethene	1.060	0.040	1	0	106	70	130				
Vinyl chloride	1.150	0.040	1	0	115	70	130				

Sample ID	ALCS1UGD-040617	Sample Type:	LCSD	TestCode:	0.25CT-TCE-	Units:	ppbV	Prep Date:	RunNo:	12114		
Client ID:	ZZZZZ	Batch ID:	R12114	TestNo:	TO-15	Analysis Date:				4/6/2017	SeqNo:	141661
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual	
1,1-Dichloroethene	0.9900	0.15	1	0	99.0	70	130	0.97	2.04	30		
Chloroethane	1.000	0.15	1	0	100	70	130	1.09	8.61	30		

Qualifiers: - Results reported are not blank corrected  
J Analyte detected below quantitation limit  
S Spike Recovery outside accepted recovery limits  
E Estimated Value above quantitation range  
ND Not Detected at the Limit of Detection  
H Holding times for preparation or analysis exceeded  
R RPD outside accepted recovery limits

CLIENT: LaBella Associates, P.C.  
 Work Order: C1704014  
 Project: 691 St Paul Street

TestCode: 0.25CT-TCE-VC

Sample ID	ALCS1UGD-040617	Sample Type	LCSD	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12114				
Client ID:	ZZZZZ	Batch ID:	R12114	TestNo: TO-15		Analysis Date: 4/6/2017	SeqNo: 141661				
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
cis-1,2-Dichloroethene	0.9600	0.15	1	0	96.0	70	130	0.97	1.04	30	
trans-1,2-Dichloroethene	0.9800	0.15	1	0	98.0	70	130	0.95	3.11	30	
Trichloroethene	1.000	0.040	1	0	100	70	130	0.97	3.05	30	
Vinyl chloride	1.060	0.040	1	0	106	70	130	0.97	8.87	30	

Sample ID	ALCS1UGD-040717	SampleType:	LCSD	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12118				
Client ID: ZZZZ	Batch ID: R12118	TestNo: TO-15	Analysis Date: 4/8/2017	SeqNo: 141732							
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	1.080	0.15	1	0	108	70	130	0.99	8.70	30	
Chloroethane	1.180	0.15	1	0	118	70	130	1.18	0	30	
cis-1,2-Dichloroethene	1.120	0.15	1	0	112	70	130	1	11.3	30	
trans-1,2-Dichloroethene	1.070	0.15	1	0	107	70	130	0.98	8.78	30	
Trichloroethene	1.100	0.040	1	0	110	70	130	1.06	3.70	30	
Vinyl chloride	1.240	0.040	1	0	124	70	130	1.15	7.53	30	

Qualifiers: . Results reported are not blank corrected E Estimated Value above quantitation range H Holding times for preparation or analysis exceeded  
 J Analyte detected below quantitation limit ND Not Detected at the Limit of Detection R RPD outside accepted recovery limits  
 S Spike Recovery outside accepted recovery limits





Date: 04-May-17

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.

Work Order: C1704014

Project: 691 St Paul Street

TestCode: 0.25CT-TCE-VC

Sample ID	C1704014-001A MS	MS	SampleType:	MS	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12114			
Client ID:	691-B19-IAQ-1	R12114	Batch ID:	R12114	TestNo: TO-15		Analysis Date: 4/6/2017	SeqNo: 141726			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	0.9100	0.15	1	0	91.0	70	130				
Chloroethane	1.040	0.15	1	0	104	70	130				
cis-1,2-Dichloroethene	1.230	0.15	1	0.33	90.0	70	130				
trans-1,2-Dichloroethene	0.9500	0.15	1	0	95.0	70	130				
Trichloroethene	1.390	0.040	1	0.4	99.0	70	130				
Vinyl chloride	1.030	0.040	1	0	103	70	130				

Sample ID	C1704014-001A MS	MSD	SampleType:	MSD	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12114			
Client ID:	691-B19-IAQ-1	R12114	Batch ID:	R12114	TestNo: TO-15		Analysis Date: 4/6/2017	SeqNo: 141727			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	1.000	0.15	1	0	100	70	130	0.91	9.42	30	
Chloroethane	1.040	0.15	1	0	104	70	130	1.04	0	30	
cis-1,2-Dichloroethene	1.320	0.15	1	0.33	99.0	70	130	1.23	7.06	30	
trans-1,2-Dichloroethene	0.9700	0.15	1	0	97.0	70	130	0.95	2.08	30	
Trichloroethene	1.370	0.040	1	0.4	97.0	70	130	1.39	1.45	30	
Vinyl chloride	1.030	0.040	1	0	103	70	130	1.03	0	30	

Qualifiers: J Results reported are not blank corrected  
S Analyte detected below quantitation limit  
S Spike Recovery outside accepted recovery limits

E Estimated Value above quantitation range  
ND Not Detected at the Limit of Detection

H Holding times for preparation or analysis exceeded  
R RPD outside accepted recovery limits



Date: 04-May-17

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.  
Work Order: C1704014  
Project: 691 St Paul Street

TestCode: 0.25CT-TCE-VC

Sample ID	AMB1UG-040617	Sample Type:	MBLK	TestCode:	0.25CT-TCE-	Units:	ppbv	Prep Date:	RunNo:	12114	
Client ID:	ZZZZZ	Batch ID:	R12114	TestNo:	TO-15			Analysis Date:	SeqNo:	141659	
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,1-Dichloroethene	< 0.15	0.15									
Chloroethane	< 0.15	0.15									
cis-1,2-Dichloroethene	< 0.15	0.15									
trans-1,2-Dichloroethene	< 0.15	0.15									
Trichloroethene	< 0.040	0.040									
Vinyl chloride	< 0.040	0.040									

Sample ID	AMB1UG-040717	Sample Type:	MBLK	TestCode:	0.25CT-TCE-	Units:	ppbV	Prep Date:	RunNo:	12118	
Client ID:	ZZZZZ	Batch ID:	R12118	TestNo:	TO-15			Analysis Date:	SeqNo:	141730	
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,1-Dichloroethene	< 0.15	0.15									
Chloroethane	< 0.15	0.15									
cis-1,2-Dichloroethene	< 0.15	0.15									
trans-1,2-Dichloroethene	< 0.15	0.15									
Trichloroethene	< 0.040	0.040									
Vinyl chloride	< 0.040	0.040									

Qualifiers: J Results reported are not blank corrected  
S Analyte detected below quantitation limit  
S Spike Recovery outside accepted recovery limits  
E Estimated Value above quantitation range  
ND Not Detected at the Limit of Detection  
H Holding times for preparation or analysis exceeded  
R RPD outside accepted recovery limits

Name	Amount	IDL#1	IDL#2	IDL#3	IDL#4	IDL#5	IDL#6	IDL#7	Average	StdDev	%Rec	IDL
Propylene	0.15	0.16	0.15	0.16	0.14	0.16	0.14	0.16	0.153	0.010	98.1	0.030
Freon 12	0.15	0.18	0.17	0.17	0.17	0.18	0.17	0.17	0.173	0.005	86.8	0.015
Chloromethane	0.15	0.19	0.18	0.16	0.18	0.18	0.2	0.17	0.180	0.013	83.3	0.041
Freon 114	0.15	0.18	0.17	0.17	0.17	0.18	0.17	0.18	0.174	0.005	86.1	0.017
Vinyl Chloride	0.15	0.17	0.16	0.16	0.15	0.16	0.15	0.15	0.157	0.008	95.5	0.024
Butane	0.15	0.18	0.16	0.17	0.18	0.18	0.19	0.19	0.179	0.011	84.0	0.034
1,3-butadiene	0.15	0.21	0.2	0.2	0.22	0.17	0.18	0.23	0.201	0.021	74.5	0.066
Bromomethane	0.15	0.18	0.2	0.21	0.18	0.22	0.16	0.21	0.194	0.021	77.2	0.068
Chloroethane	0.15	0.19	0.19	0.16	0.19	0.19	0.18	0.19	0.184	0.011	81.4	0.036
Ethanol	0.15	0.16	0.16	0.18	0.17	0.19	0.18	0.19	0.176	0.013	85.4	0.040
Acrolein	0.15	0.22	0.17	0.19	0.16	0.18	0.21	0.17	0.186	0.022	80.8	0.070
Vinyl Bromide	0.15	0.17	0.15	0.16	0.16	0.17	0.17	0.17	0.164	0.008	91.3	0.025
Freon 11	0.15	0.18	0.17	0.17	0.18	0.19	0.17	0.18	0.177	0.008	84.7	0.024
Acetone	0.15	0.2	0.17	0.18	0.15	0.15	0.18	0.14	0.167	0.021	89.7	0.067
Pentane	0.15	0.18	0.17	0.18	0.16	0.17	0.2	0.16	0.174	0.014	86.1	0.044
Isopropyl alcohol	0.15	0.22	0.2	0.19	0.2	0.19	0.21	0.19	0.200	0.012	75.0	0.036
1,1-dichloroethene	0.15	0.2	0.17	0.19	0.19	0.19	0.18	0.18	0.186	0.010	80.8	0.031
Freon 113	0.15	0.17	0.16	0.18	0.18	0.18	0.17	0.17	0.173	0.008	86.8	0.024
t-Butyl alcohol	0.15	0.21	0.2	0.2	0.21	0.2	0.2	0.18	0.200	0.010	75.0	0.031
Methylene chloride	0.15	0.2	0.18	0.19	0.18	0.2	0.19	0.17	0.187	0.011	80.2	0.035
Allyl chloride	0.15	0.18	0.17	0.16	0.18	0.18	0.2	0.18	0.179	0.012	84.0	0.038
Carbon disulfide	0.15	0.2	0.17	0.19	0.19	0.2	0.18	0.19	0.189	0.011	79.5	0.034
trans-1,2-dichloroethene	0.15	0.15	0.14	0.14	0.14	0.16	0.14	0.15	0.146	0.008	102.9	0.025
methyl tert-butyl ether	0.15	0.14	0.14	0.14	0.13	0.15	0.14	0.13	0.139	0.007	108.2	0.022
1,1-dichloroethane	0.15	0.17	0.15	0.16	0.15	0.17	0.16	0.16	0.160	0.008	93.8	0.026
Vinyl acetate	0.15	0.14	0.13	0.14	0.13	0.13	0.13	0.12	0.131	0.007	114.1	0.022
Methyl Ethyl Ketone	0.15	0.17	0.17	0.16	0.16	0.15	0.13	0.12	0.151	0.020	99.1	0.061
cis-1,2-dichloroethene	0.15	0.15	0.14	0.16	0.15	0.16	0.15	0.14	0.150	0.008	100.0	0.026
Hexane	0.15	0.12	0.14	0.13	0.13	0.13	0.12	0.12	0.127	0.008	118.0	0.024
Ethyl acetate	0.15	0.16	0.17	0.14	0.15	0.14	0.16	0.13	0.150	0.014	100.0	0.044
Chloroform	0.15	0.16	0.16	0.16	0.16	0.17	0.16	0.17	0.163	0.005	92.1	0.015
Tetrahydrofuran	0.15	0.15	0.13	0.15	0.15	0.15	0.15	0.14	0.146	0.008	102.9	0.025
1,2-dichloroethane	0.15	0.16	0.15	0.16	0.16	0.17	0.16	0.17	0.161	0.007	92.9	0.022
1,1,1-trichloroethane	0.15	0.17	0.16	0.17	0.17	0.16	0.17	0.17	0.167	0.005	89.7	0.015
Cyclohexane	0.15	0.14	0.14	0.14	0.15	0.15	0.14	0.14	0.143	0.005	105.0	0.015
Carbon tetrachloride	0.15	0.13	0.15	0.15	0.15	0.15	0.15	0.16	0.149	0.009	101.0	0.028
Benzene	0.15	0.15	0.16	0.16	0.15	0.16	0.16	0.16	0.157	0.005	95.5	0.015
Methyl methacrylate	0.15	0.15	0.15	0.14	0.14	0.14	0.15	0.11	0.140	0.014	107.1	0.044
1,4-dioxane	0.15	0.18	0.18	0.19	0.18	0.15	0.17	0.12	0.167	0.024	89.7	0.076

Confidential

1/8/2016

Name	Amount	IDL#1	IDL#2	IDL#3	IDL#4	IDL#5	IDL#6	IDL#7	Average	StdDev	%Rec	IDL
2,2,4-trimethylpentane	0.15	0.15	0.15	0.15	0.16	0.14	0.16	0.15	0.151	0.007	99.1	0.022
Heptane	0.15	0.12	0.13	0.13	0.12	0.13	0.13	0.13	0.127	0.005	118.0	0.015
Trichloroethene	0.15	0.14	0.15	0.14	0.15	0.15	0.14	0.15	0.146	0.005	102.9	0.017
1,2-dichloropropane	0.15	0.16	0.17	0.17	0.16	0.17	0.16	0.16	0.164	0.005	91.3	0.017
Bromodichloromethane	0.15	0.16	0.16	0.16	0.15	0.16	0.17	0.16	0.160	0.006	93.8	0.018
cis-1,3-dichloropropene	0.15	0.13	0.13	0.14	0.14	0.13	0.13	0.13	0.133	0.005	112.9	0.015
trans-1,3-dichloropropene	0.15	0.16	0.13	0.13	0.14	0.14	0.14	0.16	0.143	0.013	105.0	0.039
1,1,2-trichloroethane	0.15	0.16	0.15	0.16	0.15	0.16	0.18	0.17	0.161	0.011	92.9	0.034
Toluene	0.15	0.14	0.14	0.14	0.13	0.16	0.14	0.15	0.143	0.010	105.0	0.030
Methyl Isobutyl Ketone	0.15	0.18	0.18	0.18	0.18	0.16	0.18	0.15	0.173	0.013	86.8	0.039
Dibromochloromethane	0.15	0.16	0.16	0.17	0.18	0.16	0.17	0.18	0.169	0.009	89.0	0.028
Methyl Butyl Ketone	0.15	0.17	0.16	0.18	0.17	0.16	0.17	0.14	0.164	0.013	91.3	0.040
1,2-dibromoethane	0.15	0.16	0.17	0.16	0.16	0.16	0.16	0.17	0.163	0.005	92.1	0.015
Tetrachloroethylene	0.15	0.16	0.17	0.16	0.16	0.16	0.17	0.17	0.164	0.005	91.3	0.017
Chlorobenzene	0.15	0.16	0.16	0.16	0.17	0.15	0.17	0.17	0.163	0.008	92.1	0.024
1,1,1,2-tetrachloroethane	0.15	0.17	0.17	0.17	0.18	0.16	0.18	0.17	0.171	0.007	87.5	0.022
Ethylbenzene	0.15	0.13	0.14	0.14	0.14	0.12	0.14	0.13	0.134	0.008	111.7	0.025
m&p-xylene	0.3	0.25	0.25	0.25	0.23	0.25	0.25	0.25	0.247	0.008	121.4	0.024
Nonane	0.15	0.11	0.11	0.11	0.11	0.1	0.1	0.11	0.107	0.005	140.0	0.015
Styrene	0.15	0.12	0.13	0.13	0.11	0.12	0.13	0.12	0.123	0.008	122.1	0.024
Bromoform	0.15	0.15	0.15	0.16	0.15	0.15	0.17	0.16	0.156	0.008	96.3	0.025
o-xylene	0.15	0.11	0.12	0.12	0.14	0.14	0.12	0.11	0.123	0.013	122.1	0.039
Cumene	0.15	0.12	0.13	0.13	0.12	0.13	0.13	0.13	0.127	0.005	118.0	0.015
Bromofluorobenzene	1	0.88	0.9	0.9	0.87	0.89	0.89	0.9	0.890	0.012	112.4	0.036
1,1,2,2-tetrachloroethane	0.15	0.16	0.16	0.17	0.16	0.17	0.17	0.16	0.164	0.005	91.3	0.017
Propylbenzene	0.15	0.13	0.12	0.13	0.13	0.11	0.13	0.11	0.123	0.010	122.1	0.030
2-Chlorotoluene	0.15	0.13	0.13	0.13	0.14	0.13	0.12	0.13	0.130	0.006	115.4	0.018
4-ethyltoluene	0.15	0.11	0.12	0.12	0.12	0.13	0.13	0.11	0.120	0.008	125.0	0.026
1,3,5-trimethylbenzene	0.15	0.12	0.13	0.14	0.12	0.13	0.13	0.13	0.129	0.007	116.7	0.022
1,2,4-trimethylbenzene	0.15	0.12	0.13	0.12	0.12	0.13	0.12	0.12	0.123	0.005	122.1	0.015
1,3-dichlorobenzene	0.15	0.14	0.14	0.14	0.13	0.14	0.13	0.14	0.137	0.005	109.4	0.015
benzyl chloride	0.15	0.13	0.16	0.13	0.15	0.13	0.15	0.16	0.144	0.014	104.0	0.044
1,4-dichlorobenzene	0.15	0.13	0.11	0.12	0.12	0.12	0.12	0.13	0.121	0.007	123.5	0.022
1,2,3-trimethylbenzene	0.15	0.12	0.11	0.12	0.12	0.12	0.11	0.11	0.116	0.005	129.6	0.017
1,2-dichlorobenzene	0.15	0.13	0.14	0.14	0.14	0.14	0.14	0.13	0.137	0.005	109.4	0.015
1,2,4-trichlorobenzene	0.15	0.1	0.11	0.1	0.11	0.11	0.12	0.1	0.107	0.008	140.0	0.024
Naphthalene	0.15	0.13	0.13	0.14	0.11	0.12	0.14	0.12	0.127	0.011	118.0	0.035
Hexachloro-1,3-butadiene	0.15	0.16	0.17	0.17	0.17	0.16	0.16	0.16	0.164	0.005	91.3	0.017



Name	Amount	IDL#1	IDL#2	IDL#3	IDL#4	IDL#5	IDL#6	IDL#7	Average	StdDev	%Rec	IDL
Vinyl Chloride	0.1	0.11	0.11	0.09	0.09	0.1	0.09	0.1	0.099	0.009	101.4	0.028
Carbon tetrachloride	0.1	0.1	0.11	0.08	0.09	0.09	0.09	0.09	0.093	0.010	107.7	0.030
Trichloroethene	0.1	0.1	0.1	0.07	0.08	0.08	0.08	0.08	0.084	0.011	118.6	0.036
Tetrachloroethylene	0.1	0.11	0.12	0.09	0.09	0.1	0.09	0.09	0.099	0.012	101.4	0.038
Naphthalene	0.1	0.09	0.08	0.07	0.06	0.06	0.07	0.06	0.070	0.012	142.9	0.036

## GC/MS-Whole Air Calculations

### Relative Response Factor (RRF)

$$RRF = \frac{A_x * C_{is}}{A_{is} * C_x}$$

where:  $A_x$  = area of the characteristic ion for the compound being measured  
 $A_{is}$  = area of the characteristic ion for the specific internal standard of the compound being measured  
 $C_x$  = concentration of the compound being measured (ppbv)  
 $C_{is}$  = concentration of the internal standard (ppbv)

### Percent Relative Standard Deviation (%RSD)

$$\% RSD = \frac{\text{Standard deviation of RRF values} * 100}{\text{mean RRF}}$$

### Percent Difference (%D)

$$\% D = \frac{(RRF_c - \text{mean RRF}_i) * 100}{\text{mean RRF}_i}$$

where:  $RRF_c$  = relative response factor from the continuing calibration  
 $\text{mean RRF}_i$  = mean relative response factor from the initial calibration

### Sample Calculations

$$\text{ppbv} = \frac{A_x * I_s * D_f}{A_{is} * RRF}$$

where:  $A_x$  = area of the characteristic ion for the compound being measured  
 $A_{is}$  = area of the characteristic ion for the specific internal standard of the compound being measured  
 $I_s$  = Concentration of the internal standard injected (ppbv)  
 $RRF$  = relative response factor for the compound being measured  
 $D_f$  = Dilution factor

**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**SAMPLE DATA**

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-001A

Client Sample ID: 691-B19-1AQ-1  
Tag Number: 1321.1163  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
		FLD				Analyst:
Lab Vacuum In	-4			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
		TO-15				Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/6/2017 9:55:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	4/6/2017 9:55:00 PM
cis-1,2-Dichloroethene	0.33	0.15		ppbV	1	4/6/2017 9:55:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/6/2017 9:55:00 PM
Trichloroethene	0.40	0.040		ppbV	1	4/6/2017 9:55:00 PM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/6/2017 9:55:00 PM
Surr: Bromofluorobenzene	96.0	70-130		%REC	1	4/6/2017 9:55:00 PM

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-001A

Client Sample ID: 691-B19-1AQ-1  
Tag Number: 1321.1163  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/6/2017 9:55:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/6/2017 9:55:00 PM
cis-1,2-Dichloroethene	1.3	0.59		ug/m3	1	4/6/2017 9:55:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/6/2017 9:55:00 PM
Trichloroethene	2.1	0.21		ug/m3	1	4/6/2017 9:55:00 PM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/6/2017 9:55:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

Data File : C:\HPCHEM\1\DATA\AO040621.D

Vial: 21

Acq On : 6 Apr 2017 9:55 pm

Operator: RJP

Sample : C1704014-001A

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 11 14:02:25 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.57	128	25396	1.00	ppb	0.03
35) 1,4-difluorobenzene	11.95	114	113281	1.00	ppb	0.02
50) Chlorobenzene-d5	16.83	117	97086	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.47	95	63494	0.96	ppb	0.02
Spiked Amount	1.000	Range	70 - 130	Recovery	=	96.00%

## Target Compounds

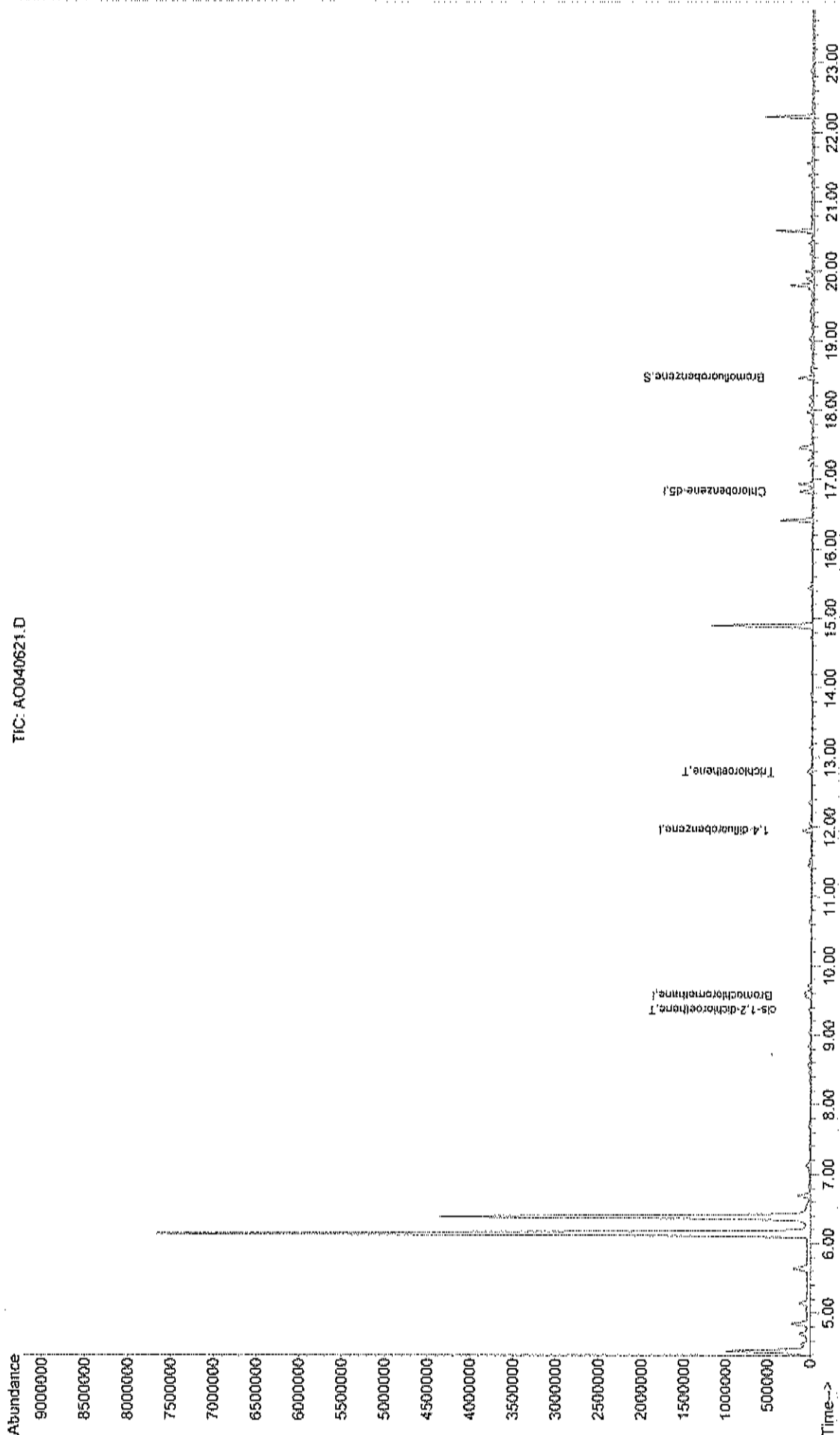
	R.T.	QIon	Response	Conc	Units	Qvalue
29) cis-1,2-dichloroethene	9.37	61	17307	0.33	ppb	96
44) Trichloroethene	12.79	130	20426	0.40	ppb	92

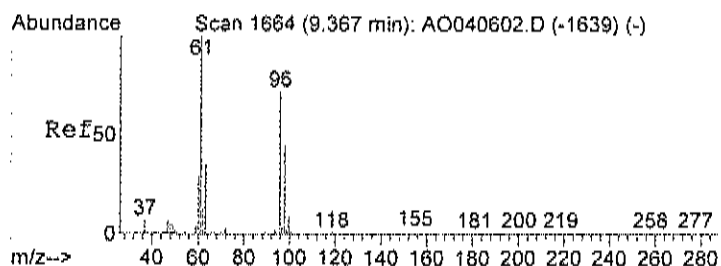
Data File : C:\HPCHEM\1\DATA\AO040621.D  
 Acq On : 6 Apr 2017 9:55 pm  
 Sample : C1704014-001A  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 11 14:03 2017

Vial: 21  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

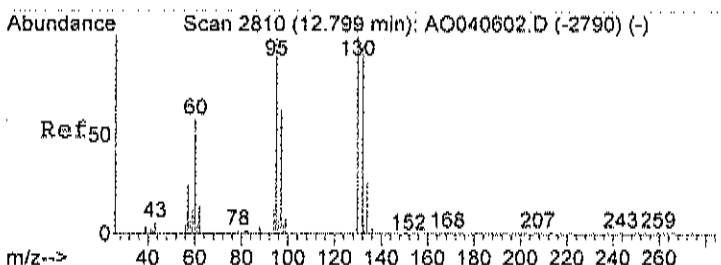
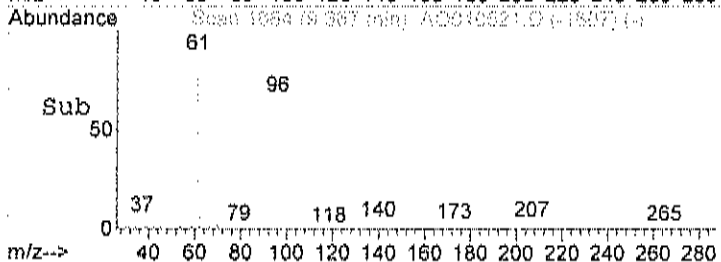
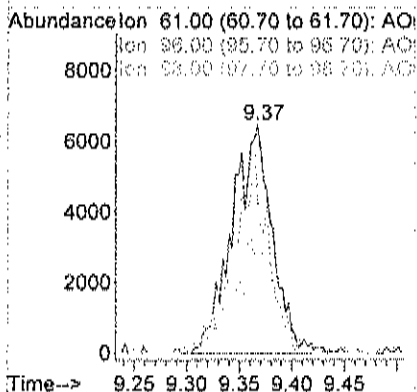
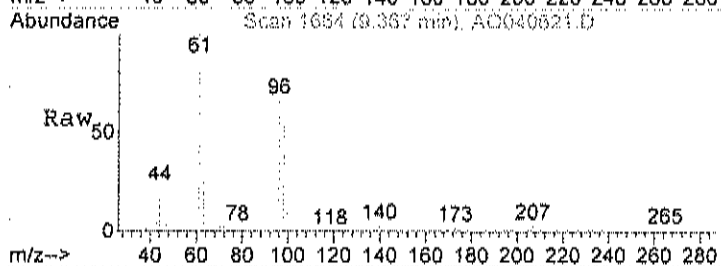
Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration





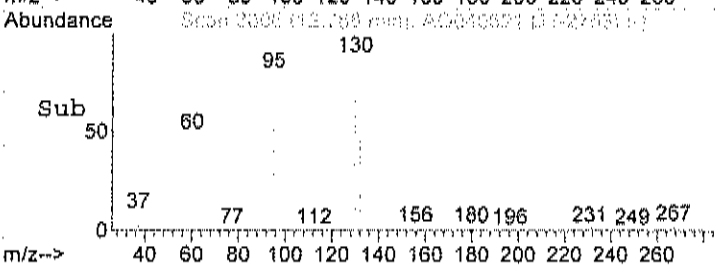
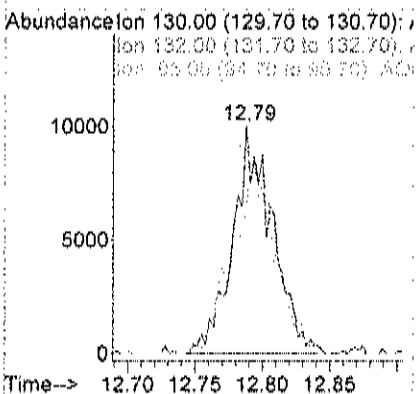
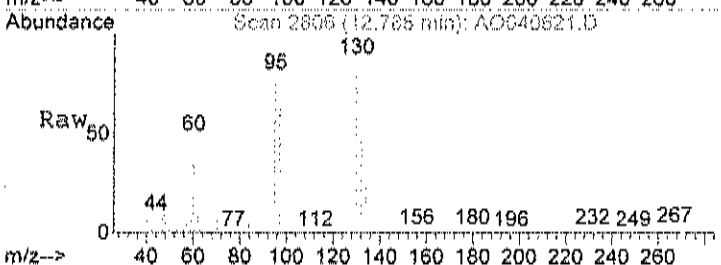
#29  
cis-1,2-dichloroethene  
Concen: 0.33 ppb  
RT: 9.37 min Scan# 1664  
Delta R.T. 0.02 min  
Lab File: AO040621.D  
Acq: 6 Apr 2017 9:55 pm

Tgt Ion:	61	Resp:	17307
Ion	Ratio	Lower	Upper
61	100		
96	74.6	58.1	98.1
98	52.6	29.3	69.3



#44  
Trichloroethene  
Concen: 0.40 ppb  
RT: 12.79 min Scan# 2806  
Delta R.T. 0.01 min  
Lab File: AO040621.D  
Acq: 6 Apr 2017 9:55 pm

Tgt Ion:	130	Resp:	20426
Ion	Ratio	Lower	Upper
130	100		
132	92.8	69.9	109.9
95	108.3	76.3	116.3





# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-002A

Client Sample ID: 691-B19-IAQ-2  
Tag Number: 1188.306  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-9			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 12:09:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 12:09:00 AM
cis-1,2-Dichloroethene	0.29	0.15		ppbV	1	4/7/2017 12:09:00 AM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 12:09:00 AM
Trichloroethene	0.39	0.040		ppbV	1	4/7/2017 12:09:00 AM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 12:09:00 AM
Surr: Bromofluorobenzene	89.0	70-130		%REC	1	4/7/2017 12:09:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-002A

Client Sample ID: 691-B19-IAQ-2  
Tag Number: 1188.306  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 12:09:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 12:09:00 AM
cis-1,2-Dichloroethene	1.1	0.59		ug/m3	1	4/7/2017 12:09:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 12:09:00 AM
Trichloroethene	2.1	0.21		ug/m3	1	4/7/2017 12:09:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 12:09:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

Data File : C:\HPCHEM\1\DATA\AO040624.D  
Acq On : 7 Apr 2017 12:09 am  
Sample : C1704014-002A  
Misc : A331\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Apr 07 07:19:04 2017

Vial: 22  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Apr 03 10:15:59 2017  
Response via : Initial Calibration  
DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	24868	1.00	ppb	0.01
35) 1,4-difluorobenzene	11.94	114	113360	1.00	ppb	0.00
50) Chlorobenzene-d5	16.83	117	102277	1.00	ppb	0.00

System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	62337	0.89	ppb	0.01
Spiked Amount	1.000	Range	70 - 130	Recovery	=	89.00%

Target Compounds

						Qvalue
29) cis-1,2-dichloroethene	9.35	61	15114	0.29	ppb	97
44) Trichloroethene	12.79	130	19716	0.39	ppb	95

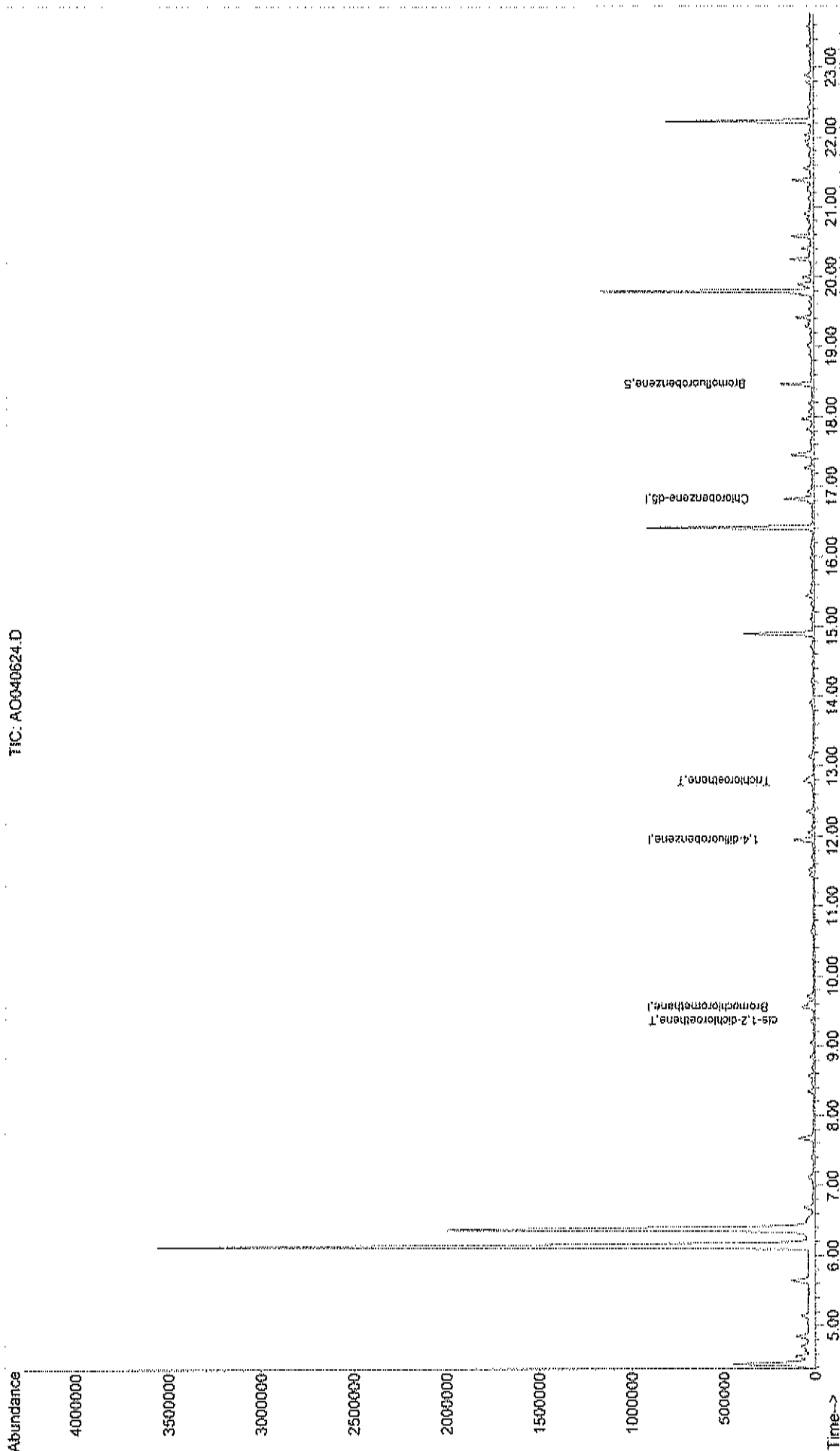
Data File : C:\HPCHEM\1\DATA\A0040624.D  
Acq On : 7 Apr 2017 12:09 am  
Sample : C1704014-002A  
Misc : A331\_LUG  
MMS Integration Params: RTEINT.P  
Quant Time: Apr 7 13:24 2017

Vial: 22  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

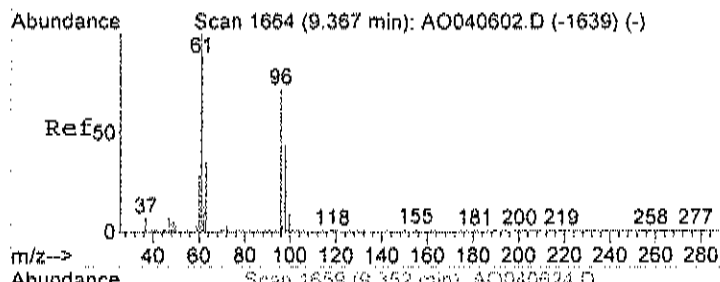
Quant Results File: A331 1UG.RES

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Method      : C:\HPCHEM\1\METHODS\A331_1UG.M (RTE Integrator)
Title       : TO-15 VOA Standards for 5 point calibration
Last Update : Thu May 04 11:27:28 2017
Response via : Initial Calibration
```

FIG: A0340624.D

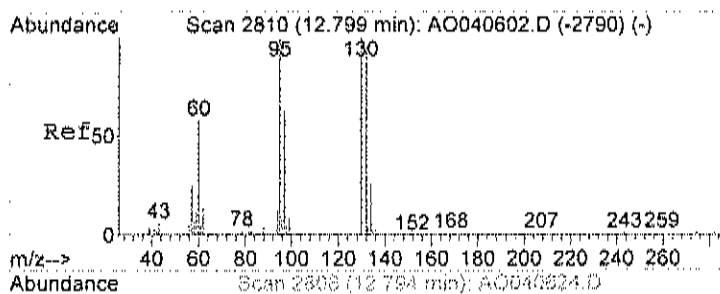
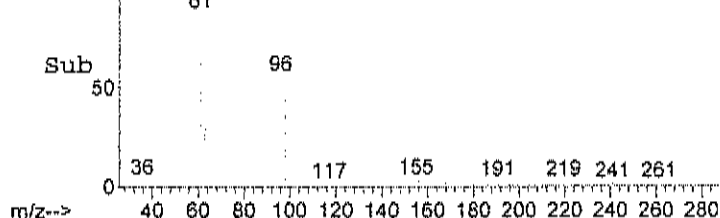
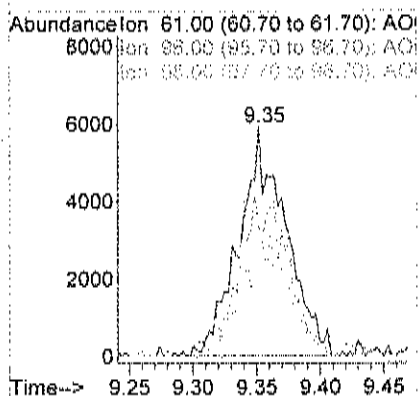
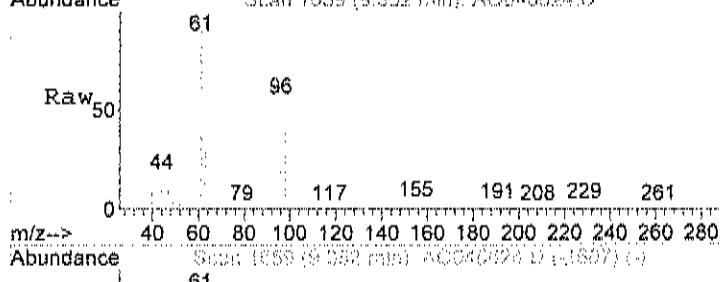






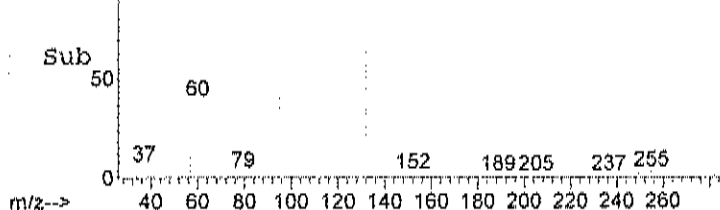
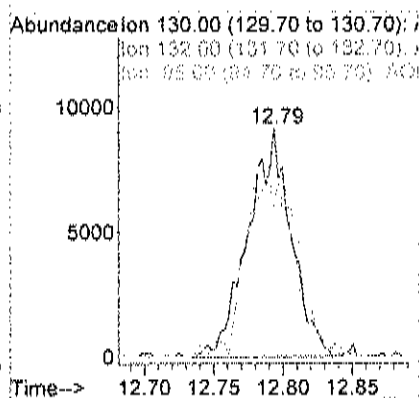
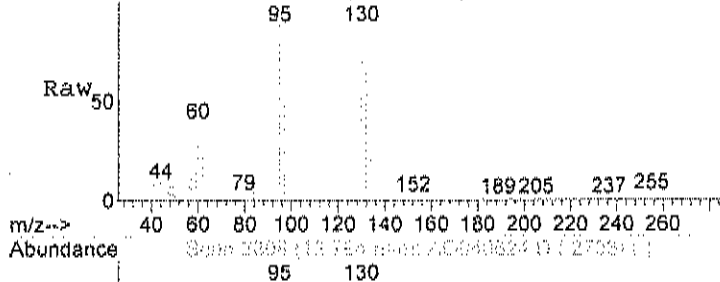
#29  
cis-1,2-dichloroethene  
Concen: 0.29 ppb  
RT: 9.35 min Scan# 1659  
Delta R.T. 0.01 min  
Lab File: AO040624.D  
Acq: 7 Apr 2017 12:09 am

Tgt Ion	Ratio	Lower	Upper
61	100		
96	75.2	58.1	98.1
98	50.4	29.3	69.3



#44  
Trichloroethene  
Concen: 0.39 ppb  
RT: 12.79 min Scan# 2808  
Delta R.T. 0.01 min  
Lab File: AO040624.D  
Acq: 7 Apr 2017 12:09 am

Tgt Ion	Ratio	Lower	Upper
130	100		
132	89.9	69.9	109.9
95	105.7	76.3	116.3



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-003A

Client Sample ID: 691-B19-SV1  
Tag Number: 362.446  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-9			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 3:23:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 3:23:00 PM
cis-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 3:23:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 3:23:00 PM
Trichloroethene	4.8	0.40		ppbV	10	4/7/2017 8:07:00 PM
Vinyl chloride	0.080	0.040		ppbV	1	4/7/2017 3:23:00 PM
Surr: Bromofluorobenzene	109	70-130		%REC	1	4/7/2017 3:23:00 PM

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte, Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-003A

Client Sample ID: 691-B19-SV1  
Tag Number: 362.446  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 3:23:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 3:23:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 3:23:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 3:23:00 PM
Trichloroethene	26	2.1		ug/m3	10	4/7/2017 8:07:00 PM
Vinyl chloride	0.20	0.10		ug/m3	1	4/7/2017 3:23:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

Page 3 of 15

Data File : C:\HPCHEM\1\DATA\AO040708.D

Vial: 8

Acq On : 7 Apr 2017 3:23 pm

Operator: RJP

Sample : C1704014-003A

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 07 15:59:23 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.55	128	20468	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.94	114	96070	1.00	ppb	0.00
50) Chlorobenzene-d5	16.82	117	92232	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	68296	1.09	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	109.00%

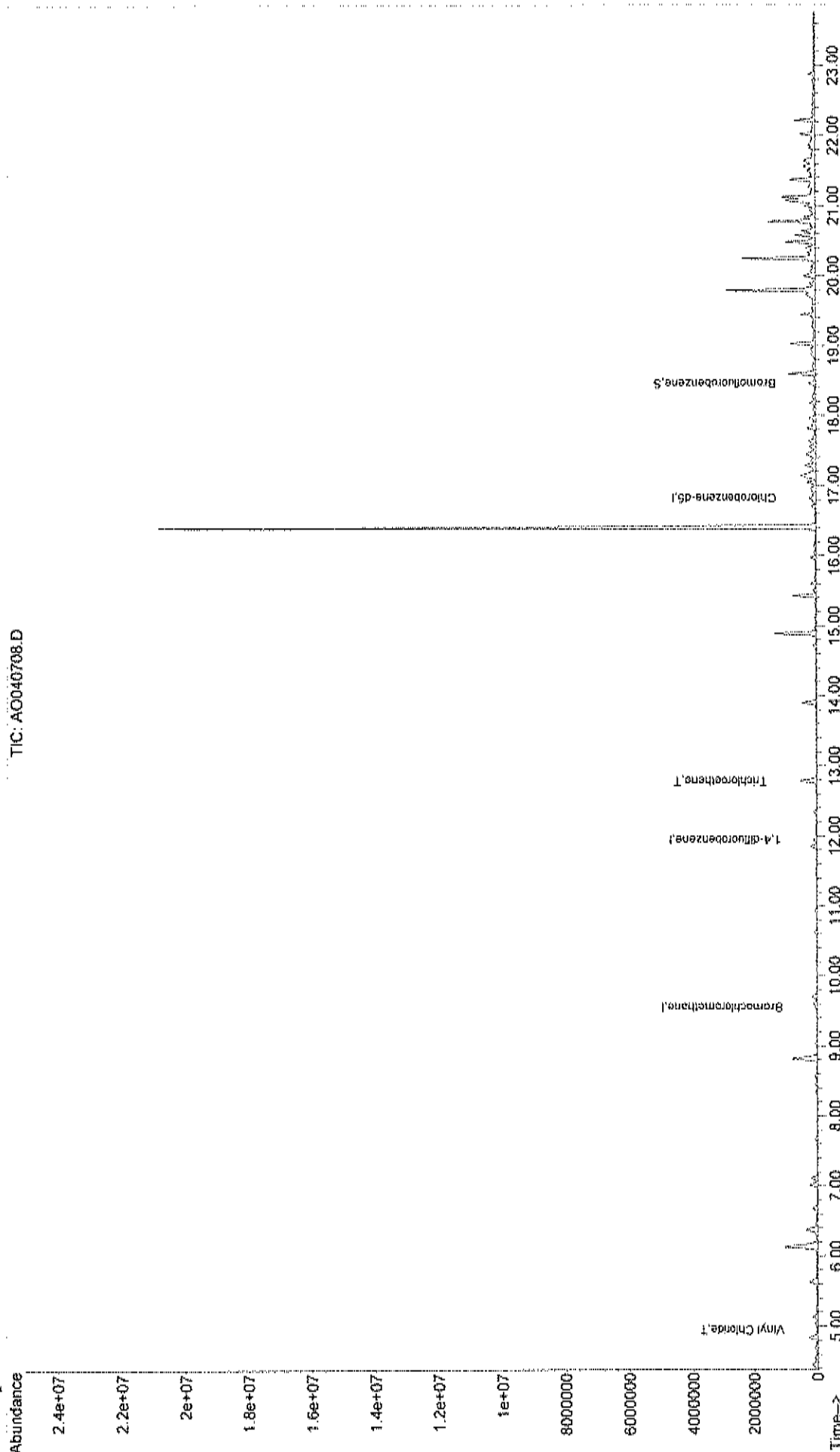
## Target Compounds

						Qvalue
6) Vinyl Chloride	4.95	62	3018	0.08	ppb	79
44) Trichloroethene	12.78	130	203569	4.71	ppb	91

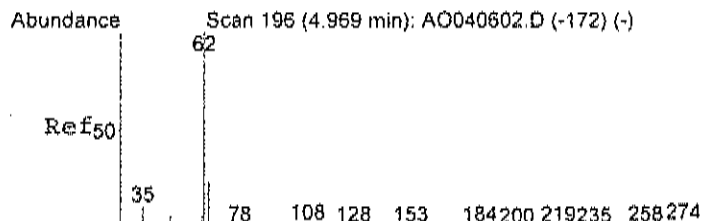
Data File : C:\HPCHEM\1\DATA\AO040708.D  
 Acq On : 7 Apr 2017 3:23 pm  
 Sample : C1704014-003A  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 11 14:07 2017

Vial: 8  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00  
 Quant Results File: A331\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration

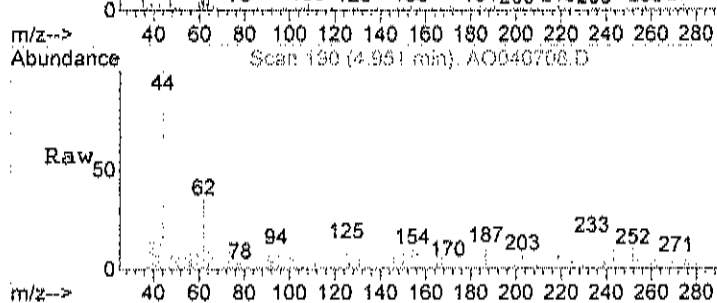




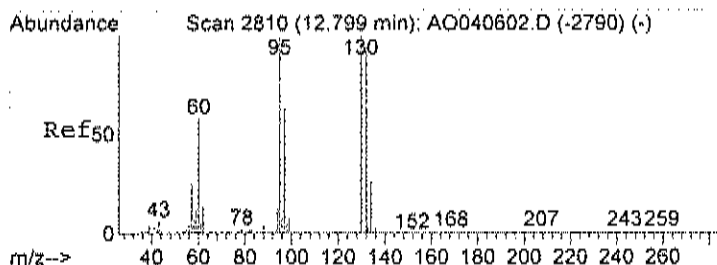
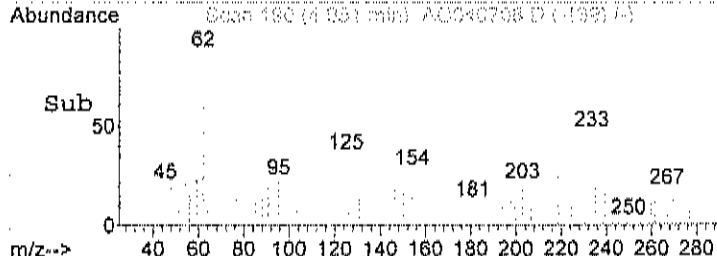
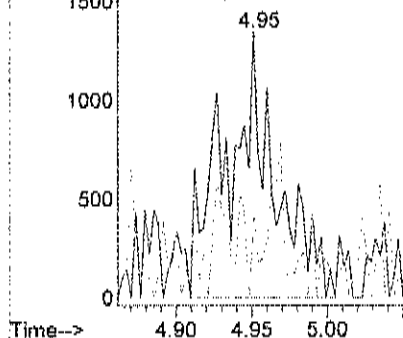


#6  
 Vinyl Chloride  
 Concen: 0.08 ppb  
 RT: 4.95 min Scan# 190  
 Delta R.T. 0.00 min  
 Lab File: AO040708.D  
 Acq: 7 Apr 2017 3:23 pm

Tgt Ion: 62 Resp: 3018  
 Ion Ratio Lower Upper  
 62 100  
 64 17.7 0.0 58.6

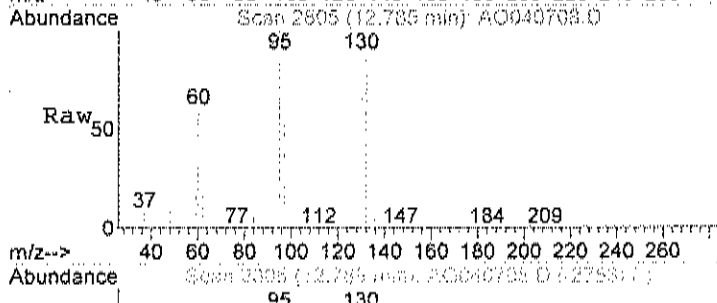


Abundance Ion 62.00 (61.70 to 62.70): AO  
 Ion 64.00 (63.70 to 64.70): AO

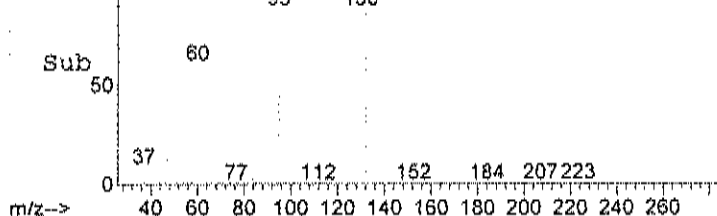
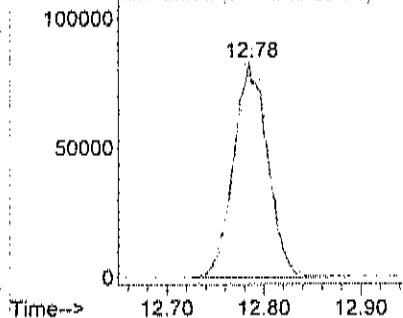


#44  
 Trichloroethene  
 Concen: 4.71 ppb  
 RT: 12.78 min Scan# 2805  
 Delta R.T. 0.01 min  
 Lab File: AO040708.D  
 Acq: 7 Apr 2017 3:23 pm

Tgt Ion: 130 Resp: 203569  
 Ion Ratio Lower Upper  
 130 100  
 132 96.4 69.9 109.9  
 95 107.5 76.3 116.3



Abundance Ion 130.00 (129.70 to 130.70):  
 Ion 132.00 (131.70 to 132.70):  
 Ion 95.00 (94.70 to 95.70): AO



Data File : C:\HPCHEM\1\DATA\AO040714.D  
 Acq On : 7 Apr 2017 8:07 pm  
 Sample : C1704014-003A 10x  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 08 08:59:49 2017

Vial: 14  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Mon Apr 03 10:15:59 2017  
 Response via : Initial Calibration  
 DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	31326	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	147339	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	127930	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.47	95	83018	0.95	ppb	0.02
Spiked Amount	1.000	Range	70 - 130	Recovery	=	95.00%

## Target Compounds

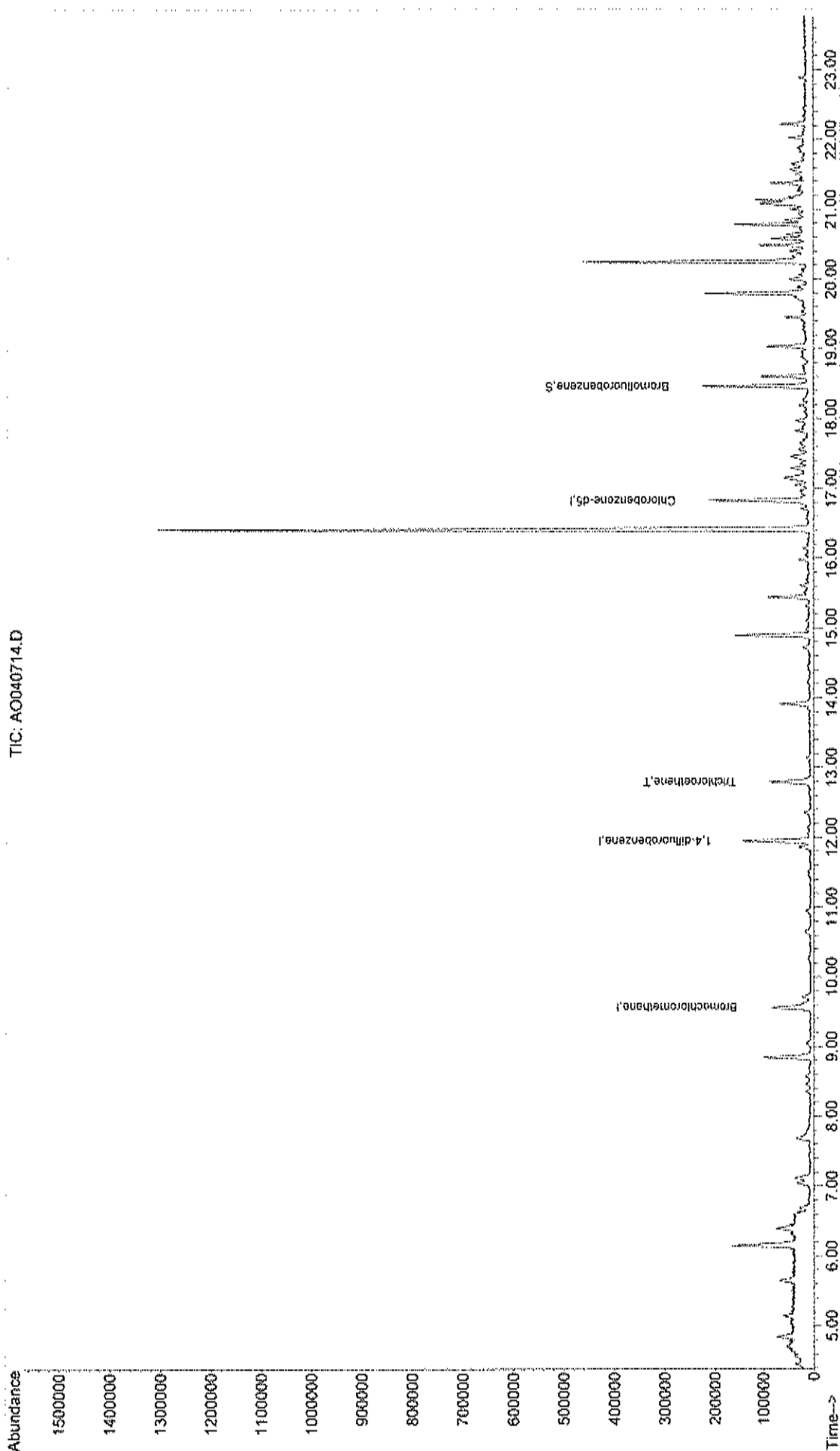
	R.T.	QIon	Response	Conc	Units	Qvalue
44) Trichloroethene	12.79	130	31747	0.48	ppb	92

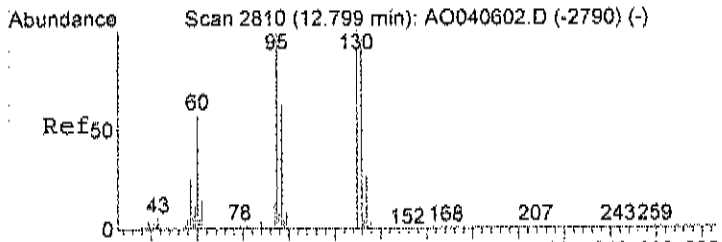
Data File : C:\HPCHEM\1\DATA\A0040714.D  
 Acq On : 7 Apr 2017 8:07 pm  
 Sample : C1704014-003A 10x  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 11 14:57 2017

Vial: 14  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

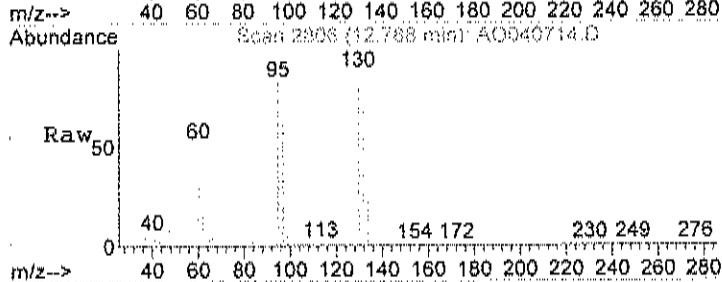
Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration





#44  
 Trichloroethene  
 Concen: 0.48 ppb  
 RT: 12.79 min Scan# 2806  
 Delta R.T. 0.01 min  
 Lab File: AO040714.D  
 Acq: 7 Apr 2017 8:07 pm

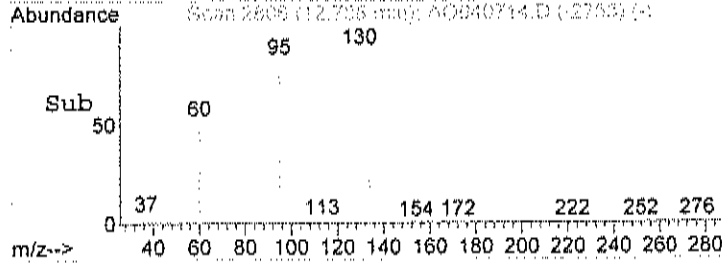
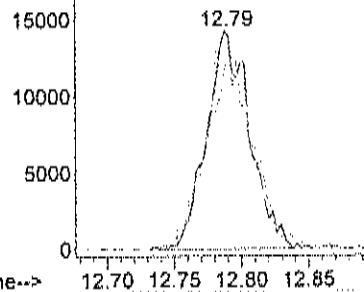
Tgt Ion	Ratio	Lower	Upper
130	100		
132	95.8	69.9	109.9
95	105.6	76.3	116.3



Abundance Ion 130.00 (129.70 to 130.70):

20000 Ion 132.00 (131.70 to 132.70):

Ion 95.00 (94.70 to 95.70): AO



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.

Client Sample ID: 691-B15-IAQ

Lab Order: C1704014

Tag Number: 87.299

Project: 691 St Paul Street

Collection Date: 4/1/2017

Lab ID: C1704014-004A

Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-3			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 12:50:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 12:50:00 AM
cis-1,2-Dichloroethene	0.40	0.15		ppbV	1	4/7/2017 12:50:00 AM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 12:50:00 AM
Trichloroethene	0.41	0.040		ppbV	1	4/7/2017 12:50:00 AM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 12:50:00 AM
Surr: Bromofluorobenzene	92.0	70-130		%REC	1	4/7/2017 12:50:00 AM

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte, Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

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# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-004A

Client Sample ID: 691-B15-1AQ  
Tag Number: 87.299  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 12:50:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 12:50:00 AM
cis-1,2-Dichloroethene	1.6	0.59		ug/m3	1	4/7/2017 12:50:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 12:50:00 AM
Trichloroethene	2.2	0.21		ug/m3	1	4/7/2017 12:50:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 12:50:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

Data File : C:\HPCHEM\1\DATA\AO040625.D  
 Acq On : 7 Apr 2017 12:50 am  
 Sample : C1704014-004A  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 07 07:19:05 2017

Vial: 23  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Mon Apr 03 10:15:59 2017  
 Response via : Initial Calibration  
 DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	25596	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	114646	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	102450	1.00	ppb	0.01

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	63996	0.92	ppb	0.01
Spiked Amount	1.000	Range	70 - 130	Recovery	=	92.00%

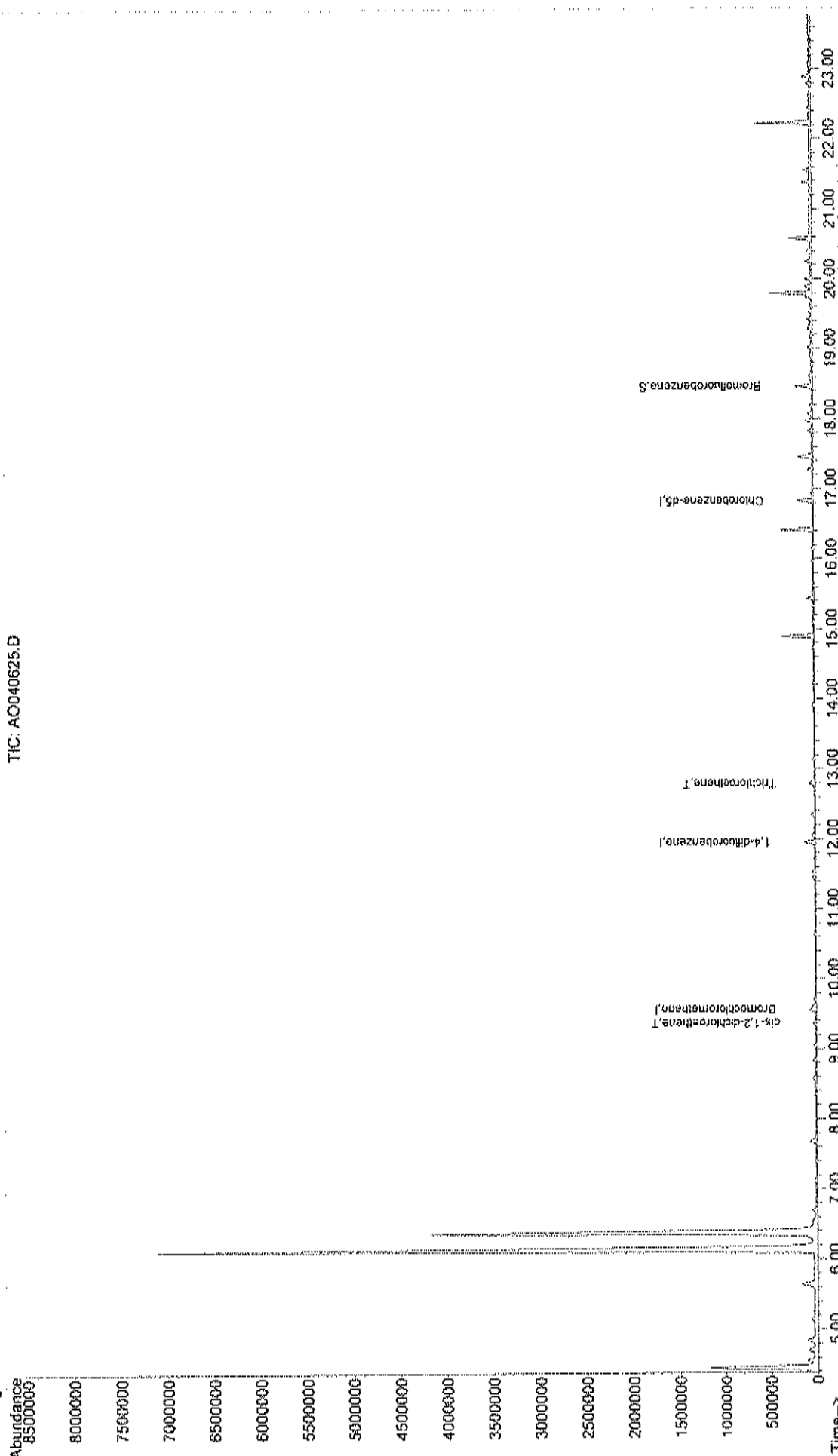
## Target Compounds

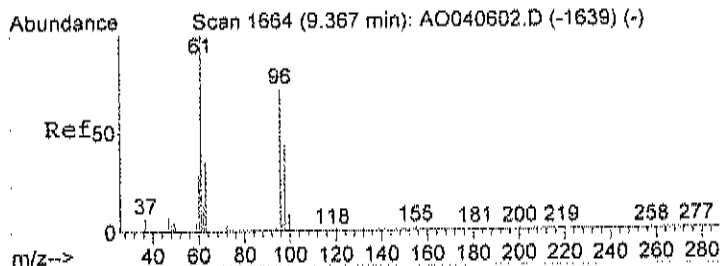
						Qvalue
29) cis-1,2-dichloroethene	9.37	61	21474	0.40	ppb	91
44) Trichloroethene	12.79	130	21262	0.41	ppb	95

Data File : C:\HPCHEM\1\DATA\A0040625.D  
 Acq On : 7 Apr 2017 12:50 am  
 Sample : C1704014-004A  
 Misc : A331\_LUG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 7 13:25 2017

Vial: 23  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00  
 Quant Results File: A331\_LUG.RES

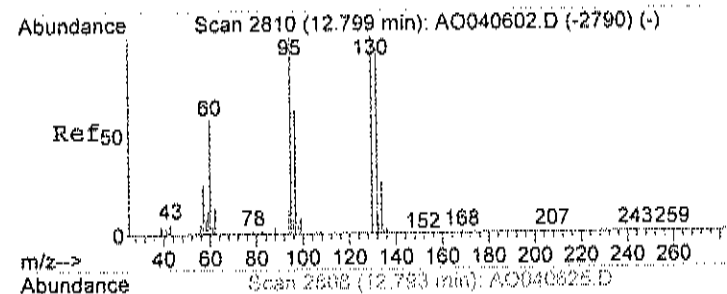
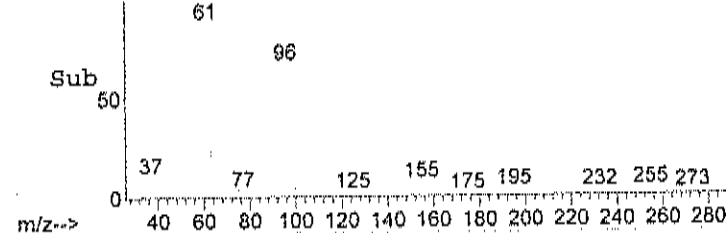
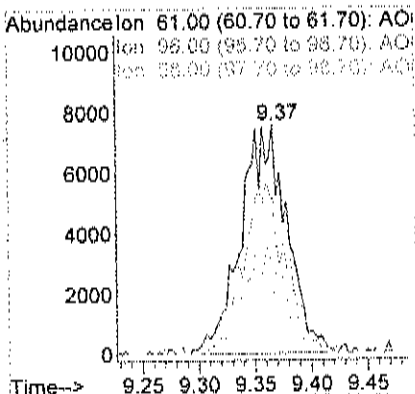
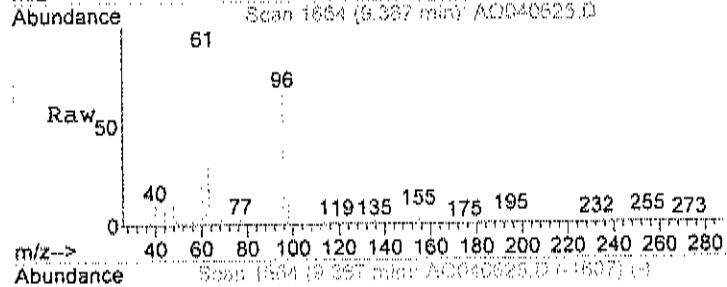
Method : C:\HPCHEM\1\METHODS\A331\_LUG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration





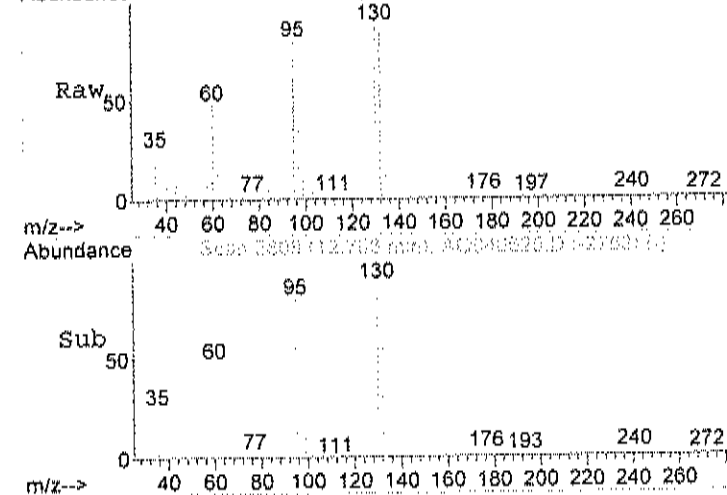
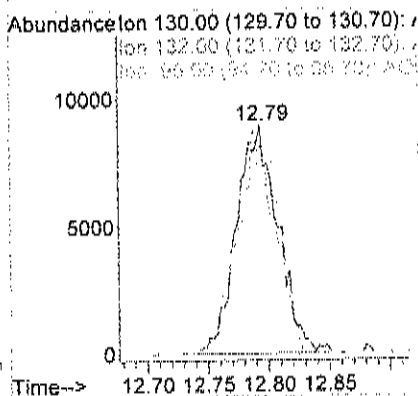
#29  
cis-1,2-dichloroethene  
Concen: 0.40 ppb  
RT: 9.37 min Scan# 1664  
Delta R.T. 0.02 min  
Lab File: AO040625.D  
Acq: 7 Apr 2017 12:50 am

Tgt	Ion: 61	Resp:	21474
Ion	Ratio	Lower	Upper
61	100		
96	71.4	58.1	98.1
98	41.9	29.3	69.3



#44  
Trichloroethene  
Concen: 0.41 ppb  
RT: 12.79 min Scan# 2808  
Delta R.T. 0.01 min  
Lab File: AO040625.D  
Acq: 7 Apr 2017 12:50 am

Tgt Ion:130	Resp:	21262	
Ion	Ratio	Lower	Upper
130	100		
132	86.9	69.9	109.9
95	102.2	76.3	116.3



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-005A

Client Sample ID: 691-B15-SV1  
Tag Number: 550.266  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
				FLD		Analyst:
Lab Vacuum In	-2			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
				TO-15		Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 4:03:00 PM
Chloroethane	0.14	0.15	J	ppbV	1	4/7/2017 4:03:00 PM
cis-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 4:03:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 4:03:00 PM
Trichloroethene	40	3.6		ppbV	90	4/7/2017 9:20:00 PM
Vinyl chloride	0.11	0.040		ppbV	1	4/7/2017 4:03:00 PM
Surr: Bromofluorobenzene	118	70-130		%REC	1	4/7/2017 4:03:00 PM

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-005A

Client Sample ID: 691-B15-SV1  
Tag Number: 550.266  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC				TO-15	Analyst: RJP	
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:03:00 PM
Chloroethane	0.37	0.40	J	ug/m3	1	4/7/2017 4:03:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:03:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:03:00 PM
Trichloroethene	220	19		ug/m3	90	4/7/2017 9:20:00 PM
Vinyl chloride	0.28	0.10		ug/m3	1	4/7/2017 4:03:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\AO040709.D  
 Acq On : 7 Apr 2017 4:03 pm  
 Sample : C1704014-005A  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 07 16:58:03 2017

Vial: 9  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Mon Apr 03 10:15:59 2017  
 Response via : Initial Calibration  
 DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	24680	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.94	114	109137	1.00	ppb	0.00
50) Chlorobenzene-d5	16.83	117	104661	1.00	ppb	0.00

System Monitoring Compounds

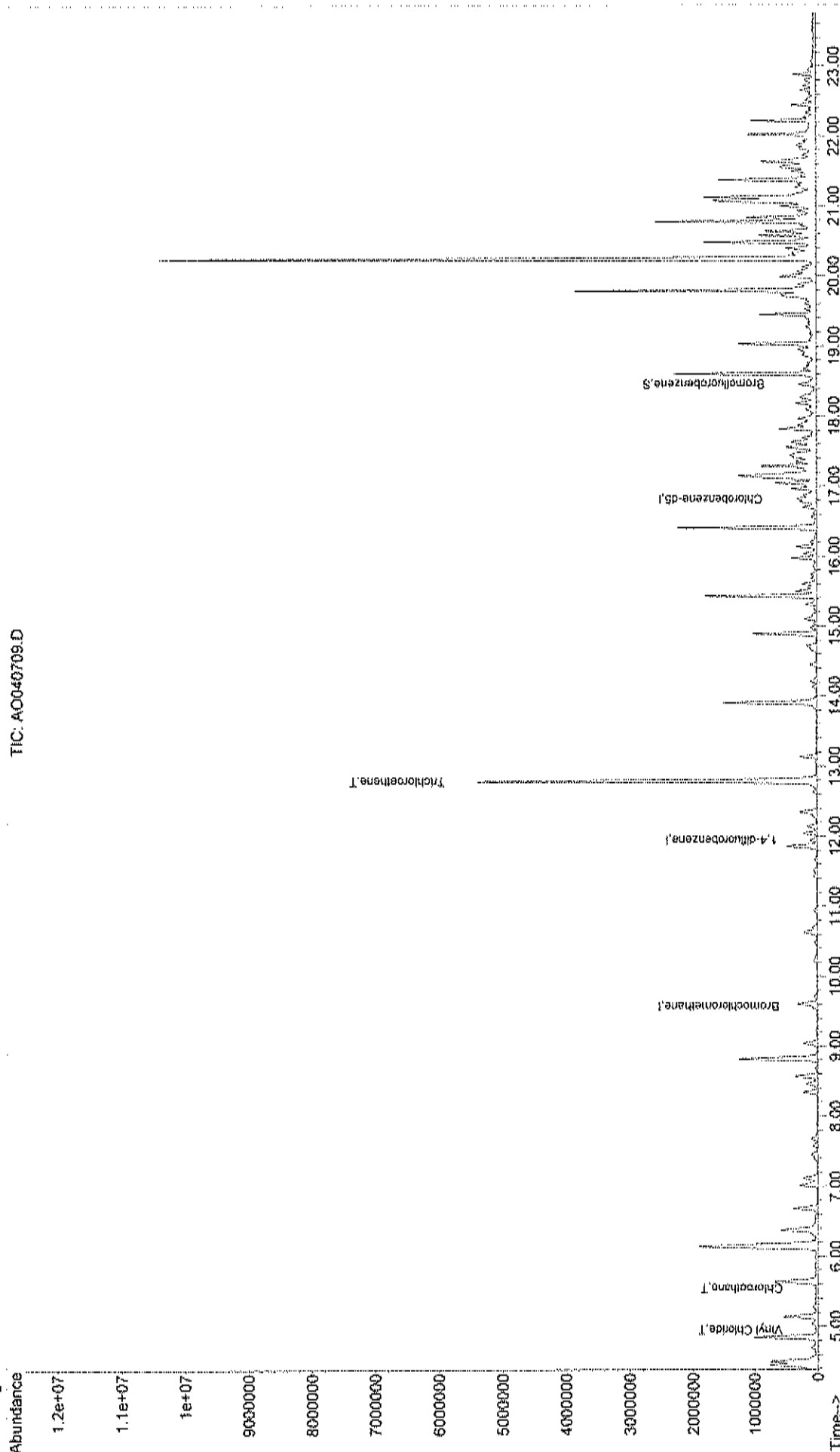
65) Bromofluorobenzene	18.46	95	84188	1.18	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	118.00%

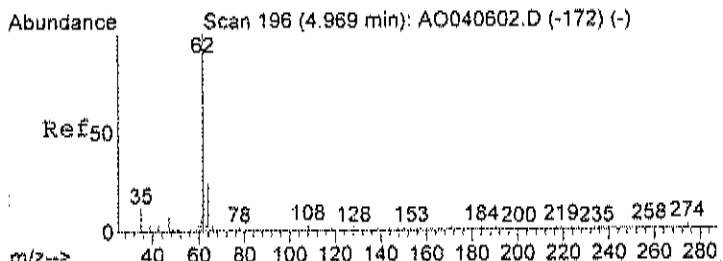
Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
6) Vinyl Chloride	4.94	62	4732	0.11	ppb	97
10) Chloroethane	5.53	64	2571	0.14	ppb	92
44) Trichloroethene	12.79	130	2088134	42.51	ppb	92

Data File : C:\HPCHEM\1\DATA\AO040709.D  
 Acq On : 7 Apr 2017 4:03 pm  
 Sample : C1704014-005A  
 Misc : A331\_IUG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 11 14:08 2017

Vial: 9  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00  
 Quant Results File: A331\_IUG.RES

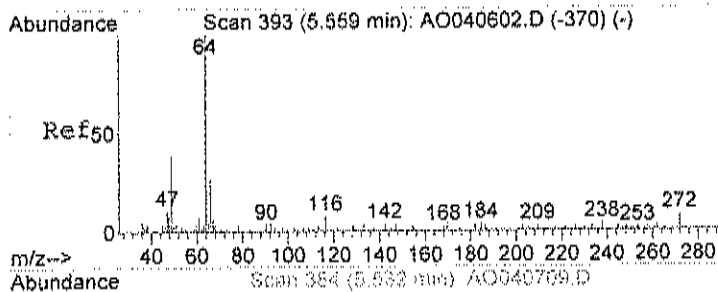
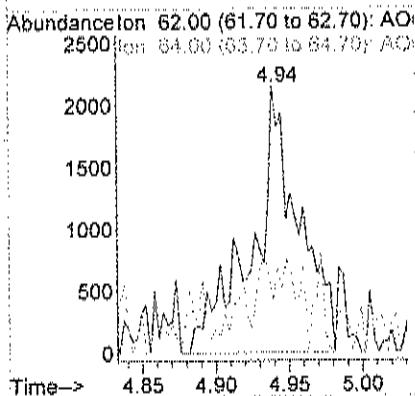
Method : C:\HPCHEM\1\METHODS\A331\_IUG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration





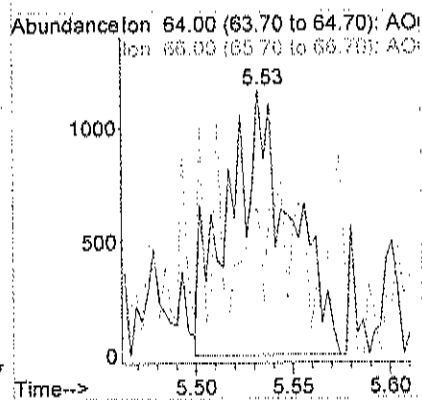
#6  
 Vinyl Chloride  
 Concen: 0.11 ppb  
 RT: 4.94 min Scan# 186  
 Delta R.T. -0.01 min  
 Lab File: AO040709.D  
 Acq: 7 Apr 2017 4:03 pm

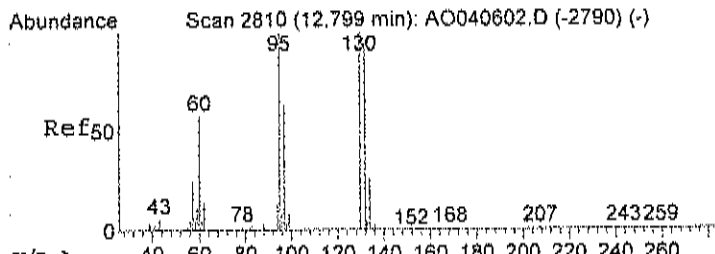
Tgt Ion	Ratio	Lower	Upper
62	100		
64	27.2	0.0	58.6



#10  
 Chloroethane  
 Concen: 0.14 ppb  
 RT: 5.53 min Scan# 384  
 Delta R.T. -0.00 min  
 Lab File: AO040709.D  
 Acq: 7 Apr 2017 4:03 pm

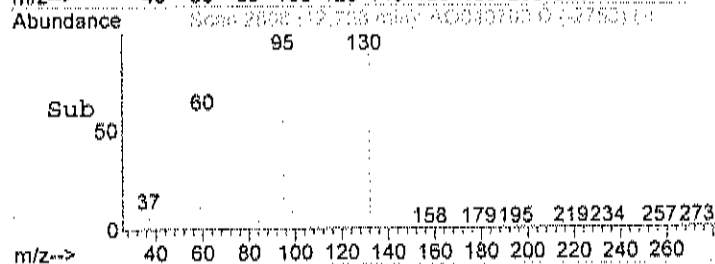
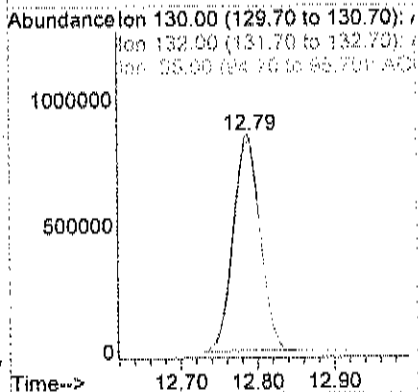
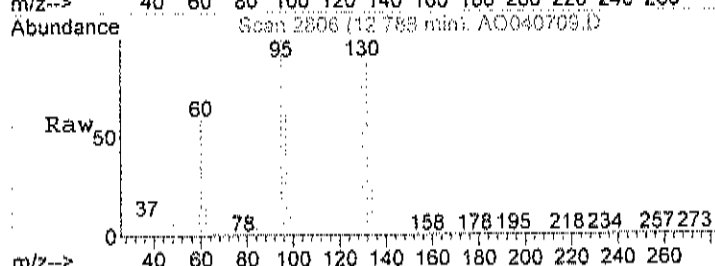
Tgt Ion	Ratio	Lower	Upper
64	100		
66	32.5	22.7	34.1





#44  
 Trichloroethene  
 Concen: 42.51 ppb  
 RT: 12.79 min Scan# 2806  
 Delta R.T. 0.01 min  
 Lab File: AO040709.D  
 Acq: 7 Apr 2017 4:03 pm

Tgt Ion	Ratio	Lower	Upper
130	100		
132	95.8	69.9	109.9
95	106.5	76.3	116.3





Data File : C:\HPCHEM\1\DATA\AO040716.D

Vial: 16

Acq On : 7 Apr 2017 9:20 pm

Operator: RJP

Sample : C1704014-005A 90x

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 08 08:59:51 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	26112	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	119025	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	102419	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	63983	0.92	ppb	0.01
Spiked Amount	1.000	Range	70 - 130	Recovery	=	92.00%

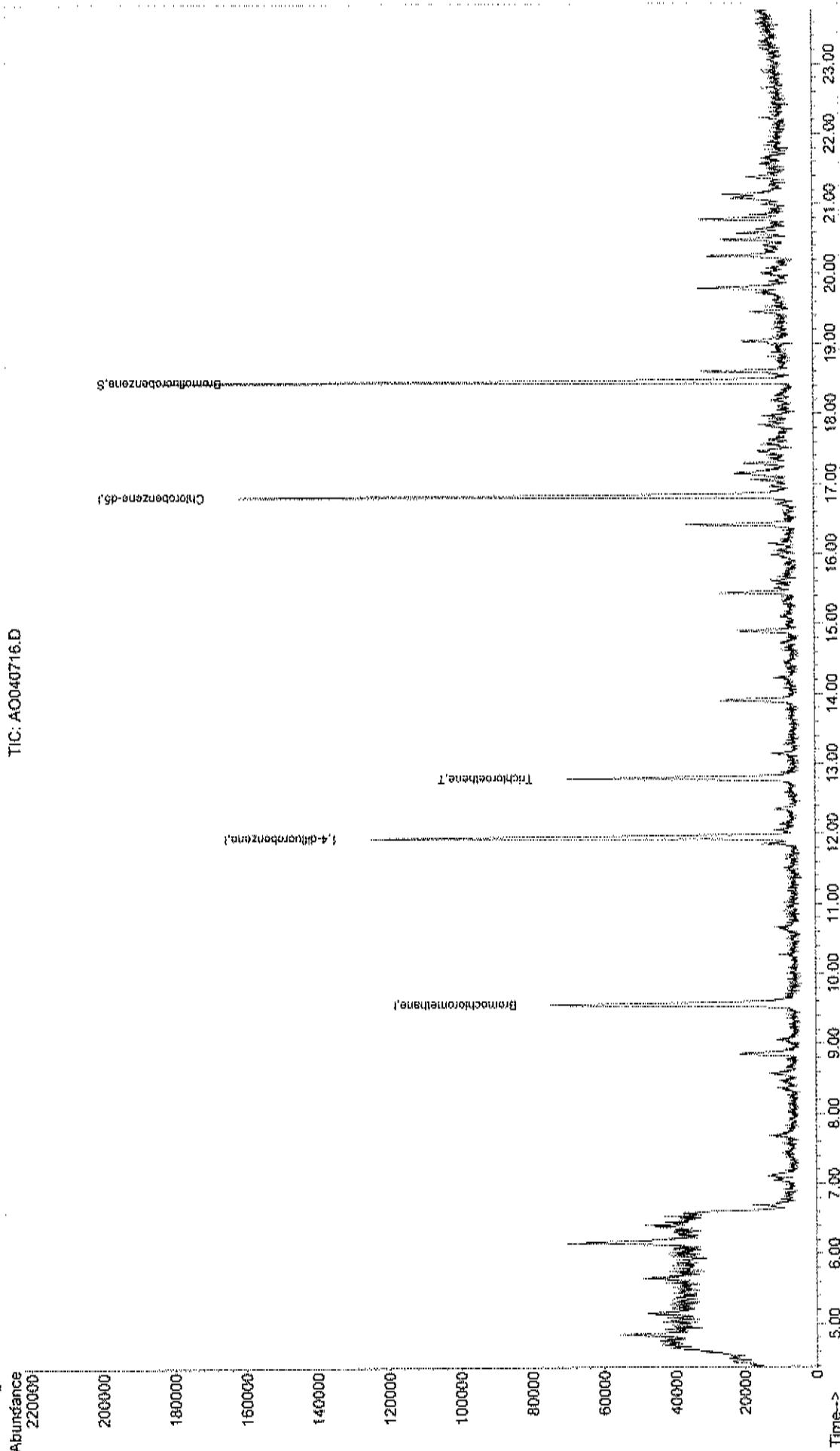
## Target Compounds

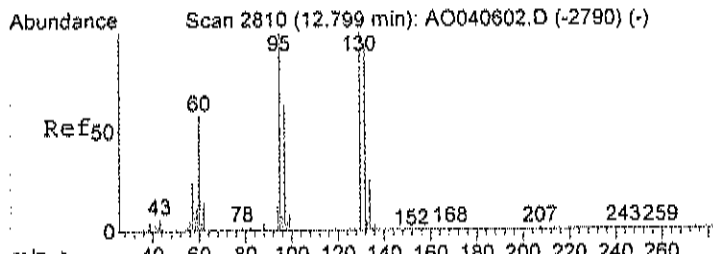
44) Trichloroethene	12.79	130	23989	0.45	ppb	Qvalue 87
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Data File : C:\HPCHEM\1\DATA\AO040716.D  
 Acq On : 7 Apr 2017 9:20 pm  
 Sample : C1704014-005A 90x  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 11 14:59 2017

Vial: 16  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00  
 Quant Results File: A331\_1UG.RES

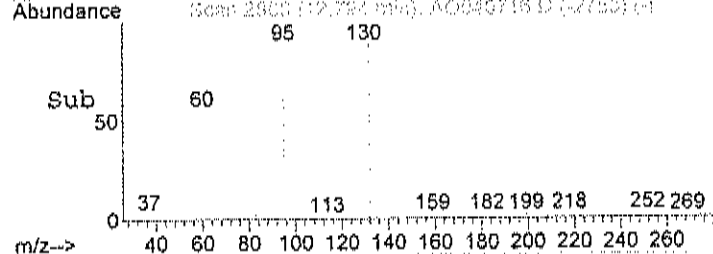
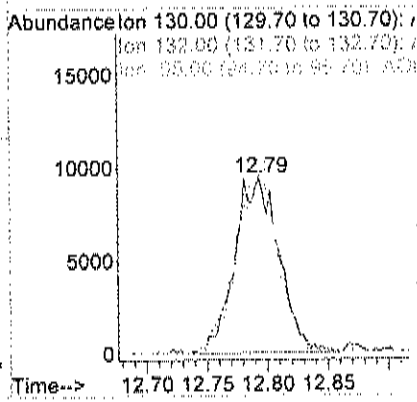
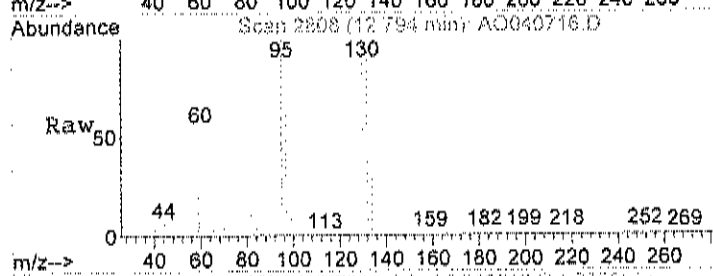
Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration





#44  
 Trichloroethene  
 Concen: 0.45 ppb  
 RT: 12.79 min Scan# 2808  
 Delta R.T. 0.01 min  
 Lab File: AO040716.D  
 Acq: 7 Apr 2017 9:20 pm

Tgt Ion: 130	Resp: 23989
Ion Ratio	Lower Upper
130	100
132	101.1 69.9 109.9
95	110.6 76.3 116.3



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-006A

Client Sample ID: 691-Outdoor-04012017  
Tag Number: 240,340  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-8			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 1:31:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 1:31:00 AM
cis-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 1:31:00 AM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 1:31:00 AM
Trichloroethene	< 0.040	0.040		ppbV	1	4/7/2017 1:31:00 AM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 1:31:00 AM
Surr: Bromofluorobenzene	88.0	70-130		%REC	1	4/7/2017 1:31:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte, Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-006A

Client Sample ID: 691-Outdoor-04012017  
Tag Number: 240.340  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 1:31:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 1:31:00 AM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 1:31:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 1:31:00 AM
Trichloroethene	< 0.21	0.21		ug/m3	1	4/7/2017 1:31:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 1:31:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection



Data File : C:\HPCHEM\1\DATA\AO040626.D  
 Acq On : 7 Apr 2017 1:31 am  
 Sample : C1704014-006A  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 07 07:19:06 2017

Vial: 24  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Mon Apr 03 10:15:59 2017  
 Response via : Initial Calibration  
 DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	25816	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	110508	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	98499	1.00	ppb	0.01

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	58874	0.88	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	88.00%

Target Compounds

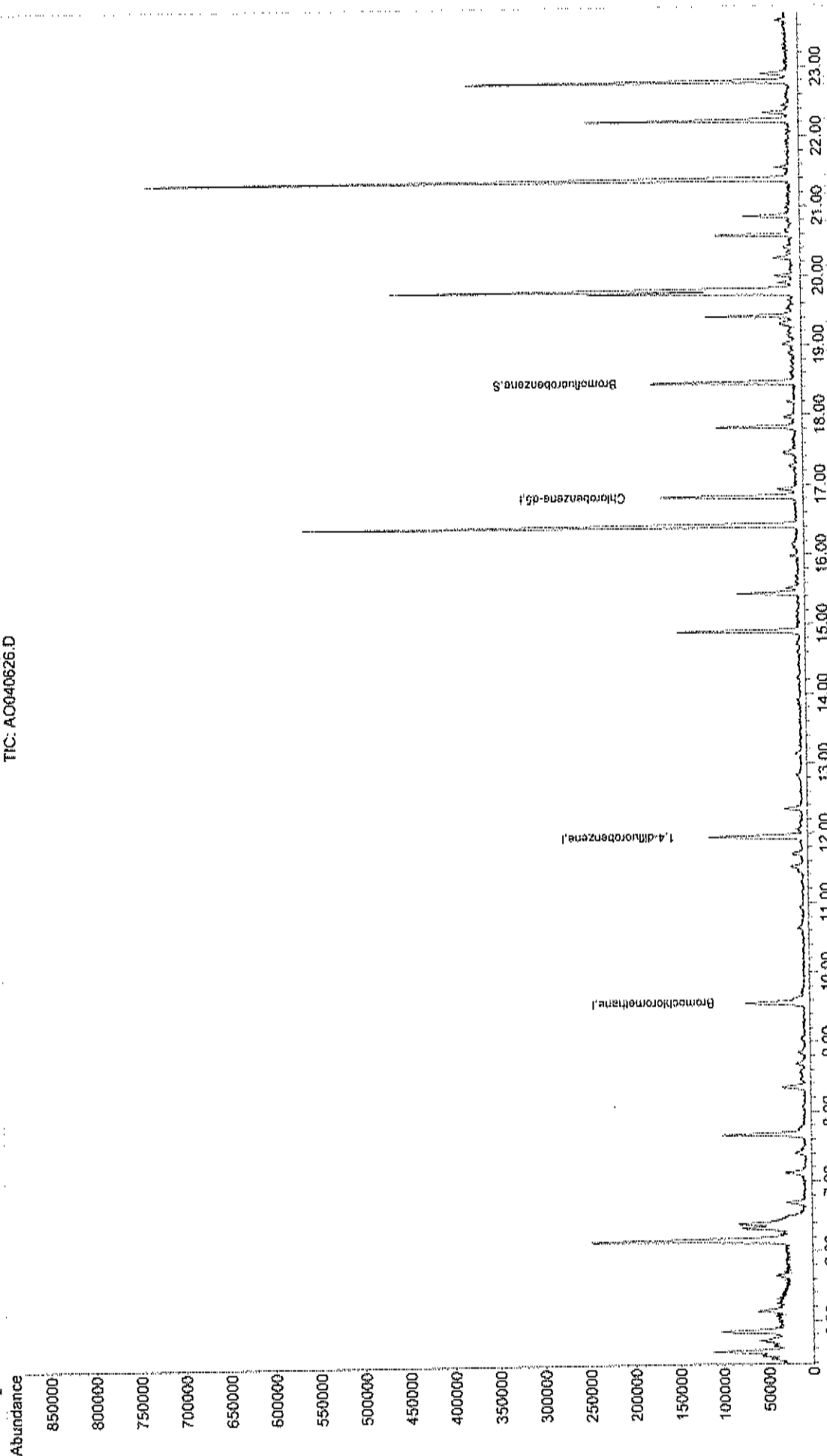
Qvalue

Data File : C:\HPCHEM\1\DATA\A0040626.D  
Acq On : 7 Apr 2017 1:31 am  
Sample : C1704014-006A  
Misc : A331\_IUG  
MS Integration Params: RTEINT.P  
Quant Time: Apr 7 13:25 2017

Vial: 24  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00  
Quant Results File: A331\_IUG.RES

Method : C:\HPCHEM\1\METHODS\A331\_IUG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Thu May 04 11:27:28 2017  
Response via : Initial Calibration

TIC: A0040626.D



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-007A

Client Sample ID: 691-Duplicate  
Tag Number: 101.299  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
FIELD PARAMETERS						Analyst:
				FLD		
Lab Vacuum In	-3			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC						Analyst: RJP
				TO-15		
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 2:12:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 2:12:00 AM
cis-1,2-Dichloroethene	0.42	0.15		ppbV	1	4/7/2017 2:12:00 AM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 2:12:00 AM
Trichloroethene	0.41	0.040		ppbV	1	4/7/2017 2:12:00 AM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 2:12:00 AM
Surr: Bromofluorobenzene	91.0	70-130		%REC	1	4/7/2017 2:12:00 AM

Qualifiers:	**	Quantitation Limit	-	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte, Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-007A

Client Sample ID: 691-Duplicate  
Tag Number: 101.299  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 2:12:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 2:12:00 AM
cis-1,2-Dichloroethene	1.7	0.59		ug/m3	1	4/7/2017 2:12:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 2:12:00 AM
Trichloroethene	2.2	0.21		ug/m3	1	4/7/2017 2:12:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 2:12:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

Page 7 of 15

Data File : C:\HPCHEM\1\DATA\AO040627.D  
Acq On : 7 Apr 2017 2:12 am  
Sample : C1704014-007A  
Misc : A331\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Apr 07 07:19:07 2017

Vial: 25  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Apr 03 10:15:59 2017  
Response via : Initial Calibration  
DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.57	128	24957	1.00	ppb	0.03
35) 1,4-difluorobenzene	11.94	114	110993	1.00	ppb	0.00
50) Chlorobenzene-d5	16.83	117	99559	1.00	ppb	0.01

System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	61990	0.91	ppb	0.01
Spiked Amount	1.000	Range	70 - 130	Recovery	=	91.00%

Target Compounds						Qvalue
29) cis-1,2-dichloroethene	9.36	61	21892	0.42	ppb	93
44) Trichloroethene	12.79	130	20378	0.41	ppb	92

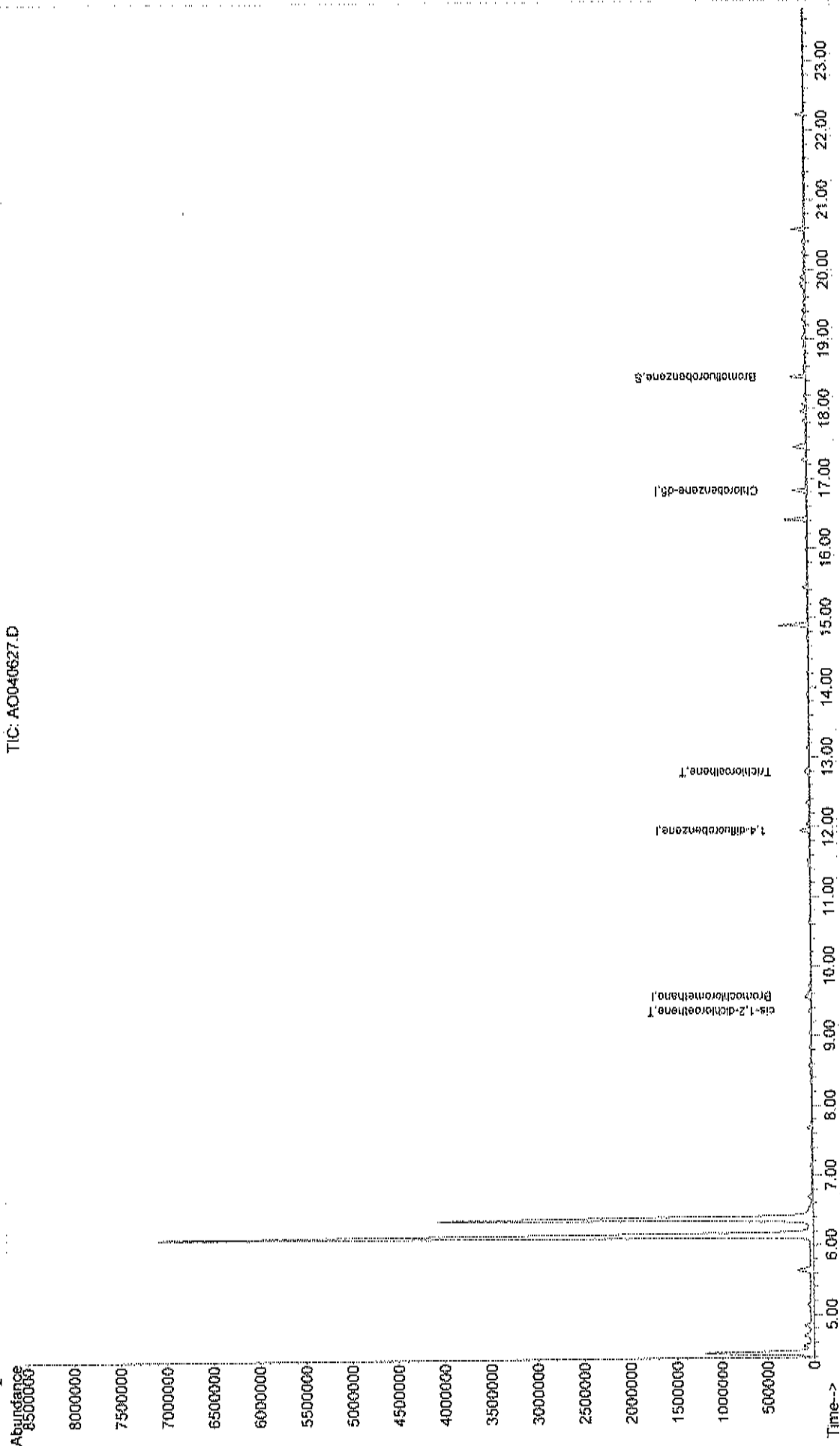


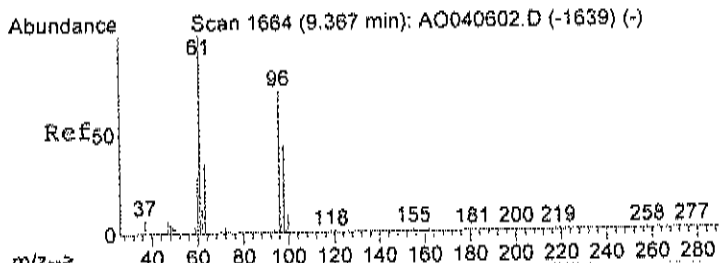
Data File : C:\HPCHEM\1\DATA\AO040627.D  
Acq On : 7 Apr 2017 2:12 am  
Sample : C1704014-007A  
Misc : A331\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Apr 7 13:26 2017

Vial: 25  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A331\_1UG.RES

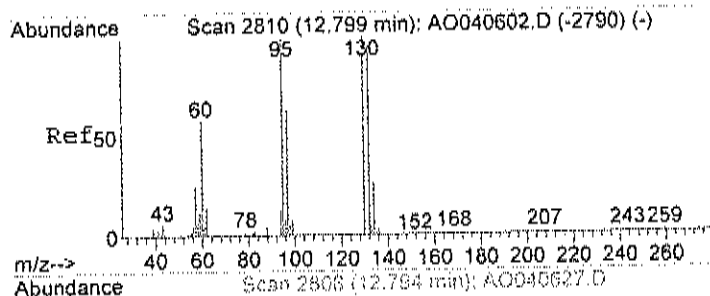
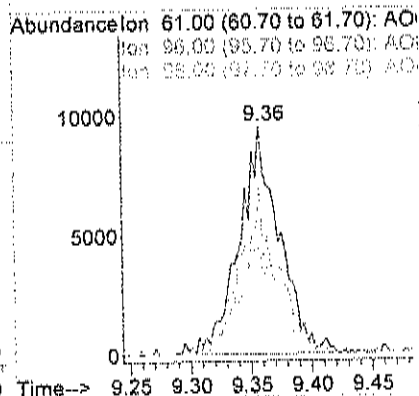
Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Thu May 04 11:27:28 2017  
Response via : Initial Calibration





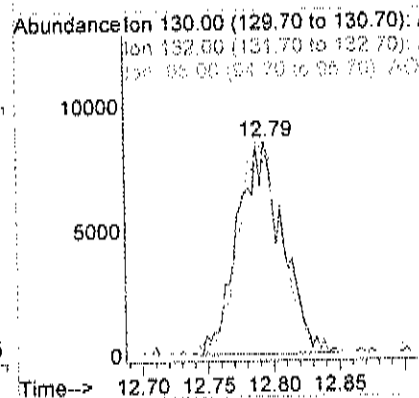
#29  
cis-1,2-dichloroethene  
Concen: 0.42 ppb  
RT: 9.36 min Scan# 1661  
Delta R.T. 0.01 min  
Lab File: AO040627.D  
Acq: 7 Apr 2017 2:12 am

Tgt Ion:	61	Resp:	21892
Ion	Ratio	Lower	Upper
61	100		
96	68.4	58.1	98.1
98	48.2	29.3	69.3



#44  
Trichloroethene  
Concen: 0.41 ppb  
RT: 12.79 min Scan# 2808  
Delta R.T. 0.01 min  
Lab File: AO040627.D  
Acq: 7 Apr 2017 2:12 am

Tgt Ion:	130	Resp:	20378
Ion	Ratio	Lower	Upper
130	100		
132	97.4	69.9	109.9
95	104.4	76.3	116.3



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-008A

Client Sample ID: 691-NE-1AQ  
Tag Number: 94.379  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-1			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 2:52:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 2:52:00 AM
cis-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 2:52:00 AM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 2:52:00 AM
Trichloroethene	< 0.040	0.040		ppbV	1	4/7/2017 2:52:00 AM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 2:52:00 AM
Surr: Bromofluorobenzene	93.0	70-130		%REC	1	4/7/2017 2:52:00 AM

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.

Client Sample ID: 691-NE-IAQ

Lab Order: C1704014

Tag Number: 94.379

Project: 691 St Paul Street

Collection Date: 4/1/2017

Lab ID: C1704014-008A

Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 2:52:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 2:52:00 AM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 2:52:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 2:52:00 AM
Trichloroethene	< 0.21	0.21		ug/m3	1	4/7/2017 2:52:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 2:52:00 AM

## Qualifiers:

\*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

Data File : C:\HPCHEM\1\DATA\AO040628.D

Vial: 26

Acq On : 7 Apr 2017 2:52 am

Operator: RJP

Sample : C1704014-008A

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 07 07:19:08 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	25088	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	108028	1.00	ppb	0.02
50) Chlorobenzene-d5	16.83	117	95812	1.00	ppb	0.01

## System Monitoring Compounds

65) Bromofluorobenzene	18.47	95	60915	0.93	ppb	0.02
Spiked Amount	1.000	Range	70 - 130	Recovery	=	93.00%

Target Compounds

Qvalue

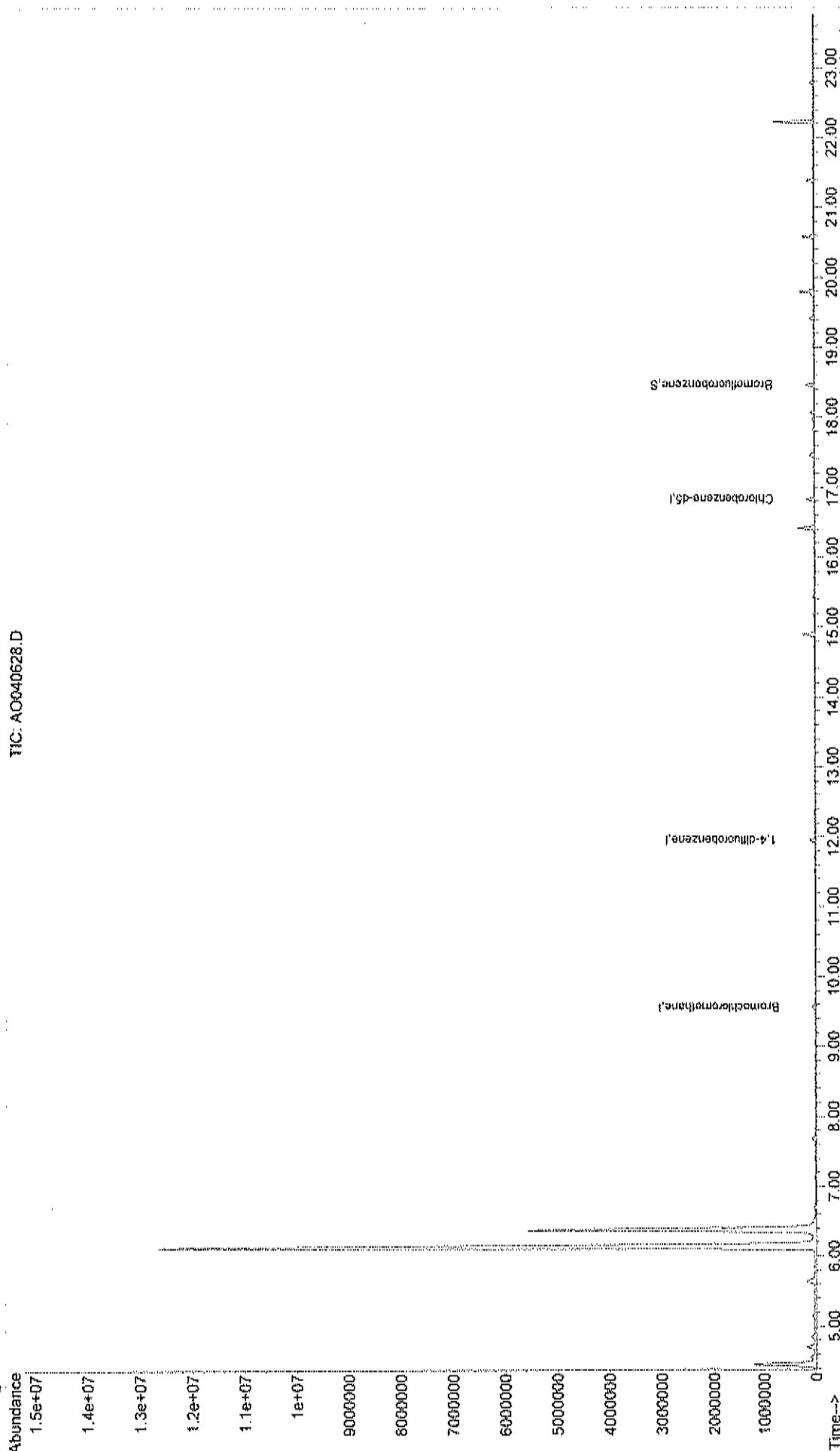


Data File : C:\HPCHEM\1\DATA\AO040628.D  
 Acq On : 7 Apr 2017 2:52 am  
 Sample : C1704014-008A  
 Misc : A331\_IUG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 7 13:27 2017

Vial: 26  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_IUG.RES

Method : C:\HPCHEM\1\METHODS\A331\_IUG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-009A

Client Sample ID: 691-NE-SVI  
Tag Number: 1186.145  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						Analyst:
Lab Vacuum In	-3			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						Analyst: RJP
		TO-15				
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 5:28:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 5:28:00 PM
cis-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 5:28:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 5:28:00 PM
Trichloroethene	0.72	0.040		ppbV	1	4/7/2017 5:28:00 PM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 5:28:00 PM
Surr: Bromofluorobenzene	113	70-130		%REC	1	4/7/2017 5:28:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte, Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-009A

Client Sample ID: 691-NE-SV1  
Tag Number: 1186.145  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 5:28:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 5:28:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 5:28:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 5:28:00 PM
Trichloroethene	3.9	0.21		ug/m3	1	4/7/2017 5:28:00 PM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 5:28:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

Data File : C:\HPCHEM\1\DATA\AO040710.D

Vial: 10

Acq On : 7 Apr 2017 5:28 pm

Operator: RJP

Sample : C1704014-009A

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 08 08:58:21 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	27394	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	125972	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	121869	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	94050m	1.13	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	113.00%

## Target Compounds

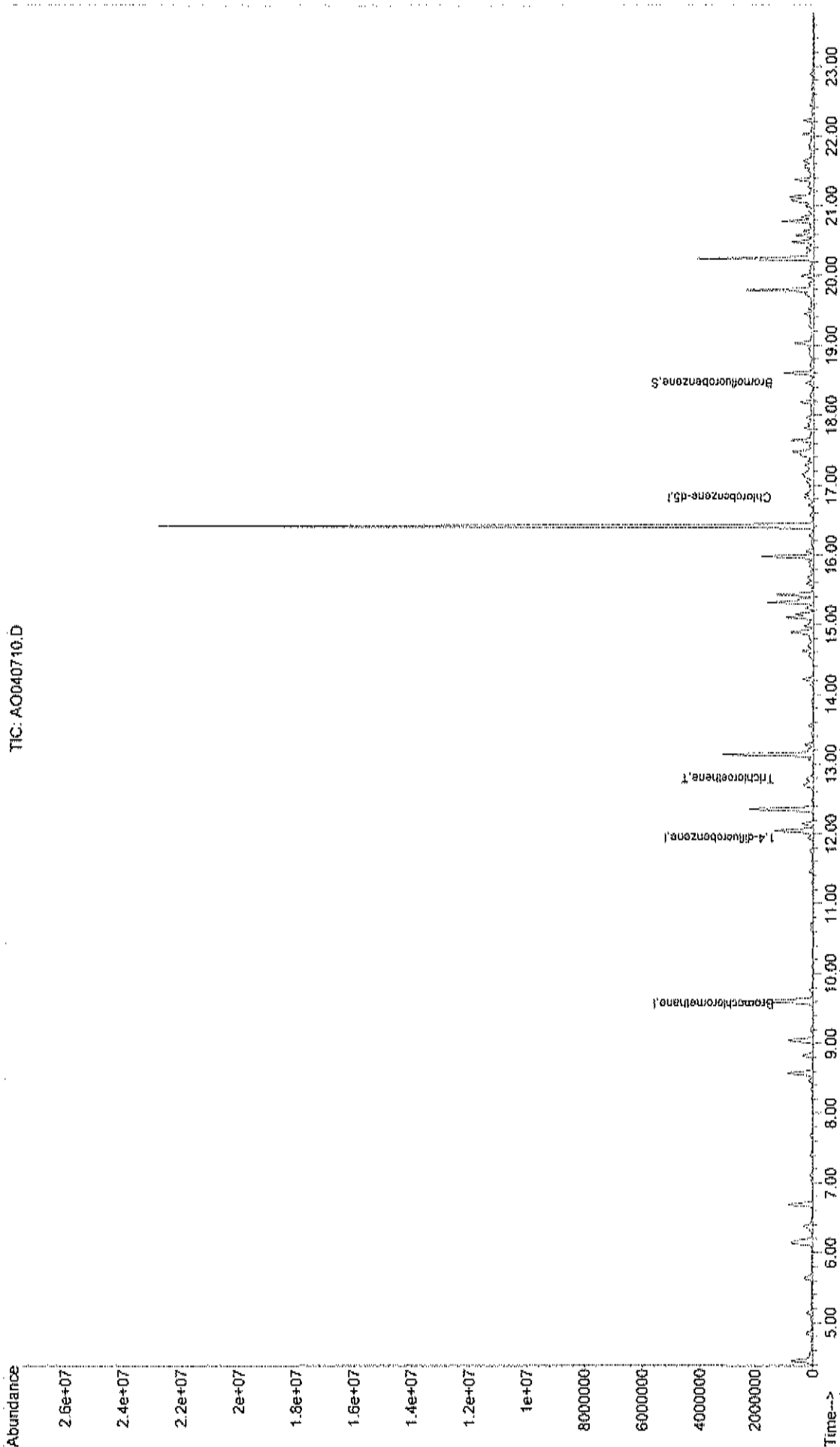
44) Trichloroethene	12.79	130	40805	0.72	ppb	Qvalue 86
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Data File : C:\HPCHEM\1\DATA\AO040710.D  
 Acq On : 7 Apr 2017 5:28 pm  
 Sample : C1704014-009A  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 11 14:10 2017

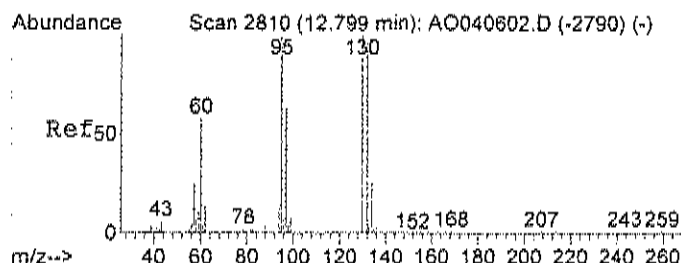
Vial: 10  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration

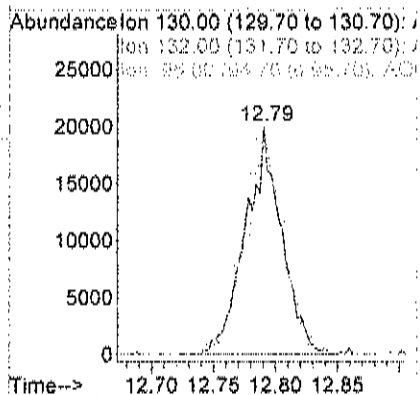






#44  
 Trichloroethene  
 Concen: 0.72 ppb  
 RT: 12.79 min Scan# 2807  
 Delta R.T. 0.01 min  
 Lab File: AO040710.D  
 Acq: 7 Apr 2017 5:28 pm

Tgt Ion: 130 Resp: 40805  
 Ion Ratio Lower Upper  
 130 100  
 132 99.9 69.9 109.9  
 95 113.5 76.3 116.3



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.

Client Sample ID: 691-SB1-1AQ

Lab Order: C1704014

Tag Number: 170.387

Project: 691 St Paul Street

Collection Date: 4/1/2017

Lab ID: C1704014-010A

Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-5			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 3:33:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 3:33:00 AM
cis-1,2-Dichloroethene	0.66	0.15		ppbV	1	4/7/2017 3:33:00 AM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 3:33:00 AM
Trichloroethene	0.24	0.040		ppbV	1	4/7/2017 3:33:00 AM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 3:33:00 AM
Surr: Bromofluorobenzene	92.0	70-130		%REC	1	4/7/2017 3:33:00 AM

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-010A

Client Sample ID: 691-SB1-IAQ  
Tag Number: 170.387  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 3:33:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 3:33:00 AM
cis-1,2-Dichloroethene	2.6	0.59		ug/m3	1	4/7/2017 3:33:00 AM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 3:33:00 AM
Trichloroethene	1.3	0.21		ug/m3	1	4/7/2017 3:33:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 3:33:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

Data File : C:\HPCHEM\1\DATA\AO040629.D

Vial: 27

Acq On : 7 Apr 2017 3:33 am

Operator: RJP

Sample : C1704014-010A

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 07 07:19:09 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	24485	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	107880	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	98198	1.00	ppb	0.01

## System Monitoring Compounds

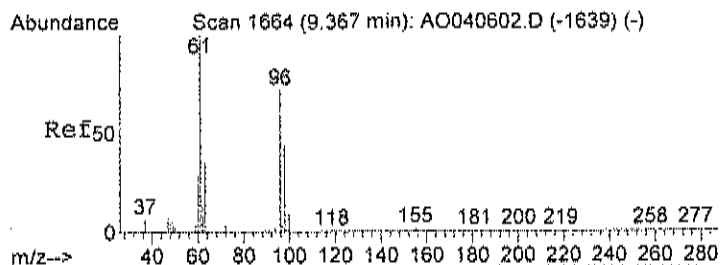
65) Bromofluorobenzene	18.47	95	61676	0.92	ppb	0.02
Spiked Amount	1.000	Range	70 - 130	Recovery	=	92.00%

## Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
29) cis-1,2-dichloroethene	9.36	61	33781	0.66	ppb	98
44) Trichloroethene	12.79	130	11462	0.24	ppb	# 81

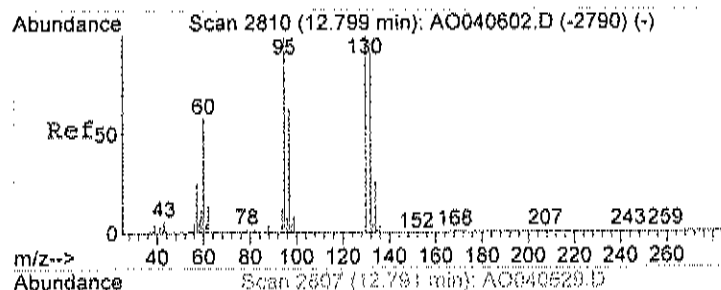
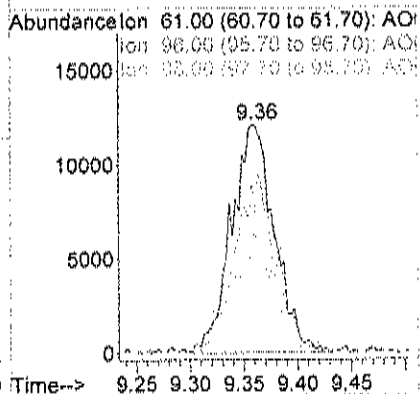






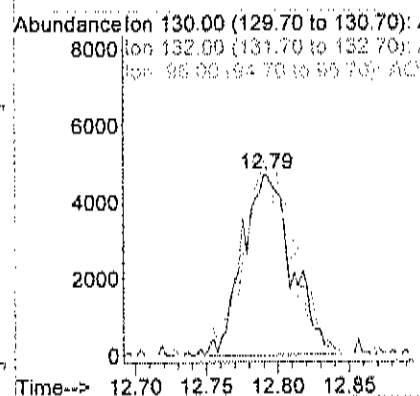
#29  
 cis-1,2-dichloroethene  
 Concen: 0.66 ppb  
 RT: 9.36 min Scan# 1661  
 Delta R.T. 0.01 min  
 Lab File: AO040629.D  
 Acq: 7 Apr 2017 3:33 am

Tgt Ion	Ratio	Lower	Upper
61	100		
96	75.3	58.1	98.1
98	48.7	29.3	69.3



#44  
 Trichloroethene  
 Concen: 0.24 ppb  
 RT: 12.79 min Scan# 2807  
 Delta R.T. 0.01 min  
 Lab File: AO040629.D  
 Acq: 7 Apr 2017 3:33 am

Tgt Ion	Ratio	Lower	Upper
130	100		
132	100.4	69.9	109.9
95	122.7	76.3	116.3#



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-011A

Client Sample ID: 691-SB1-SVI  
Tag Number: 203.372  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>		<b>FLD</b>		<b>Analyst:</b>		
Lab Vacuum In	-6			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>		<b>TO-15</b>		<b>Analyst: RJP</b>		
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 6:09:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 6:09:00 PM
cis-1,2-Dichloroethene	13	1.5		ppbV	10	4/7/2017 11:10:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 6:09:00 PM
Trichloroethene	0.74	0.040		ppbV	1	4/7/2017 6:09:00 PM
Vinyl chloride	0.15	0.040		ppbV	1	4/7/2017 6:09:00 PM
Surr: Bromofluorobenzene	109	70-130		%REC	1	4/7/2017 6:09:00 PM

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-011A

Client Sample ID: 691-SB1-SV1  
Tag Number: 203.372  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 6:09:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 6:09:00 PM
cis-1,2-Dichloroethene	52	5.9		ug/m3	10	4/7/2017 11:10:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 6:09:00 PM
Trichloroethene	4.0	0.21		ug/m3	1	4/7/2017 6:09:00 PM
Vinyl chloride	0.38	0.10		ug/m3	1	4/7/2017 6:09:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

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Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\AO040711.D  
 Acq On : 7 Apr 2017 6:09 pm  
 Sample : C1704014-011A  
 Misc : A331\_1UG

Vial: 11  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 08 08:58:50 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	30170	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.94	114	132762	1.00	ppb	0.00
50) Chlorobenzene-d5	16.83	117	121408	1.00	ppb	0.00

System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	90648	1.09	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	109.00%

Target Compounds

						Qvalue
6) Vinyl Chloride	4.95	62	7786	0.15	ppb	99
29) cis-1,2-dichloroethene	9.35	61	931425	14.87	ppb	97
44) Trichloroethene	12.79	130	44293	0.74	ppb	86

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

AO040711.D A331\_1UG.M

Thu May 04 11:29:33 2017

MSD1

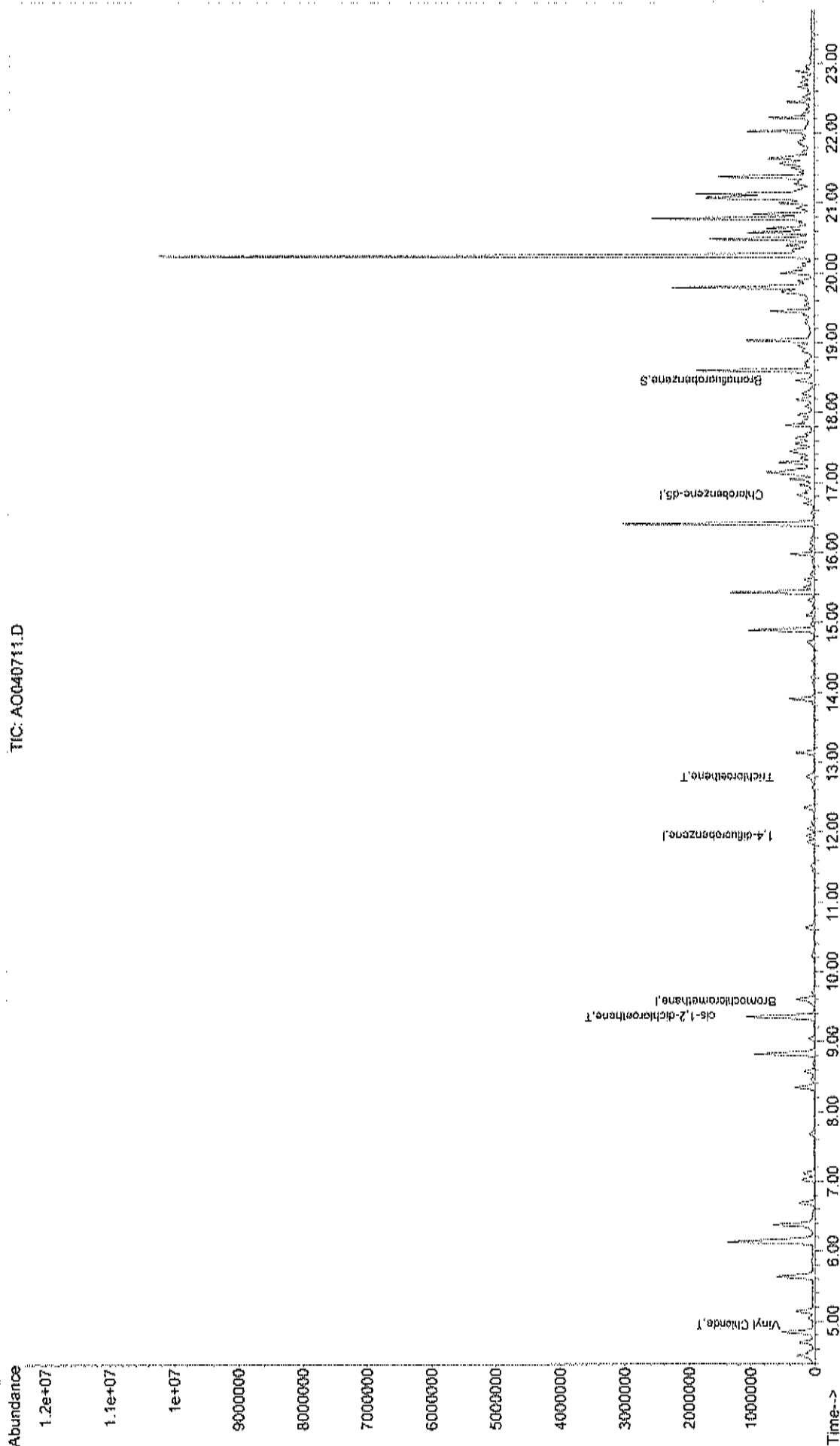
Page 1

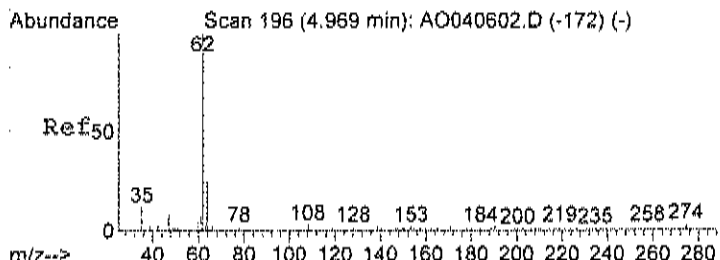
Data File : C:\HPCHEM\1\DATA\AO040711.D  
 Acq On : 7 Apr 2017 6:09 pm  
 Sample : C1704014-011A  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 11 14:13 2017

Vial: 11  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

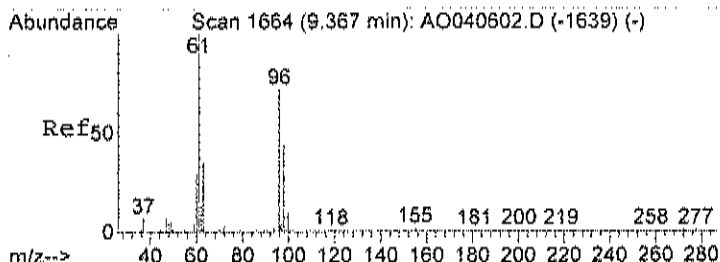
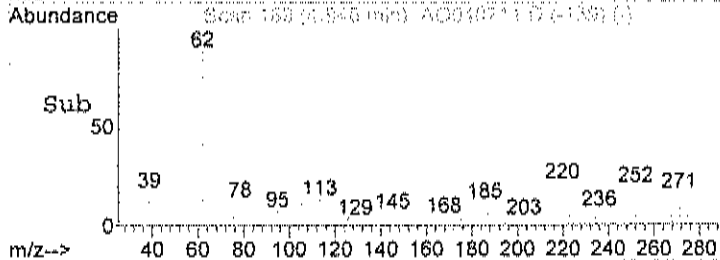
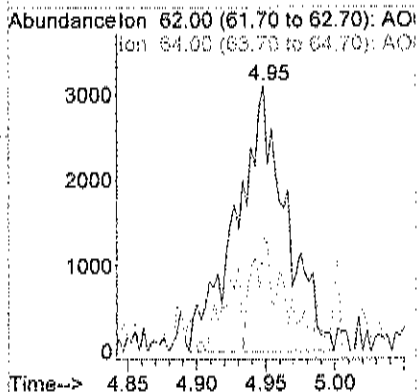
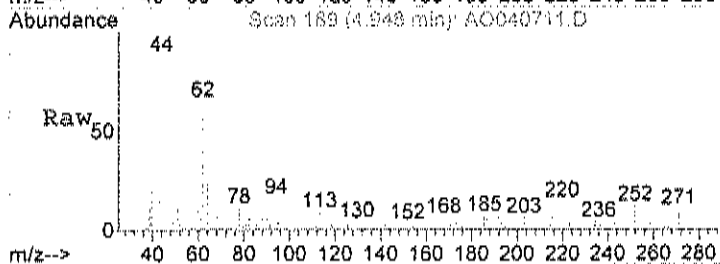
Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration





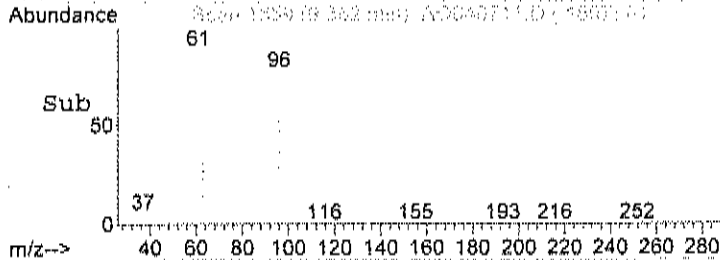
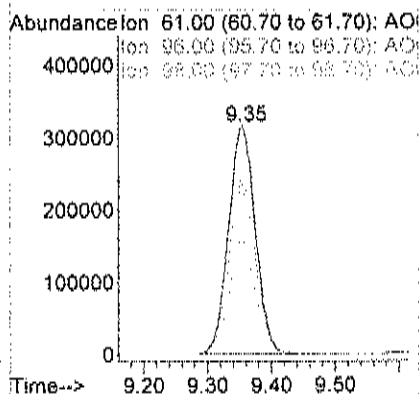
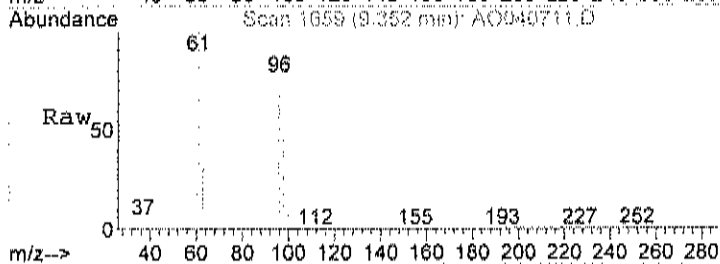
#6  
 Vinyl Chloride  
 Concen: 0.15 ppb  
 RT: 4.95 min Scan# 189  
 Delta R.T. 0.00 min  
 Lab File: AO040711.D  
 Acq: 7 Apr 2017 6:09 pm

Tgt Ion: 62 Resp: 7786  
 Ion Ratio Lower Upper  
 62 100  
 64 28.2 0.0 58.6

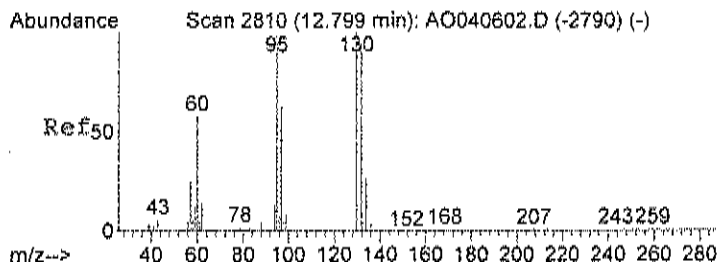


#29  
 cis-1,2-dichloroethene  
 Concen: 14.87 ppb  
 RT: 9.35 min Scan# 1659  
 Delta R.T. 0.01 min  
 Lab File: AO040711.D  
 Acq: 7 Apr 2017 6:09 pm

Tgt Ion: 61 Resp: 931425  
 Ion Ratio Lower Upper  
 61 100  
 96 74.8 58.1 98.1  
 98 48.9 29.3 69.3

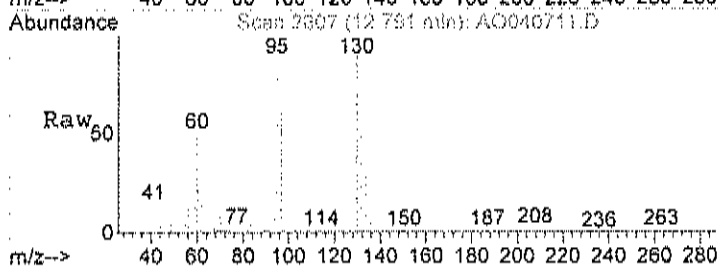






#44  
 Trichloroethene  
 Concen: 0.74 ppb  
 RT: 12.79 min Scan# 2807  
 Delta R.T. 0.01 min  
 Lab File: AO040711.D  
 Acq: 7 Apr 2017 6:09 pm

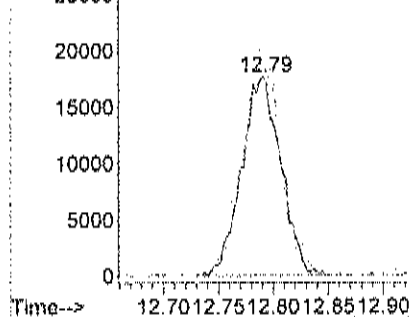
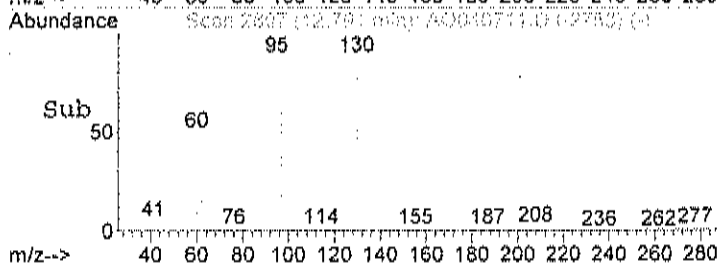
Tgt Ion	Resp	Lower	Upper
130	44293	100	100
132	101.5	69.9	109.9
95	111.9	76.3	116.3



Abundance Ion 130.00 (129.70 to 130.70):

Ion 132.00 (131.70 to 132.70):

Ion 95.00 (94.70 to 95.70):



Data File : C:\HPCHEM\1\DATA\AO040719.D

Vial: 19

Acq On : 7 Apr 2017 11:10 pm

Operator: RJP

Sample : C1704014-011A 10x

Inst : MSD #1

Misc : A331\_1UG

Multiplier: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 08 08:59:54 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	21591	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	93800	1.00	ppb	0.02
50) Chlorobenzene-d5	16.83	117	84591	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	53420	0.93	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	93.00%

## Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
29) cis-1,2-dichloroethene	9.35	61	58406	1.30	ppb	96

Data File : C:\HPCHEM\1\DATA\AO040719.D  
Acq On : 7 Apr 2017 11:10 pm  
Sample : C1704014-011A 10x  
Misc : A331\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Apr 11 15:00 2017

Vial: 19

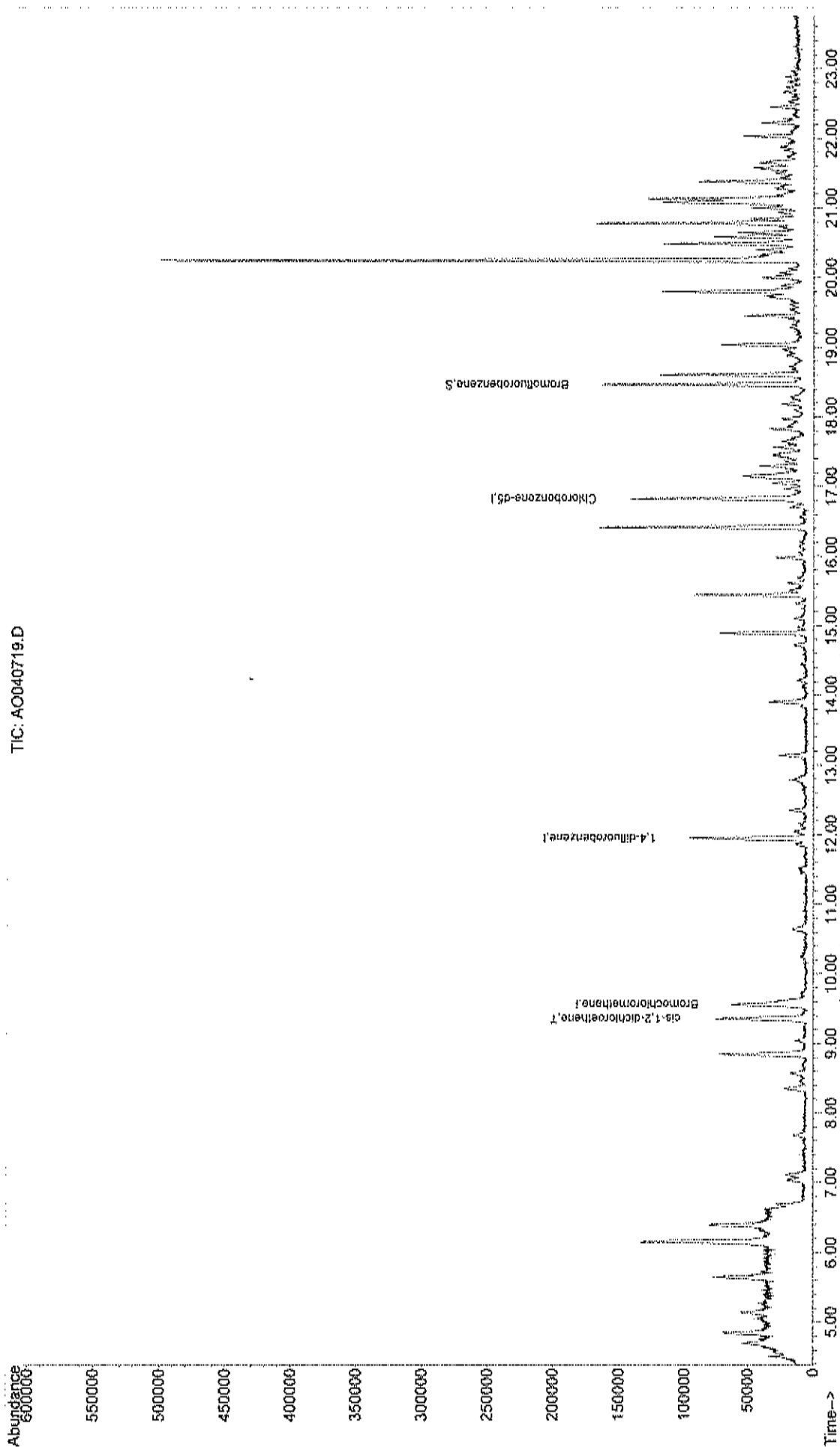
Operator: RJP

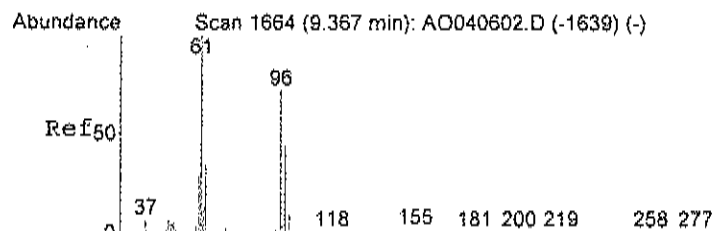
Inst : MSD #1

Multiplr: 1.00

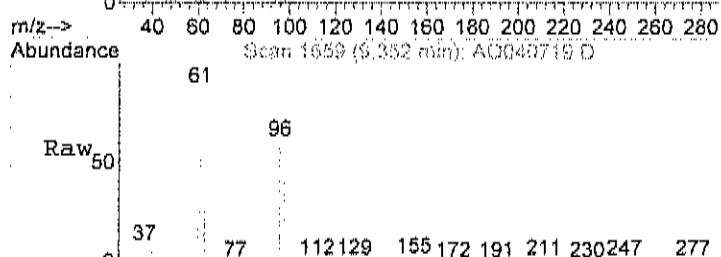
Quant Results File: A331\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Thu May 04 11:27:28 2017  
Response via : Initial Calibration

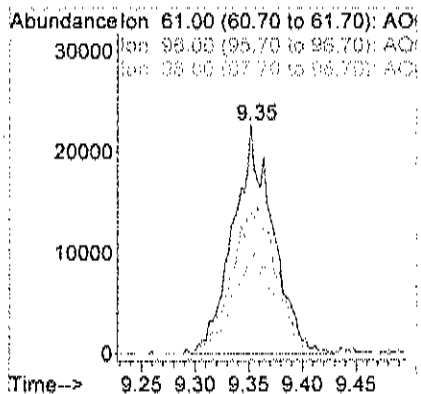
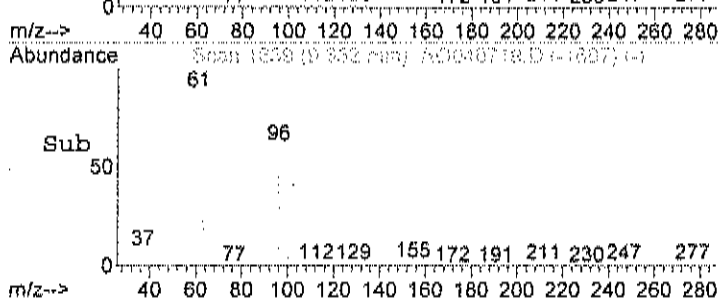




#29  
 cis-1,2-dichloroethene  
 Concen: 1.30 ppb  
 RT: 9.35 min Scan# 1659  
 Delta R.T. 0.01 min  
 Lab File: AO040719.D  
 Acq: 7 Apr 2017 11:10 pm



Tgt Ion	Ratio	Lower	Upper
61	100		
96	73.9	58.1	98.1
98	46.7	29.3	69.3



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-012A

Client Sample ID: 691-SB5A-1AQ  
Tag Number: 367.251  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-2			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 4:13:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 4:13:00 AM
cis-1,2-Dichloroethene	3.3	1.5		ppbV	10	4/7/2017 1:50:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 4:13:00 AM
Trichloroethene	1.2	0.040		ppbV	1	4/7/2017 4:13:00 AM
Vinyl chloride	< 0.040	0.040		ppbV	1	4/7/2017 4:13:00 AM
Surr: Bromofluorobenzene	103	70-130		%REC	1	4/7/2017 4:13:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.

Client Sample ID: 691-SB5A-IAQ

Lab Order: C1704014

Tag Number: 367.251

Project: 691 St Paul Street

Collection Date: 4/1/2017

Lab ID: C1704014-012A

Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:13:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 4:13:00 AM
cis-1,2-Dichloroethene	13	5.9		ug/m3	10	4/7/2017 1:50:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:13:00 AM
Trichloroethene	6.5	0.21		ug/m3	1	4/7/2017 4:13:00 AM
Vinyl chloride	< 0.10	0.10		ug/m3	1	4/7/2017 4:13:00 AM

## Qualifiers:

- \*\* Quantitation Limit
- B Analyte detected in the associated Method Blank
- H Holding times for preparation or analysis exceeded
- JN Non-routine analyte, Quantitation estimated,
- S Spike Recovery outside accepted recovery limits

- . Results reported are not blank corrected
- E Estimated Value above quantitation range
- J Analyte detected below quantitation limit
- ND Not Detected at the Limit of Detection

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Data File : C:\HPCHEM\1\DATA\AO040630.D

Vial: 28

Acq On : 7 Apr 2017 4:13 am

Operator: RJP

Sample : C1704014-012A

Inst : MSD #1

Misc : A331\_1UG

Multiplx: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 07 07:19:10 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	25278	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	111053	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	100304	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.47	95	70272	1.03	ppb	0.02
Spiked Amount	1.000	Range	70 - 130	Recovery	=	103.00%

## Target Compounds

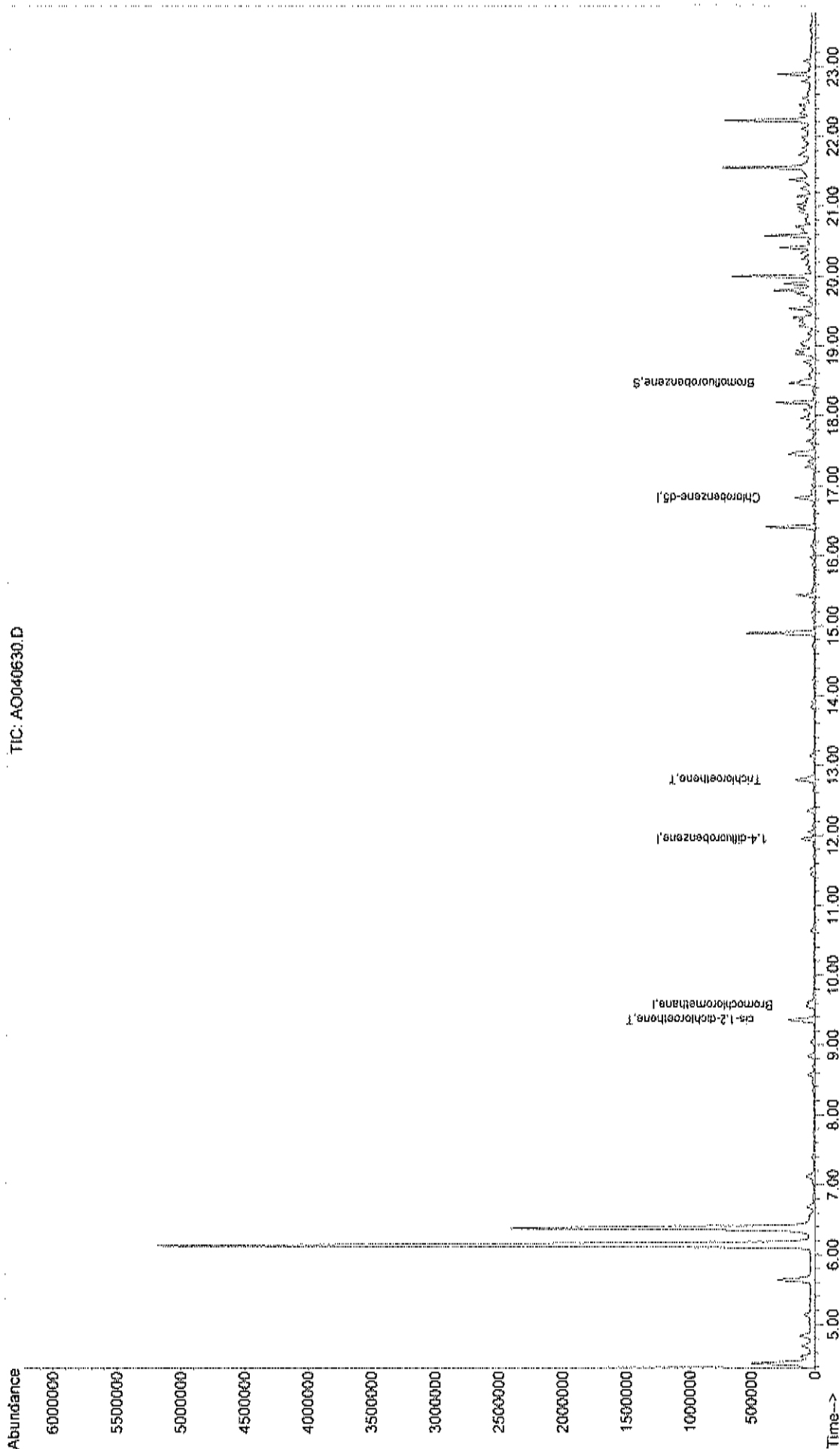
						Qvalue
29) cis-1,2-dichloroethene	9.36	61	184718	3.52	ppb	92
44) Trichloroethene	12.79	130	60313	1.21	ppb	88

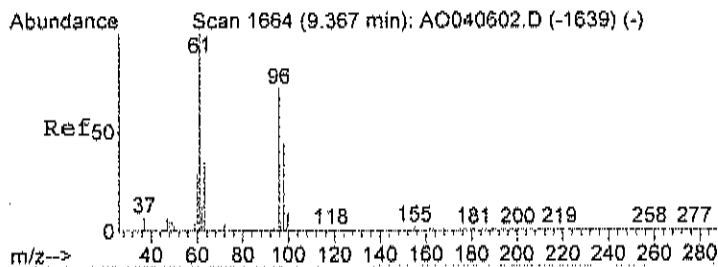
Data File : C:\HPCHEM\1\DATA\AO040630.D  
 Acq On : 7 Apr 2017 4:13 am  
 Sample : C1704014-012A  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 7 13:29 2017

Vial: 28  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

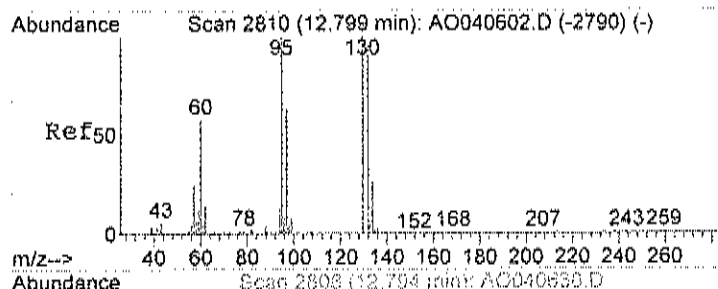
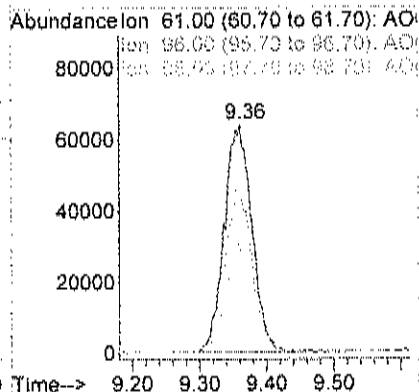
Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration





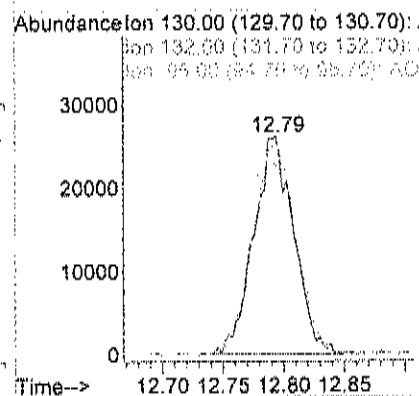
#29  
cis-1,2-dichloroethene  
Concen: 3.52 ppb  
RT: 9.36 min Scan# 1662  
Delta R.T. 0.01 min  
Lab File: AO040630.D  
Acq: 7 Apr 2017 4:13 am

Tgt Ion	Ratio	Lower	Upper
61	100		
96	70.4	58.1	98.1
98	45.4	29.3	69.3



#44  
Trichloroethene  
Concen: 1.21 ppb  
RT: 12.79 min Scan# 2808  
Delta R.T. 0.01 min  
Lab File: AO040630.D  
Acq: 7 Apr 2017 4:13 am

Tgt Ion	Ratio	Lower	Upper
130	100		
132	98.4	69.9	109.9
95	111.1	76.3	116.3



Data File : C:\HPCHEM\1\DATA\AO040706.D

Vial: 6

Acq On : 7 Apr 2017 1:50 pm

Operator: RJP

Sample : C1704014-012A 10X

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 07 14:58:05 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	20808	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.94	114	86837	1.00	ppb	0.00
50) Chlorobenzene-d5	16.82	117	77606	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	48257	0.91	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	91.00%

## Target Compounds

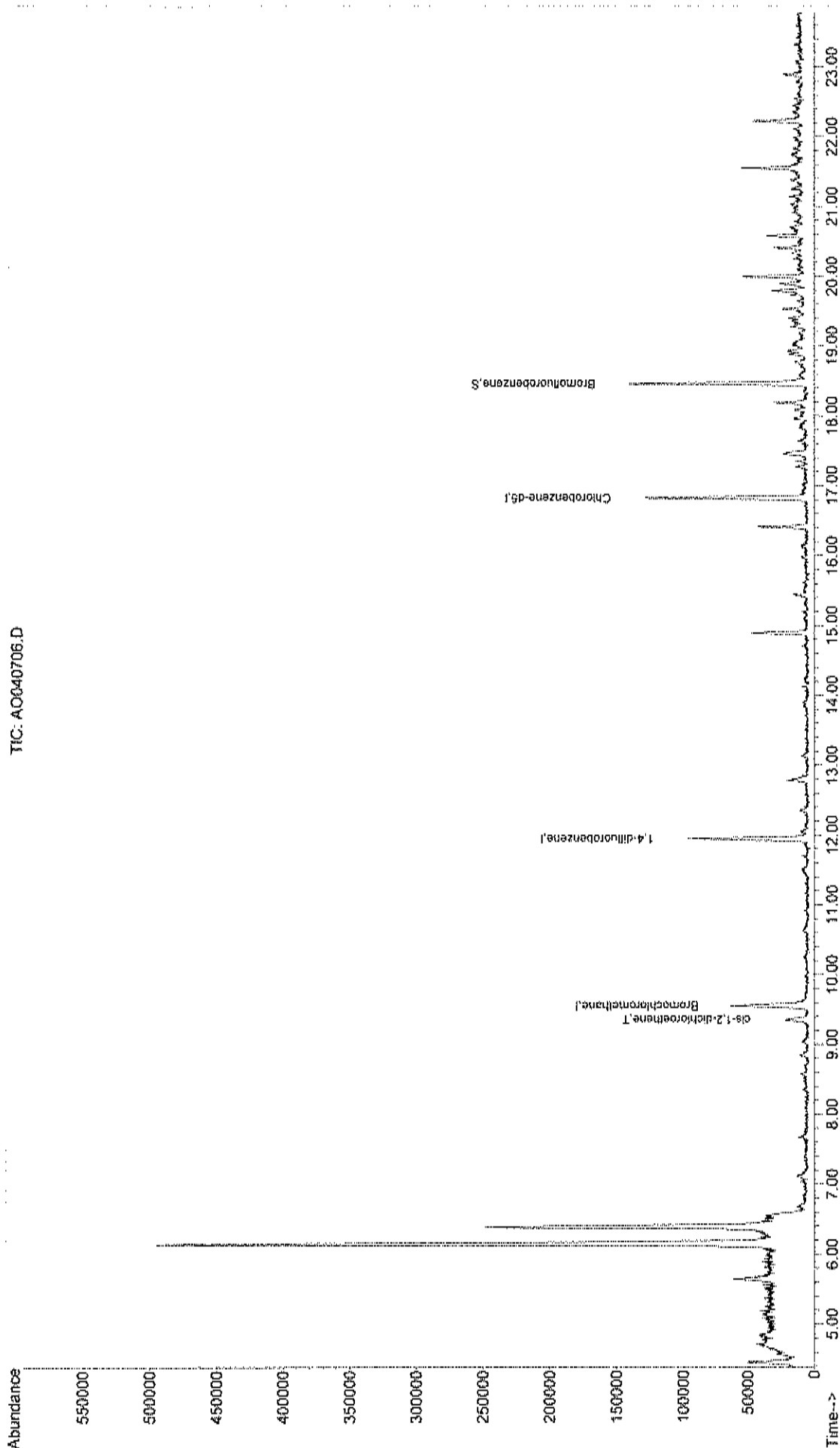
						Qvalue
29) cis-1,2-dichloroethene	9.35	61	14150	0.33	ppb	91

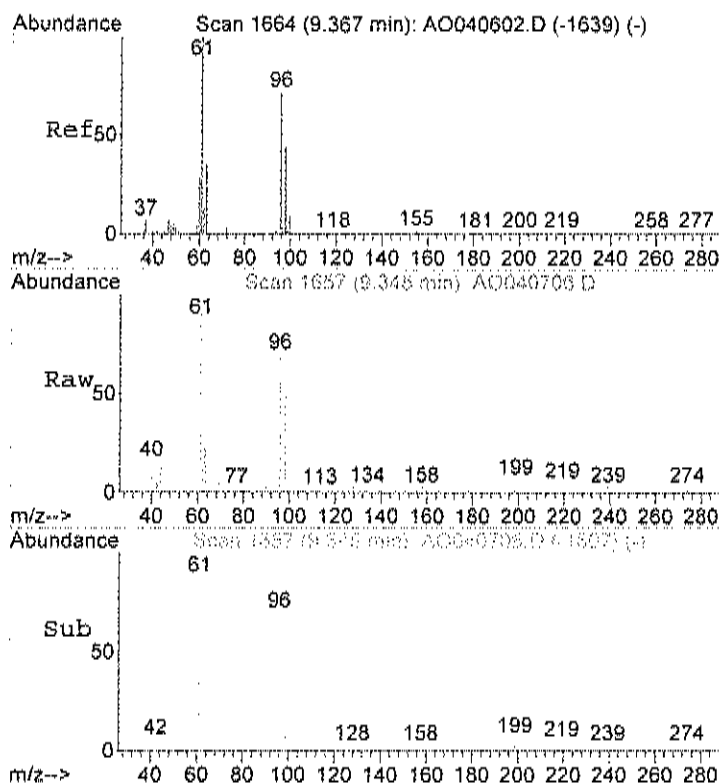
Data File : C:\HPCHEM\1\DATA\A0040706.D  
 Acq On : 7 Apr 2017 1:50 pm  
 Sample : C1704014-012A 10X  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 11 14:56 2017

Vial: 6  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

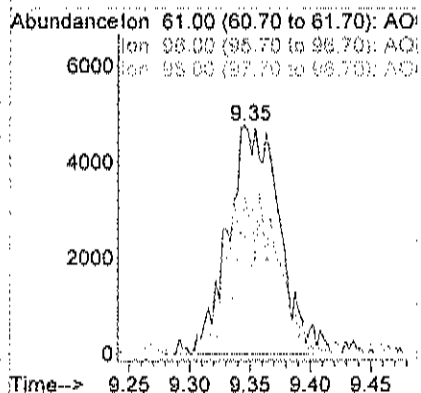
Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration





#29  
 cis-1,2-dichloroethene  
 Concen: 0.33 ppb  
 RT: 9.35 min Scan# 1657  
 Delta R.T. -0.00 min  
 Lab File: AO040706.D  
 Acq: 7 Apr 2017 1:50 pm

Tgt Ion	61	Resp	14150
Ion	Ratio	Lower	Upper
61	100		
96	67.0	58.1	98.1
98	47.3	29.3	69.3





# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-013A

Client Sample ID: 691-SB5B-1AQ  
Tag Number: 285.344  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-3			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 4:53:00 AM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 4:53:00 AM
cis-1,2-Dichloroethene	3.1	1.5		ppbV	10	4/7/2017 2:27:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 4:53:00 AM
Trichloroethene	1.3	0.040		ppbV	1	4/7/2017 4:53:00 AM
Vinyl chloride	0.060	0.040		ppbV	1	4/7/2017 4:53:00 AM
Surr: Bromofluorobenzene	102	70-130		%REC	1	4/7/2017 4:53:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
/ Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-013A

Client Sample ID: 691-SB5B-1AQ  
Tag Number: 285.344  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:53:00 AM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 4:53:00 AM
cis-1,2-Dichloroethene	12	5.9		ug/m3	10	4/7/2017 2:27:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 4:53:00 AM
Trichloroethene	6.8	0.21		ug/m3	1	4/7/2017 4:53:00 AM
Vinyl chloride	0.15	0.10		ug/m3	1	4/7/2017 4:53:00 AM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

Data File : C:\HPCHEM\1\DATA\AO040631.D

Vial: 29

Acq On : 7 Apr 2017 4:53 am

Operator: RJP

Sample : C1704014-013A

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 07 07:19:11 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	25079	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	113761	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	101873	1.00	ppb	0.01

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	70702	1.02	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	102.00%

## Target Compounds

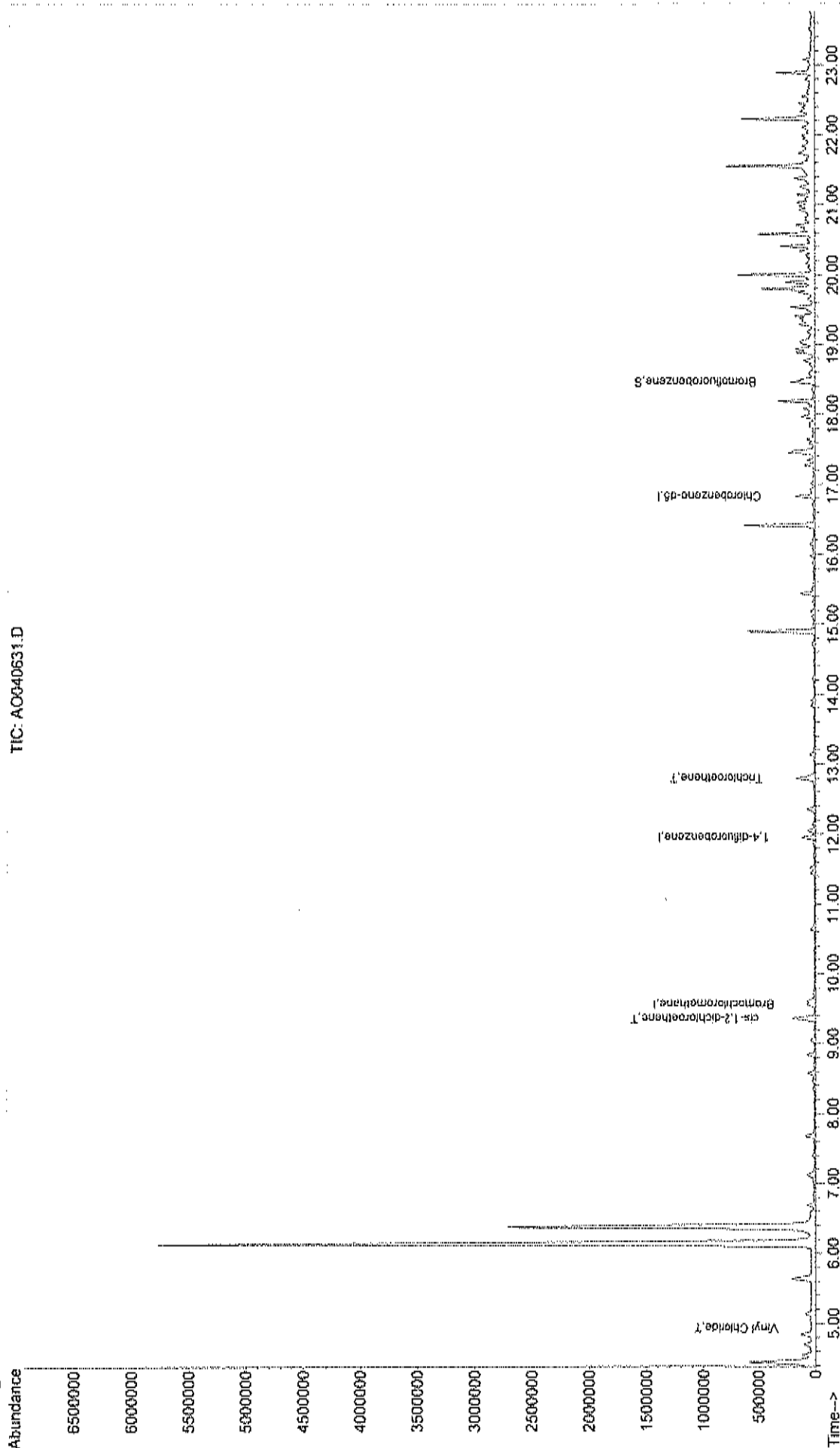
						Qvalue
6) Vinyl Chloride	4.94	62	2495	0.06	ppb	70
29) cis-1,2-dichloroethene	9.36	61	175403	3.37	ppb	94
44) Trichloroethene	12.79	130	65059	1.27	ppb	92

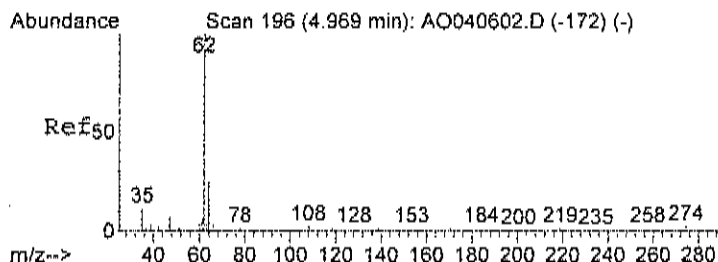
Data File : C:\HPCHEM\1\DATA\A0040631.D  
 Acq On : 7 Apr 2017 4:53 am  
 Sample : C1704014-013A  
 Misc : A331\_IUG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 7 13:31 2017

Vial: 29  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_IUG.RES

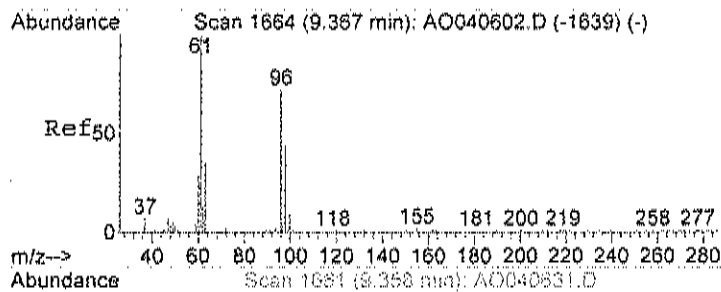
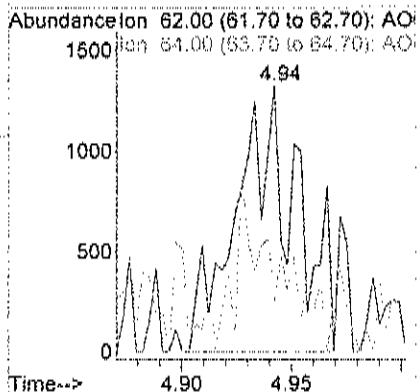
Method : C:\HPCHEM\1\METHODS\A331\_IUG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration





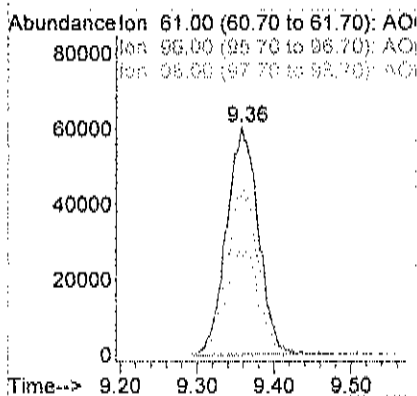
#6  
 Vinyl Chloride  
 Concen: 0.06 ppb  
 RT: 4.94 min Scan# 187  
 Delta R.T. -0.01 min  
 Lab File: AO040631.D  
 Acq: 7 Apr 2017 4:53 am

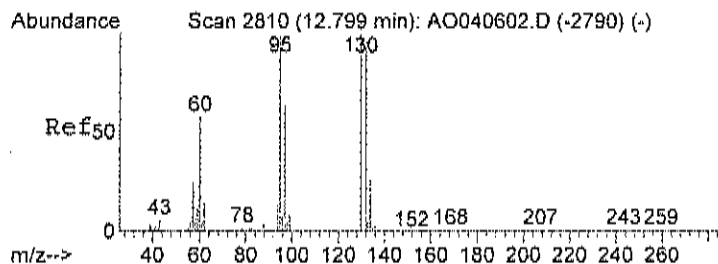
Tgt Ion: 62 Resp: 2495  
 Ion Ratio Lower Upper  
 62 100  
 64 44.7 0.0 58.6



#29  
 cis-1,2-dichloroethene  
 Concen: 3.37 ppb  
 RT: 9.36 min Scan# 1661  
 Delta R.T. 0.01 min  
 Lab File: AO040631.D  
 Acq: 7 Apr 2017 4:53 am

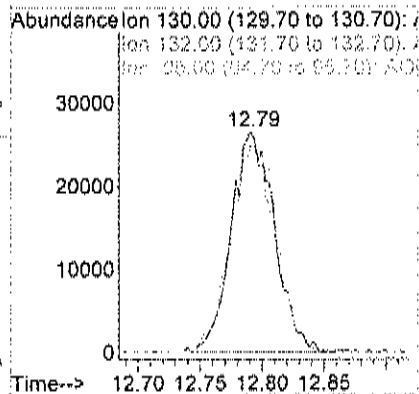
Tgt Ion: 61 Resp: 175403  
 Ion Ratio Lower Upper  
 61 100  
 96 72.2 58.1 98.1  
 98 46.4 29.3 69.3





#44  
 Trichloroethene  
 Concen: 1.27 ppb  
 RT: 12.79 min Scan# 2807  
 Delta R.T. 0.01 min  
 Lab File: AO040631.D  
 Acq: 7 Apr 2017 4:53 am

Tgt Ion	Ratio	Lower	Upper
130	100		
132	94.1	69.9	109.9
95	107.1	76.3	116.3





Data File : C:\HPCHEM\1\DATA\AO040707.D

Vial: 7

Acq On : 7 Apr 2017 2:27 pm

Operator: RJP

Sample : C1704014-013A 10X

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 07 14:58:20 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	19443	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	85901	1.00	ppb	0.01
50) Chlorobenzene-d5	16.82	117	76902	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	48147	0.92	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	92.00%

## Target Compounds

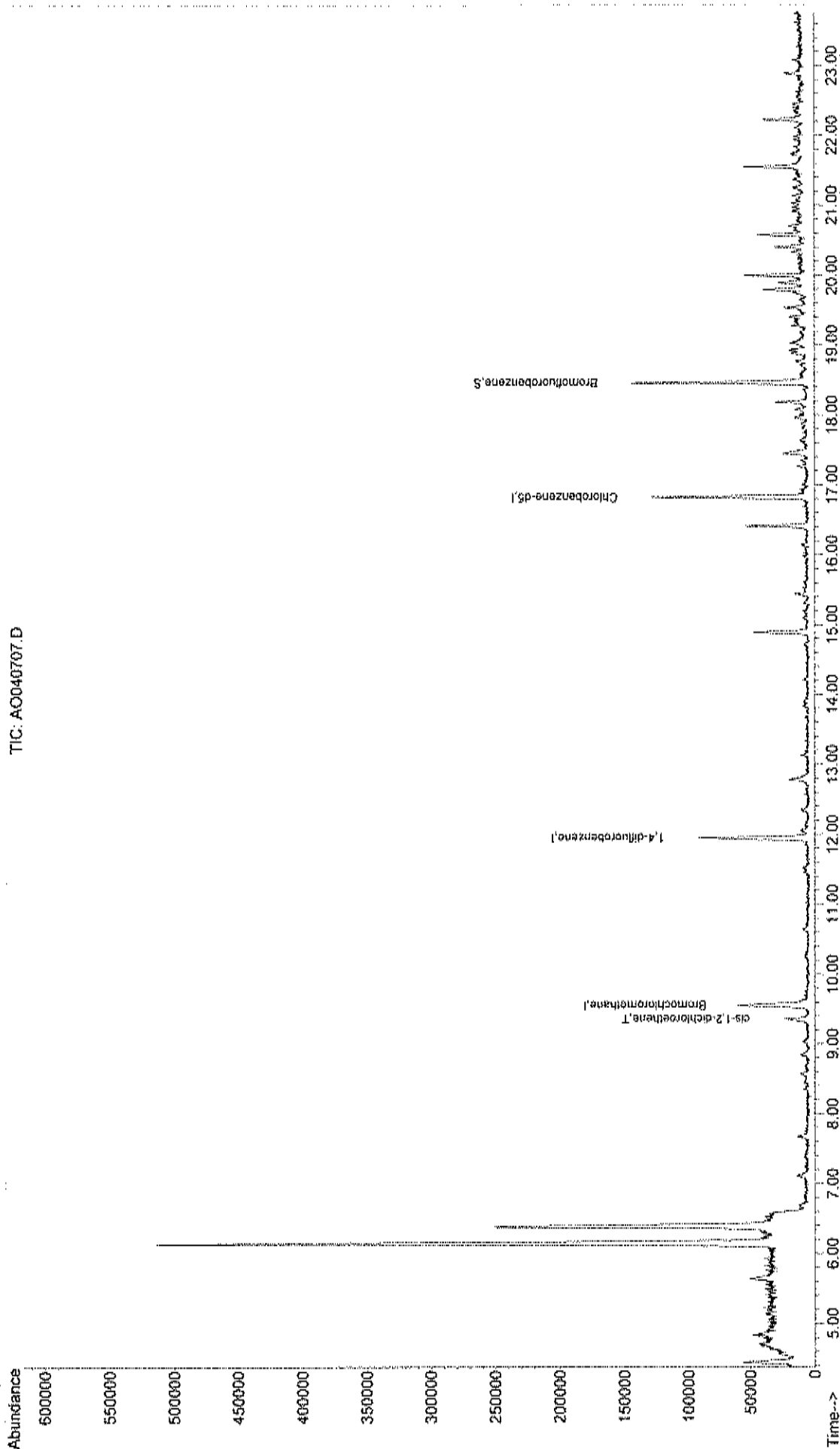
29) cis-1,2-dichloroethene	9.35	61	12708	0.31	ppb	Qvalue 96
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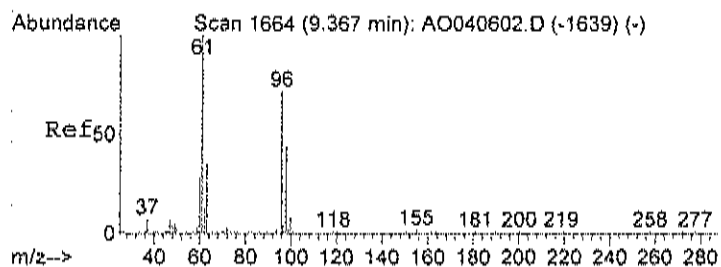
Data File : C:\HPCHEM\1\DATA\AO040707.D  
 Acq On : 7 Apr 2017 2:27 pm  
 Sample : C1704014-013A 10X  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 11 14:57 2017

Vial: 7  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

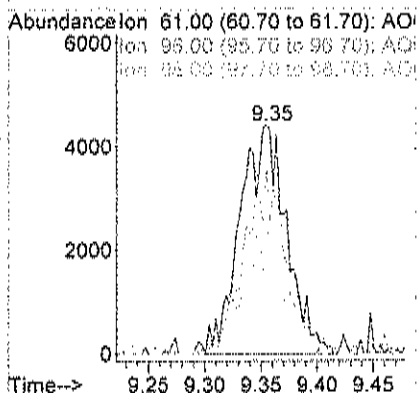
Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration





#29  
 cis-1,2-dichloroethene  
 Concen: 0.31 ppb  
 RT: 9.35 min Scan# 1660  
 Delta R.T. 0.01 min  
 Lab File: AO040707.D  
 Acq: 7 Apr 2017 2:27 pm

Tgt Ion	Ratio	Lower	Upper
61	100		
96	73.9	58.1	98.1
98	47.4	29.3	69.3



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-014A

Client Sample ID: 691-SB5A-SVI  
Tag Number: 475.1170  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-2			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 6:50:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 6:50:00 PM
cis-1,2-Dichloroethene	44	14		ppbV	90	4/8/2017 1:00:00 AM
trans-1,2-Dichloroethene	0.41	0.15		ppbV	1	4/7/2017 6:50:00 PM
Trichloroethene	3.7	0.36		ppbV	9	4/8/2017 12:23:00 AM
Vinyl chloride	0.19	0.040		ppbV	1	4/7/2017 6:50:00 PM
Surr: Bromofluorobenzene	107	70-130		%REC	1	4/7/2017 6:50:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-014A

Client Sample ID: 691-SB5A-SVI  
Tag Number: 475.1170  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DP	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 6:50:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 6:50:00 PM
cis-1,2-Dichloroethene	170	55		ug/m3	90	4/8/2017 1:00:00 AM
trans-1,2-Dichloroethene	1.6	0.59		ug/m3	1	4/7/2017 6:50:00 PM
Trichloroethene	20	1.9		ug/m3	9	4/8/2017 12:23:00 AM
Vinyl chloride	0.49	0.10		ug/m3	1	4/7/2017 6:50:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

Data File : C:\HPCHEM\1\DATA\AO040712.D

Vial: 12

Acq On : 7 Apr 2017 6:50 pm

Operator: RJP

Sample : C1704014-014A

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 08 08:59:04 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	31361	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	139148	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	131462	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	96106	1.07	ppb	0.01
Spiked Amount	1.000	Range	70 - 130	Recovery	=	107.00%

## Target Compounds

						Qvalue
6) Vinyl Chloride	4.95	62	10310m	0.19	ppb	
24) trans-1,2-dichloroethene	8.15	61	27880	0.41	ppb	96
29) cis-1,2-dichloroethene	9.35	61	2758473	42.35	ppb	98
44) Trichloroethene	12.79	130	321836	5.14	ppb	92

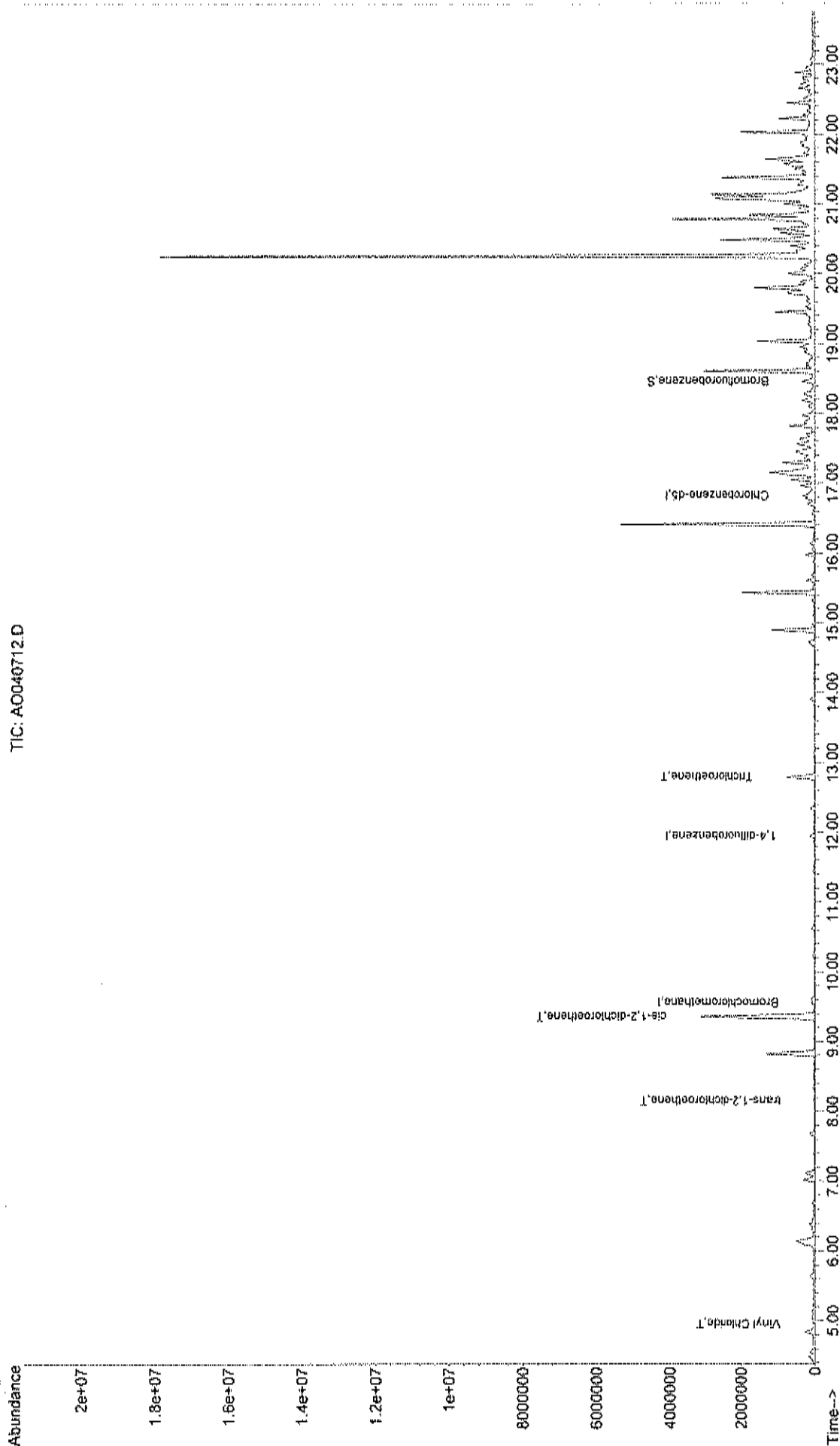


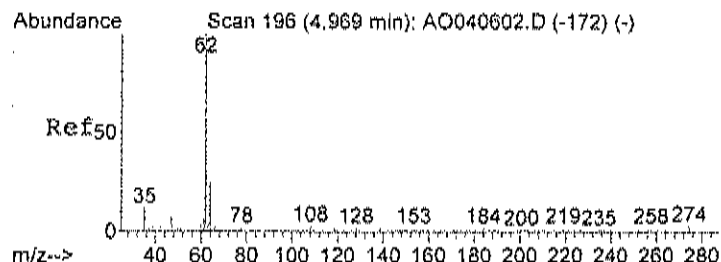
Data File : C:\HPCHEM\1\DATA\A0040712.D  
Acq On : 7 Apr 2017 6:50 pm  
Sample : C1704014-014A  
Misc : A331\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Apr 11 14:14 2017

Vial: 12  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A331\_1UG.RES

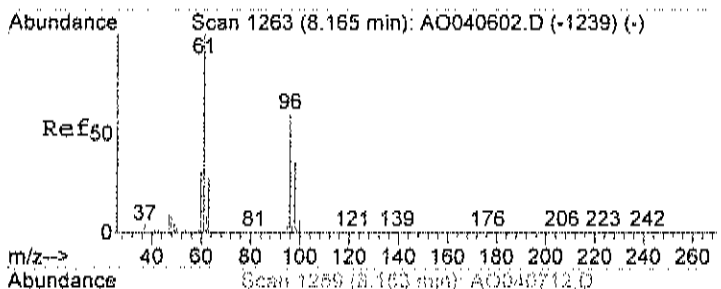
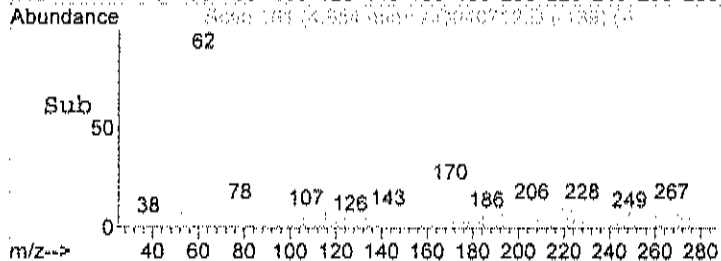
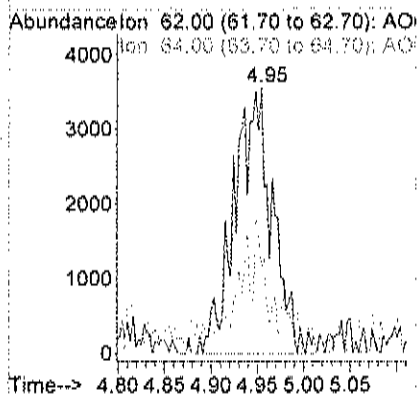
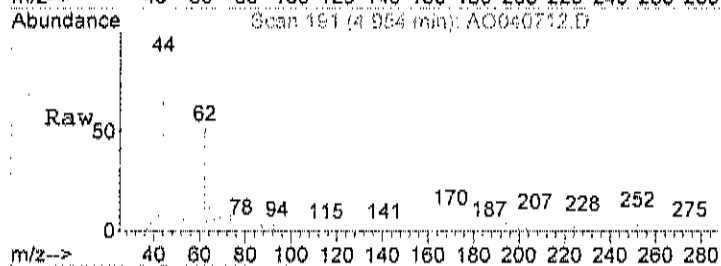
Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Thu May 04 11:27:28 2017  
Response via : Initial Calibration





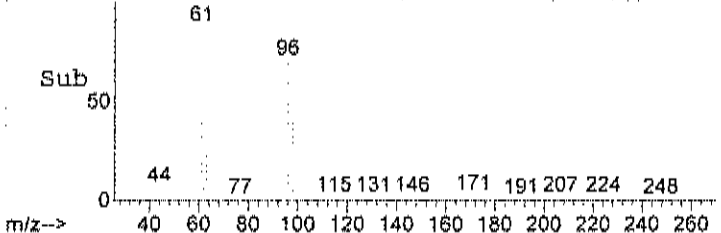
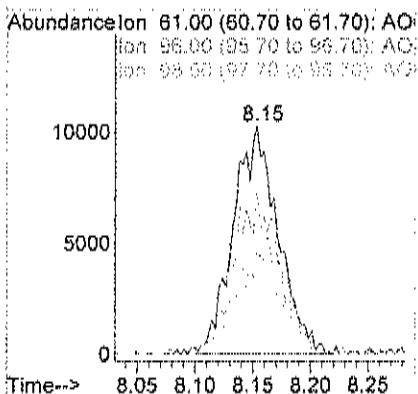
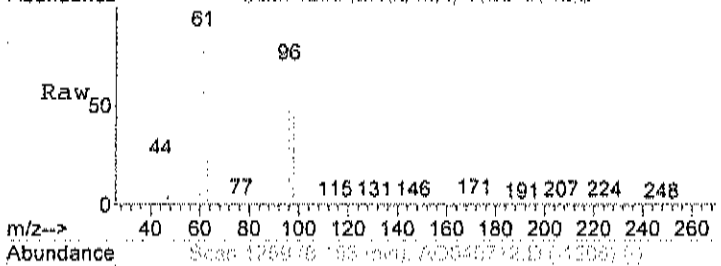
#6  
 Vinyl Chloride  
 Concen: 0.19 ppb m  
 RT: 4.95 min Scan# 191  
 Delta R.T. 0.01 min  
 Lab File: AO040712.D  
 Acq: 7 Apr 2017 6:50 pm

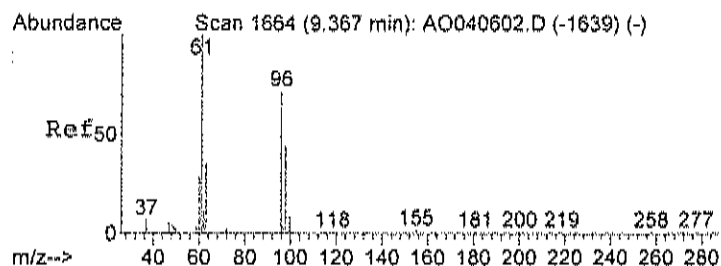
Tgt Ion: 62 Resp: 10310  
 Ion Ratio Lower Upper  
 62 100  
 64 45.0 0.0 58.6



#24  
 trans-1,2-dichloroethene  
 Concen: 0.41 ppb  
 RT: 8.15 min Scan# 1259  
 Delta R.T. 0.01 min  
 Lab File: AO040712.D  
 Acq: 7 Apr 2017 6:50 pm

Tgt Ion: 61 Resp: 27880  
 Ion Ratio Lower Upper  
 61 100  
 96 72.0 48.2 88.2  
 98 47.7 29.0 69.0

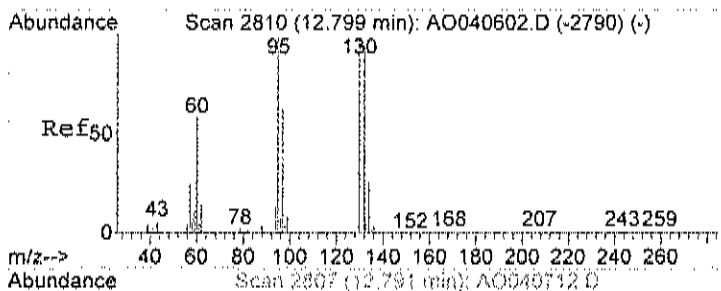
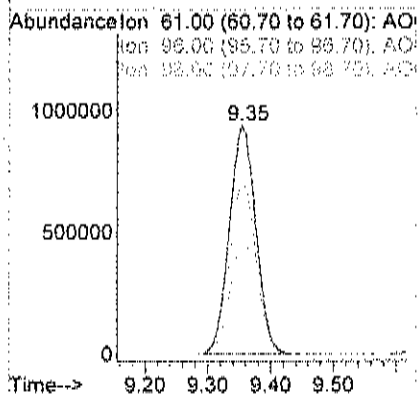




#29  
cis-1,2-dichloroethene  
Concen: 42.35 ppb  
RT: 9.35 min Scan# 1660  
Delta R.T. 0.01 min  
Lab File: AO040712.D  
Acq: 7 Apr 2017 6:50 pm

Tgt Ion: 61 Resp: 2758473

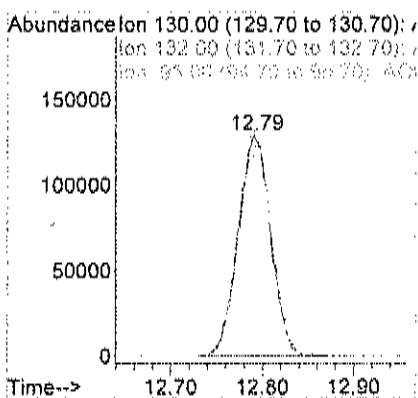
Ion	Ratio	Lower	Upper
61	100		
96	75.1	58.1	98.1
98	48.9	29.3	69.3



#44  
Trichloroethene  
Concen: 5.14 ppb  
RT: 12.79 min Scan# 2807  
Delta R.T. 0.01 min  
Lab File: AO040712.D  
Acq: 7 Apr 2017 6:50 pm

Tgt Ion: 130 Resp: 321836

Ion	Ratio	Lower	Upper
130	100		
132	96.4	69.9	109.9
95	104.9	76.3	116.3



Data File : C:\HPCHEM\1\DATA\AO040721.D

Vial: 21

Acq On : 8 Apr 2017 12:23 am

Operator: RJP

Sample : C1704014-014A 9x

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 08 08:59:56 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	20532	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	89989	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	83171	1.00	ppb	0.01

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	53731	0.95	ppb	0.01
Spiked Amount	1.000	Range	70 - 130	Recovery	=	95.00%

## Target Compounds

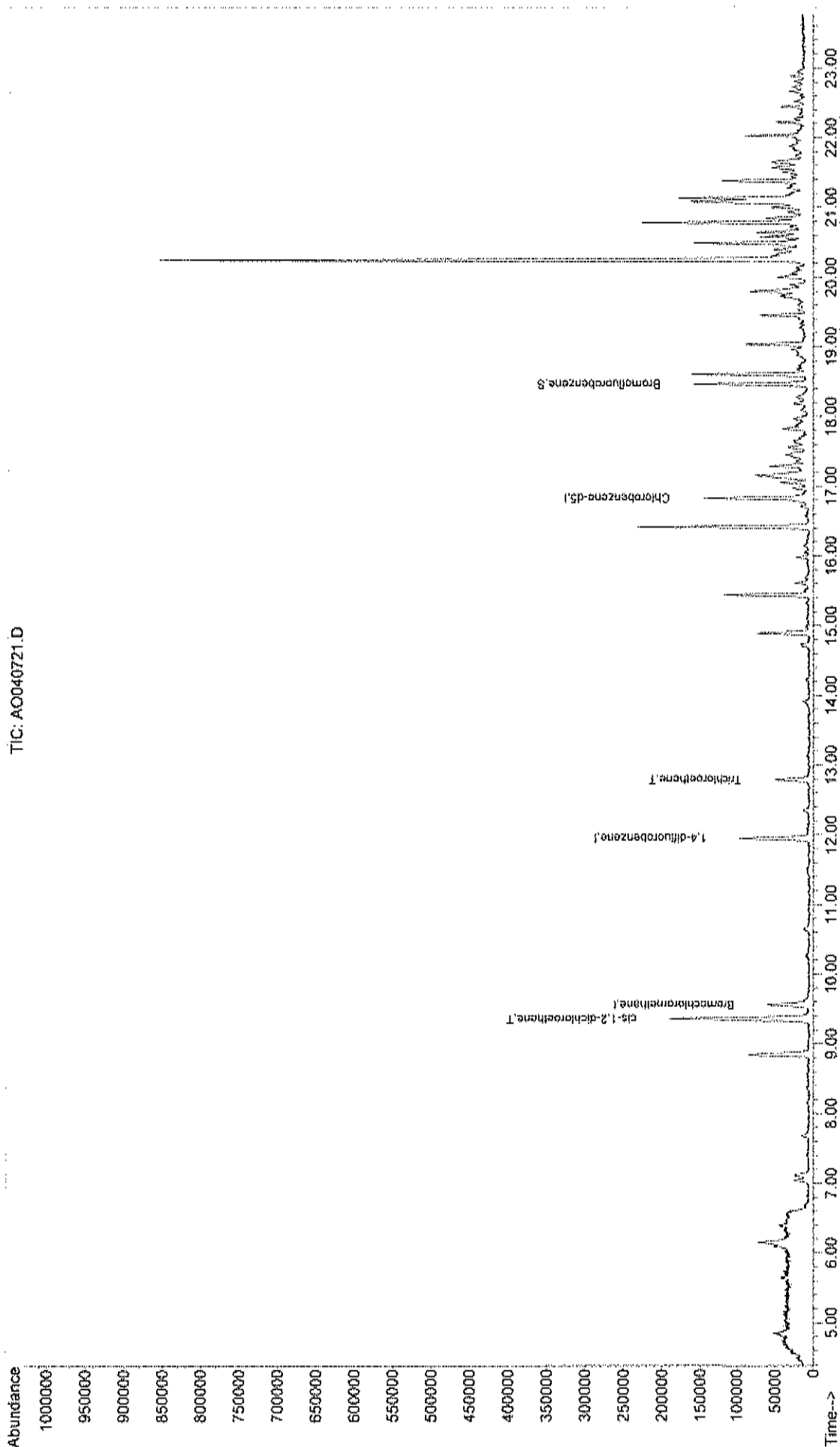
						Qvalue
29) cis-1,2-dichloroethene	9.35	61	156424	3.67	ppb	94
44) Trichloroethene	12.79	130	16628	0.41	ppb	94

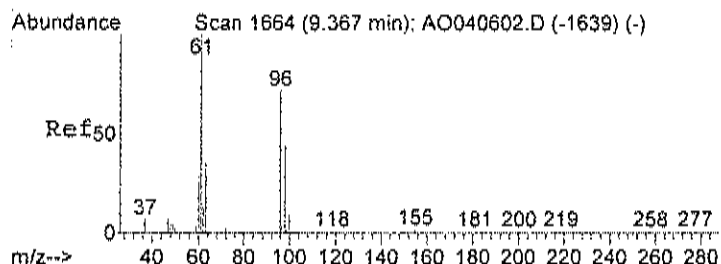
Data File : C:\HPCHEM\1\DATA\AO040721.D  
 Acq On : 8 Apr 2017 12:23 am  
 Sample : C1704014-014A 9x  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 11 15:01 2017

Vial: 21  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

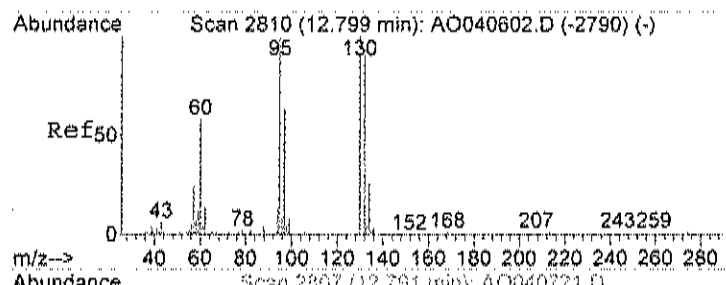
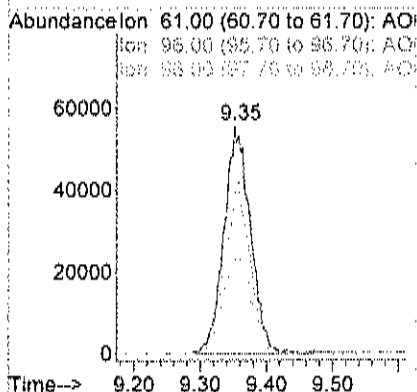
Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration





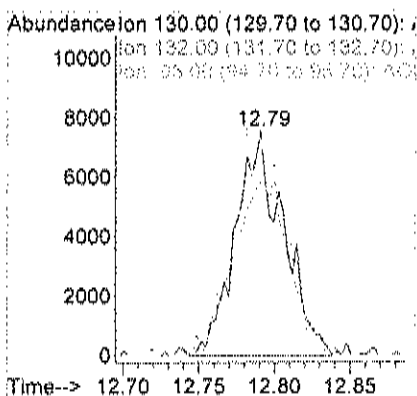
#29  
cis-1,2-dichloroethene  
Concen: 3.67 ppb  
RT: 9.35 min Scan# 1659  
Delta R.T. 0.01 min  
Lab File: AO040721.D  
Acq: 8 Apr 2017 12:23 am

Tgt Ion	Ratio	Lower	Upper
61	100		
96	71.9	58.1	98.1
98	46.2	29.3	69.3



#44  
Trichloroethene  
Concen: 0.41 ppb  
RT: 12.79 min Scan# 2807  
Delta R.T. 0.01 min  
Lab File: AO040721.D  
Acq: 8 Apr 2017 12:23 am

Tgt Ion	Ratio	Lower	Upper
130	100		
132	91.5	69.9	109.9
95	105.8	76.3	116.3





Data File : C:\HPCHEM\1\DATA\AO040722.D Vial: 22  
Acq On : 8 Apr 2017 1:00 am Operator: RJP  
Sample : C1704014-014A 90x Inst : MSD #1  
Misc : A331\_1UG Multiplr: 1.00  
MS Integration Params: RTEINT.P  
Quant Time: Apr 08 08:59:57 2017 Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Apr 03 10:15:59 2017  
Response via : Initial Calibration  
DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	20827	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.94	114	88198	1.00	ppb	0.00
50) Chlorobenzene-d5	16.83	117	80228	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	48670	0.89	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	89.00%

## Target Compounds

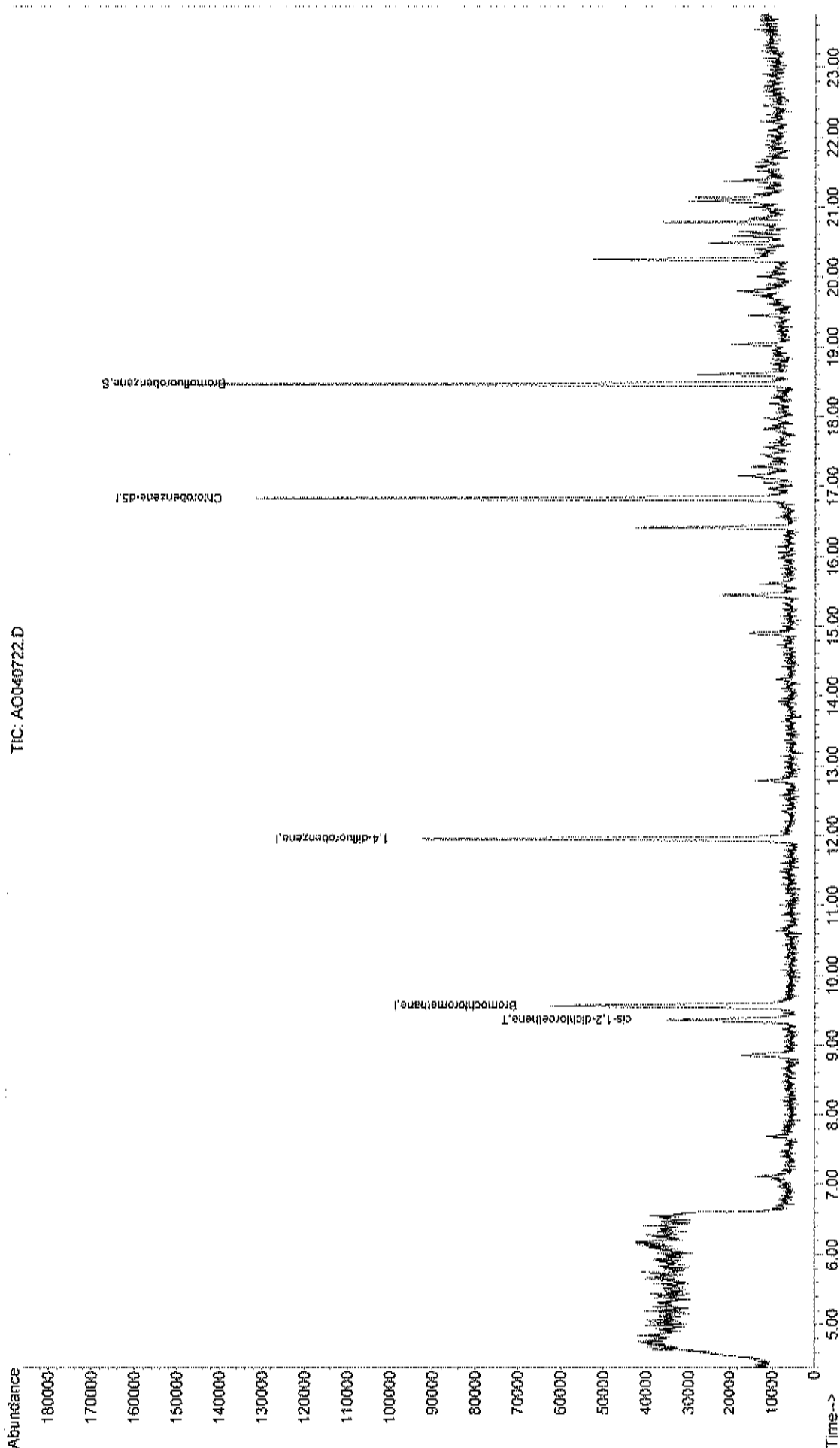
29) cis-1,2-dichloroethene	9.36	61	21151	0.49	ppb	Qvalue 96
----------------------------	------	----	-------	------	-----	-----------

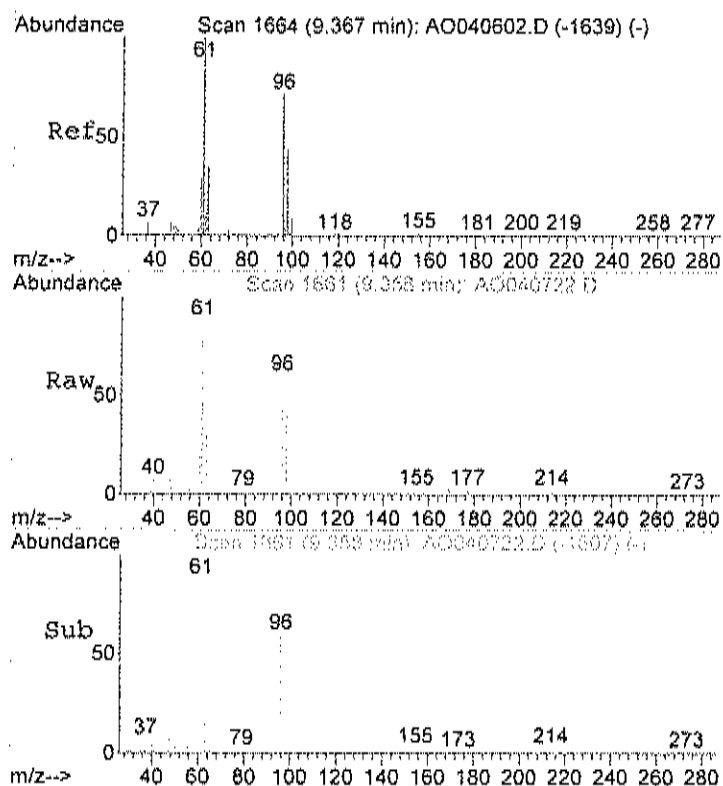
Data File : C:\HPCHEM\1\DATA\AO040722.D  
 Acq On : 8 Apr 2017 1:00 am  
 Sample : C1704014-014A 90x  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 11 15:02 2017

Vial: 22  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

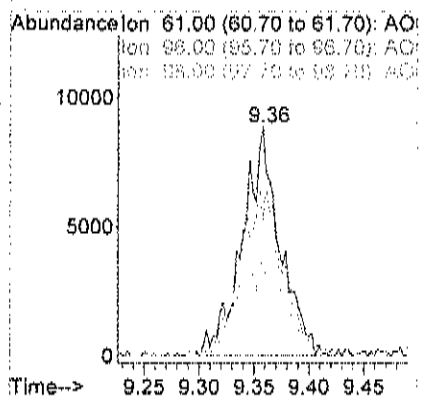
Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration





#29  
 cis-1,2-dichloroethene  
 Concen: 0.49 ppb  
 RT: 9.36 min Scan# 1661  
 Delta R.T. 0.01 min  
 Lab File: AO040722.D  
 Acq: 8 Apr 2017 1:00 am

Tgt Ion	Resp	Ion	Ratio	Lower	Upper
61	100	61	100		
96	83.8	96	83.8	58.1	98.1
98	48.9	98	48.9	29.3	69.3



# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-015A

Client Sample ID: 691-SB5B-SVI  
Tag Number: 467.1167  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-1			"Hg		4/5/2017
Lab Vacuum Out	-30			"Hg		4/5/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	4/7/2017 7:30:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	4/7/2017 7:30:00 PM
cis-1,2-Dichloroethene	5.8	1.5		ppbV	10	4/8/2017 1:37:00 AM
trans-1,2-Dichloroethene	0.18	0.15		ppbV	1	4/7/2017 7:30:00 PM
Trichloroethene	11	0.40		ppbV	10	4/8/2017 1:37:00 AM
Vinyl chloride	0.11	0.040		ppbV	1	4/7/2017 7:30:00 PM
Surr: Bromofluorobenzene	116	70-130		%REC	1	4/7/2017 7:30:00 PM

Qualifiers: \*\* Quantitation Limit  
B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
JN Non-routine analyte. Quantitation estimated.  
S Spike Recovery outside accepted recovery limits  
Results reported are not blank corrected  
E Estimated Value above quantitation range  
J Analyte detected below quantitation limit  
ND Not Detected at the Limit of Detection

# Centek Laboratories, LLC

Date: 04-May-17

CLIENT: LaBella Associates, P.C.  
Lab Order: C1704014  
Project: 691 St Paul Street  
Lab ID: C1704014-015A

Client Sample ID: 691-SB5B-SV1  
Tag Number: 467.1167  
Collection Date: 4/1/2017  
Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	4/7/2017 7:30:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	4/7/2017 7:30:00 PM
cis-1,2-Dichloroethene	23	5.9		ug/m3	10	4/8/2017 1:37:00 AM
trans-1,2-Dichloroethene	0.71	0.59		ug/m3	1	4/7/2017 7:30:00 PM
Trichloroethene	57	2.1		ug/m3	10	4/8/2017 1:37:00 AM
Vinyl chloride	0.28	0.10		ug/m3	1	4/7/2017 7:30:00 PM

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

Data File : C:\HPCHEM\1\DATA\AO040713.D

Vial: 13

Acq On : 7 Apr 2017 7:30 pm

Operator: RJP

Sample : C1704014-015A

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 08 08:59:20 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	31915	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	147799	1.00	ppb	0.02
50) Chlorobenzene-d5	16.83	117	139968	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.47	95	110484	1.16	ppb	0.02
Spiked Amount	1.000	Range	70 - 130	Recovery	=	116.00%

## Target Compounds

						Qvalue
6) Vinyl Chloride	4.95	62	6094	0.11	ppb	95
24) trans-1,2-dichloroethene	8.15	61	12259	0.18	ppb	94
29) cis-1,2-dichloroethene	9.35	61	537761	8.11	ppb	98
44) Trichloroethene	12.79	130	949109	14.27	ppb	92

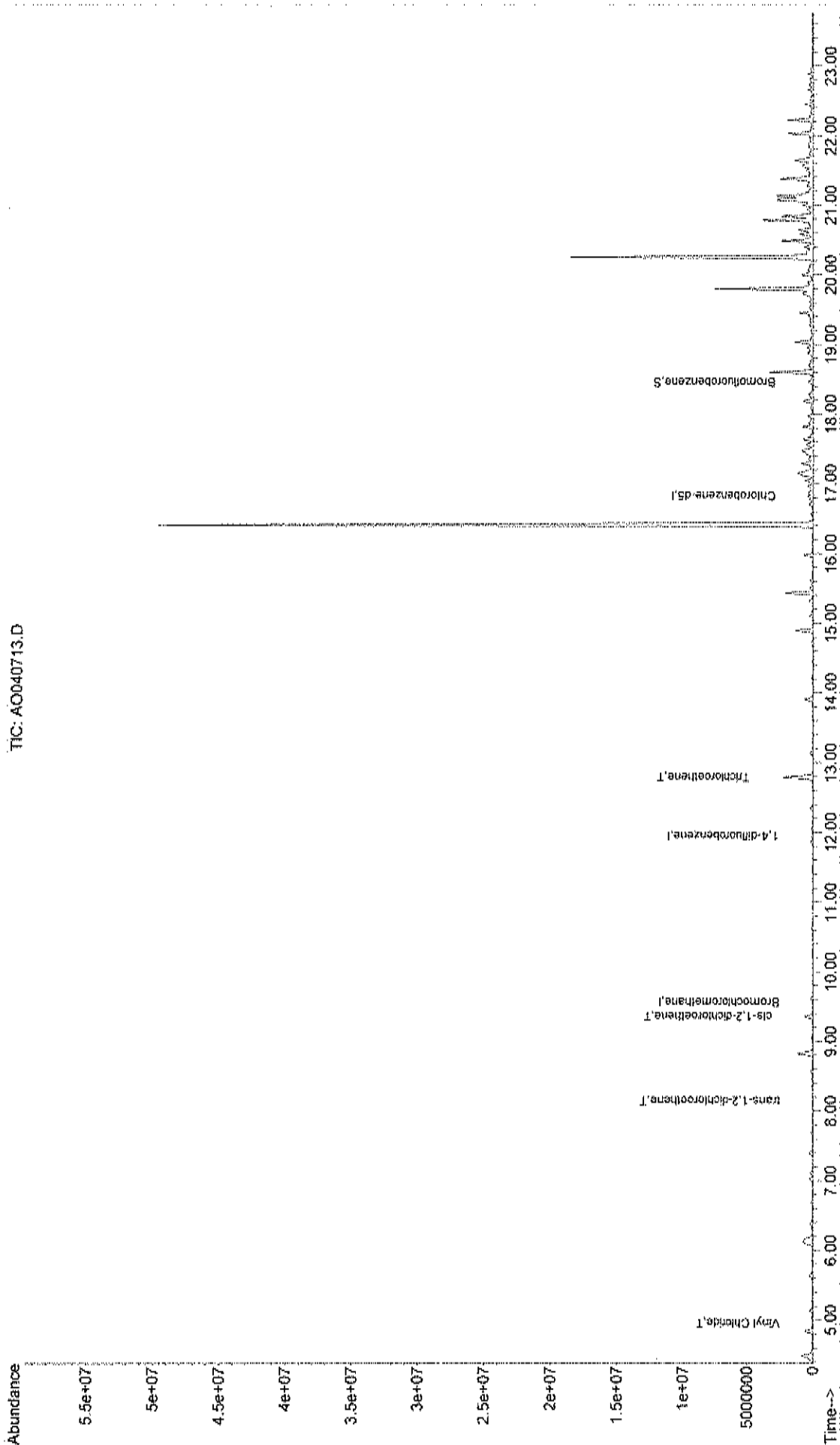


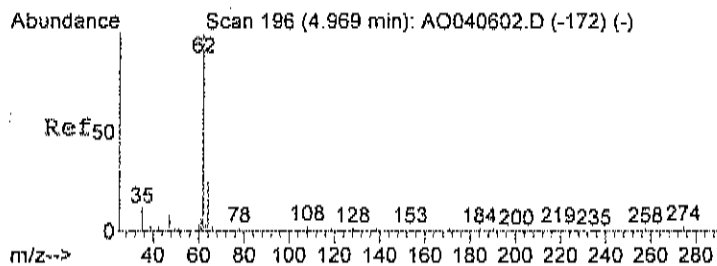
Data File : C:\HPCHEM\1\DATA\AO040713.D  
 Acq On : 7 Apr 2017 7:30 pm  
 Sample : C1704014-015A  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 11 14:16 2017

Vial: 13  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

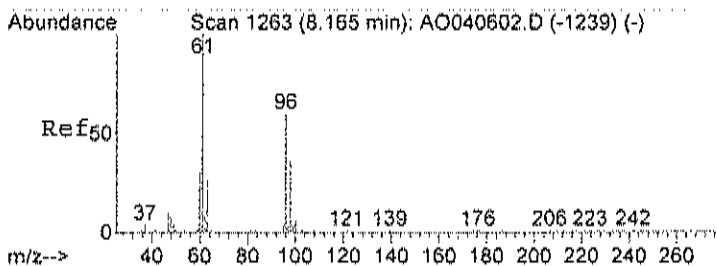
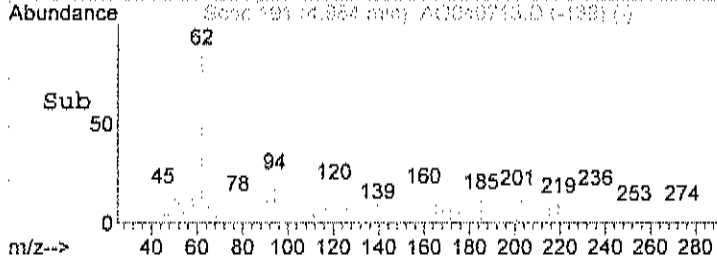
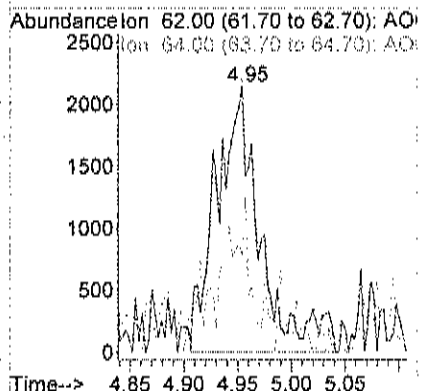
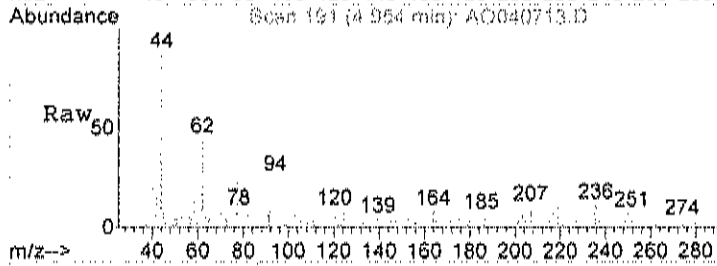
Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration





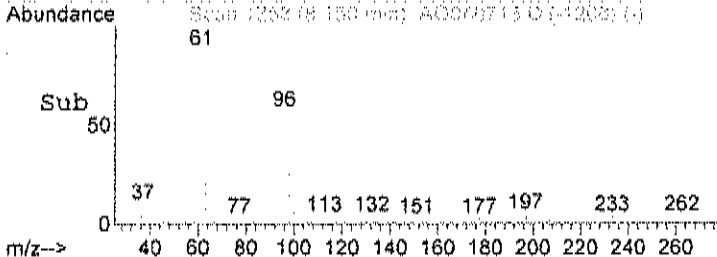
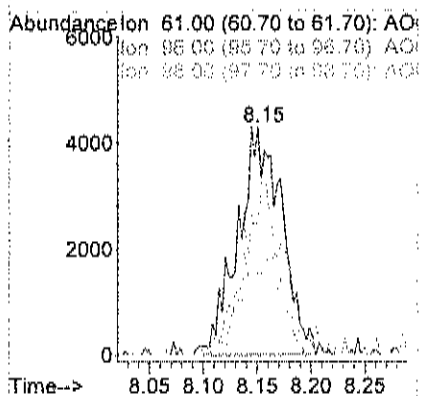
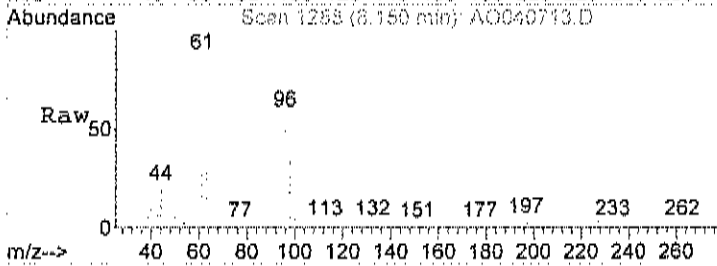
#6  
 Vinyl Chloride  
 Concen: 0.11 ppb  
 RT: 4.95 min Scan# 191  
 Delta R.T. 0.01 min  
 Lab File: AO040713.D  
 Acq: 7 Apr 2017 7:30 pm

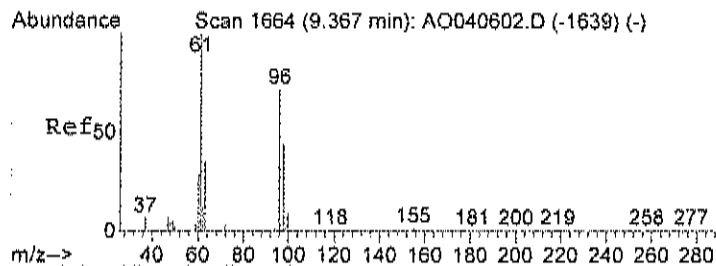
Tgt Ion: 62 Resp: 6094  
 Ion Ratio Lower Upper  
 62 100  
 64 31.0 0.0 58.6



#24  
 trans-1,2-dichloroethene  
 Concen: 0.18 ppb  
 RT: 8.15 min Scan# 1258  
 Delta R.T. 0.01 min  
 Lab File: AO040713.D  
 Acq: 7 Apr 2017 7:30 pm

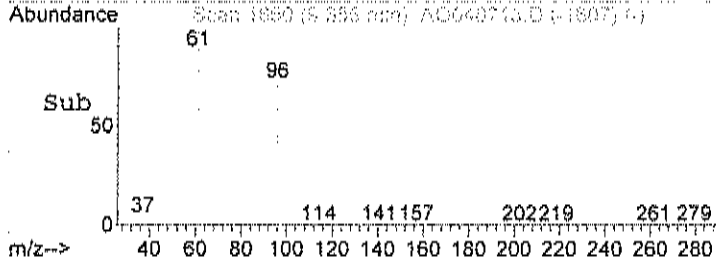
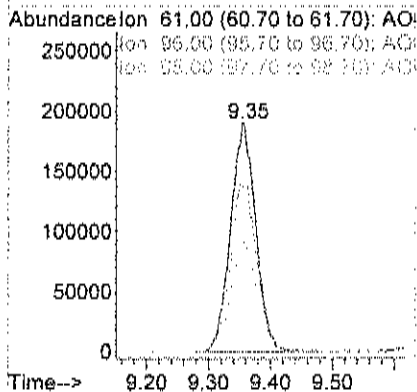
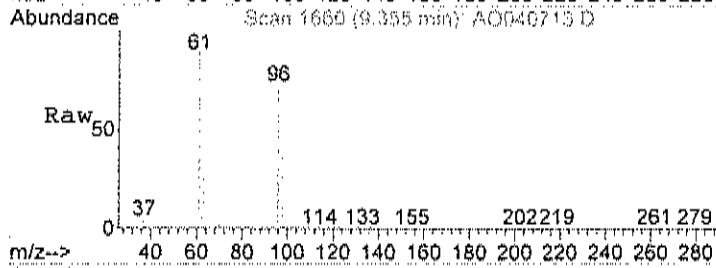
Tgt Ion: 61 Resp: 12259  
 Ion Ratio Lower Upper  
 61 100  
 96 65.3 48.2 88.2  
 98 42.2 29.0 69.0





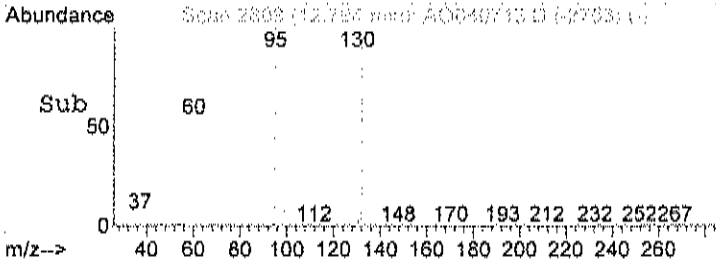
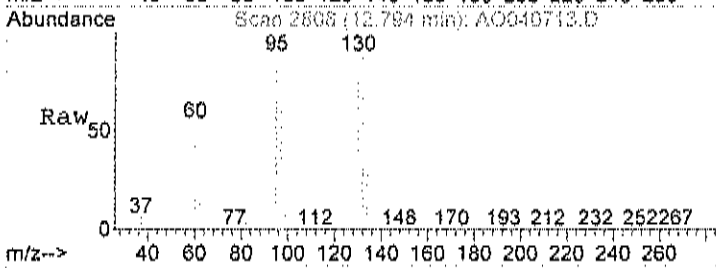
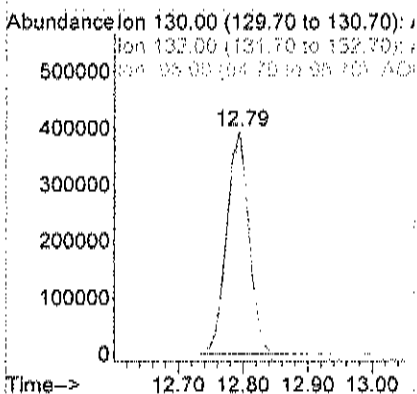
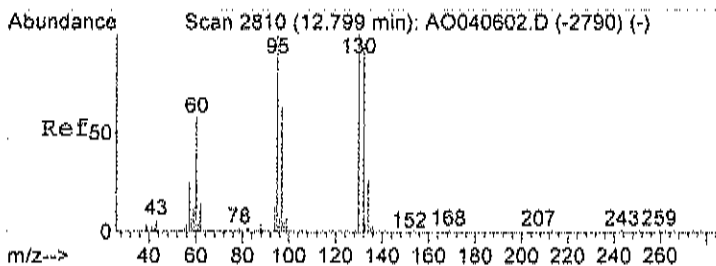
#29  
cis-1,2-dichloroethene  
Concen: 8.11 ppb  
RT: 9.35 min Scan# 1660  
Delta R.T. 0.01 min  
Lab File: AO040713.D  
Acq: 7 Apr 2017 7:30 pm

Tgt Ion	Ratio	Lower	Upper
61	100		
96	76.0	58.1	98.1
98	49.0	29.3	69.3



#44  
Trichloroethene  
Concen: 14.27 ppb  
RT: 12.79 min Scan# 2808  
Delta R.T. 0.01 min  
Lab File: AO040713.D  
Acq: 7 Apr 2017 7:30 pm

Tgt Ion	Ratio	Lower	Upper
130	100		
132	96.1	69.9	109.9
95	104.5	76.3	116.3



Data File : C:\HPCHEM\1\DATA\AO040723.D

Vial: 23

Acq On : 8 Apr 2017 1:37 am

Operator: RJP

Sample : C1704014-015A 10x

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 08 08:59:58 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.57	128	20780	1.00	ppb	0.03
35) 1,4-difluorobenzene	11.95	114	88679	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	79938	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	51379	0.94	ppb	0.01
Spiked Amount	1.000	Range	70 - 130	Recovery	=	94.00%

## Target Compounds

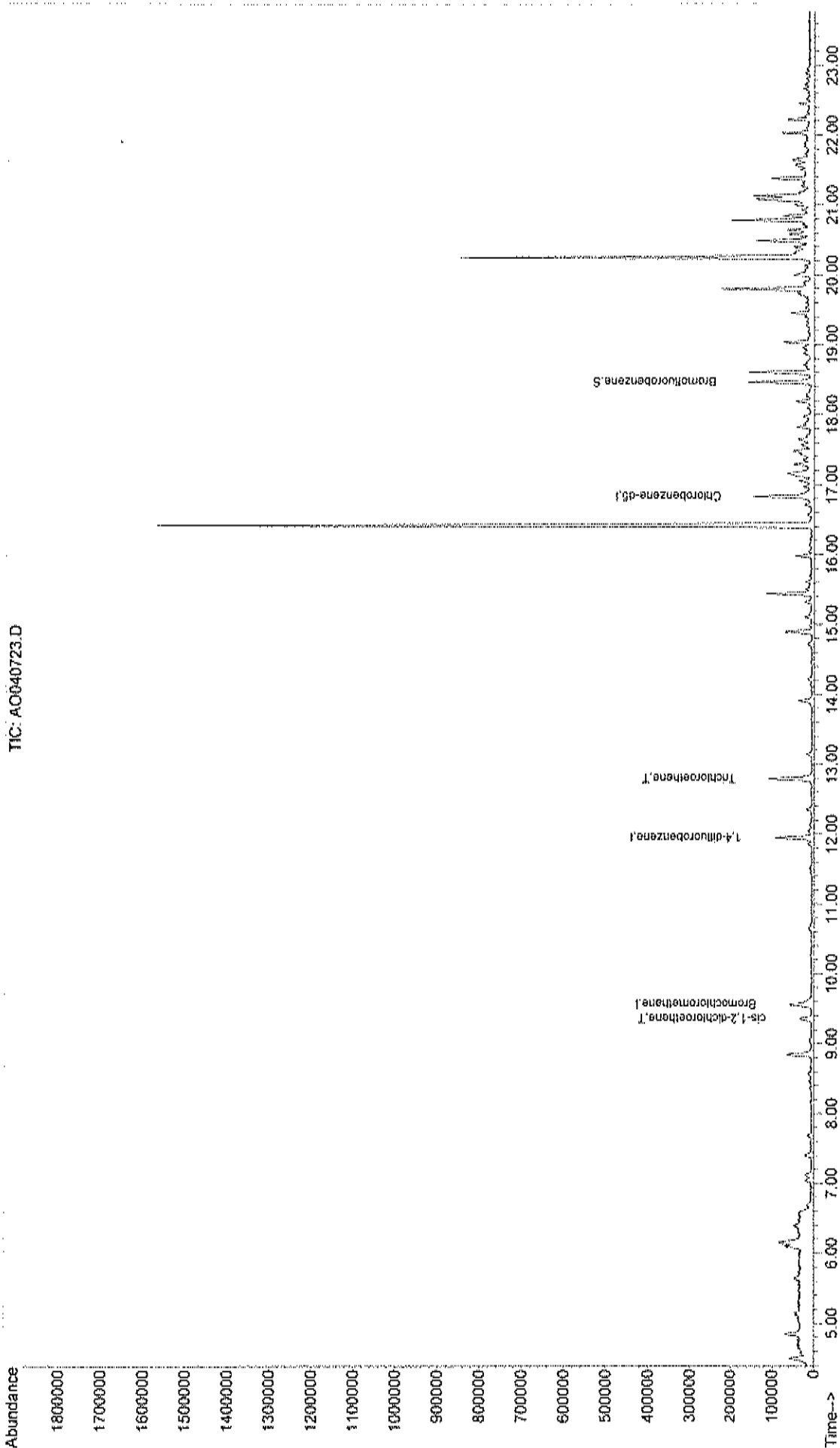
						Qvalue
29) cis-1,2-dichloroethene	9.35	61	25038	0.58	ppb	97
44) Trichloroethene	12.79	130	42344	1.06	ppb	94

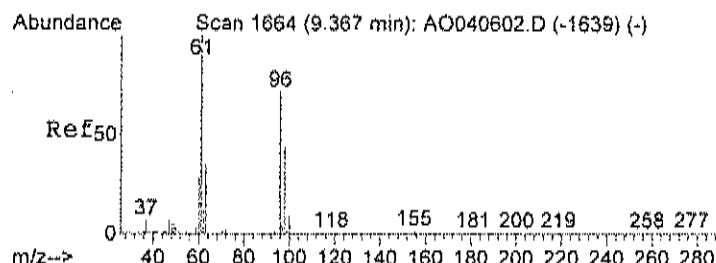
Data File : C:\HPCHEM\1\DATA\AO040723.D  
 Acq On : 8 Apr 2017 1:37 am  
 Sample : C1704014-015A 10x  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 11 15:03 2017

Vial: 23  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

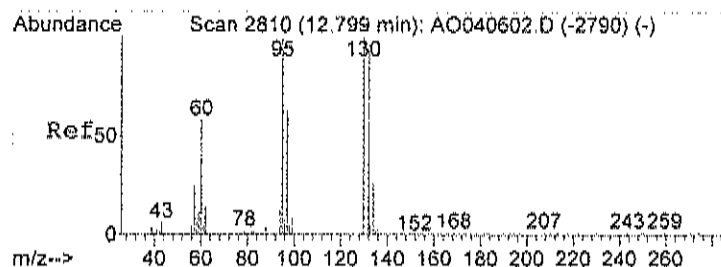
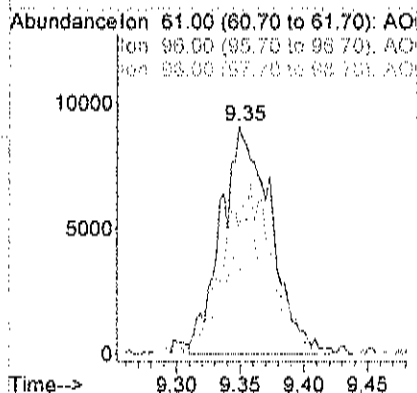
Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration





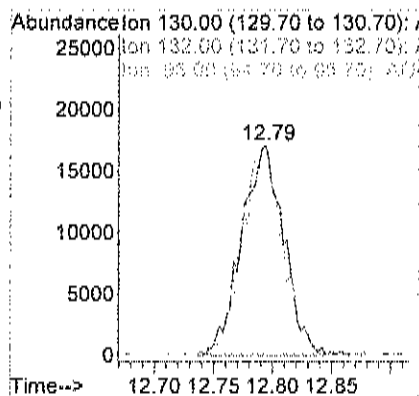
#29  
cis-1,2-dichloroethene  
Concen: 0.58 ppb  
RT: 9.35 min Scan# 1658  
Delta R.T. 0.00 min  
Lab File: AO040723.D  
Acq: 8 Apr 2017 1:37 am

Tgt Ion: 61 Resp: 25038  
Ion Ratio Lower Upper  
61 100  
96 73.3 58.1 98.1  
98 49.2 29.3 69.3



#44  
Trichloroethene  
Concen: 1.06 ppb  
RT: 12.79 min Scan# 2808  
Delta R.T. 0.02 min  
Lab File: AO040723.D  
Acq: 8 Apr 2017 1:37 am

Tgt Ion: 130 Resp: 42344  
Ion Ratio Lower Upper  
130 100  
132 93.3 69.9 109.9  
95 104.5 76.3 116.3





**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**STANDARDS DATA**

**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**INITIAL CALIBRATION**

## Response Factor Report MSD #1

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:13:15 2017  
 Response via : Initial Calibration

## Calibration Files

0.04 =AO033112.D 0.10 =AO033111.D 0.15 =AO033110.D  
 0.30 =AO033109.D 0.50 =AO033108.D 0.75 =AO033107.D

Compound		0.04	0.10	0.15	0.30	0.50	0.75	Avg	%RSD
-----									
1) I	Bromochloromethane	-----ISTD-----							
2) T	Propylene			1.544	1.513	1.429	1.374	1.425	4.95
3) T	Freon 12			9.344	8.388	8.315	8.416	8.232	6.60
4) T	Chloromethane			1.409	1.204	1.135	1.140	1.163	9.33
5) T	Freon 114			7.169	6.615	6.534	6.541	6.407	6.09
6) T	Vinyl Chloride	1.962	2.151	1.926	1.794	1.697	1.723	1.763	11.15
7) T	Butane			2.243	1.863	1.888	1.759	1.798	11.56
8) T	1,3-butadiene			1.385	1.332	1.172	1.185	1.165	12.44
9) T	Bromomethane			2.616	2.776	2.328	2.341	2.375	8.93
10) T	Chloroethane			0.809	0.822	0.754	0.711	0.733	7.64
11) T	Ethanol			0.675	0.699	0.533	0.570	0.576	12.29
12) T	Acrolein			0.825	0.551	0.639	0.600	0.604	16.23
13) T	Vinyl Bromide			2.774	2.435	2.310	2.408	2.385	7.43
14) T	Freon 11			9.323	8.425	8.442	8.349	8.237	6.73
15) T	Acetone			0.812	0.684	0.652	0.612	0.645	11.74
16) T	Pentane			1.425	1.259	1.237	1.280	1.262	5.84
17) T	Isopropyl alcoh			2.299	2.338	2.046	2.193	2.143	5.78
18) T	1,1-dichloroeth			1.456	1.286	1.356	1.296	1.300	5.86
19) T	Freon 113			3.463	3.043	3.112	3.108	3.107	4.91
20) T	t-Butyl alcohol			3.473	3.294	2.983	3.062	3.171	4.82
21) T	Methylene chlor			1.418	1.293	1.316	1.229	1.268	5.89
22) T	Allyl chloride			1.815	1.621	1.533	1.658	1.615	5.58
23) T	Carbon disulfid			4.523	4.199	4.065	4.123	4.058	5.82
24) T	trans-1,2-dichl			2.380	2.240	2.149	2.167	2.189	4.11
25) T	methyl tert-but			4.329	4.263	4.015	4.265	4.148	3.02
26) T	1,1-dichloroeth			2.962	2.740	2.777	2.701	2.710	4.53
27) T	Vinyl acetate			4.485	3.391	3.300	3.441	3.495	11.70
28) T	Methyl Ethyl Ke			0.644	0.662	0.679	0.661	0.665	2.29
29) T	cis-1,2-dichlor			2.251	2.073	2.073	2.093	2.077	3.80
30) T	Hexane			1.947	1.942	1.910	1.981	1.907	2.73
31) T	Ethyl acetate			4.499	4.398	4.247	4.265	4.254	3.36
32) T	Chloroform			4.143	3.779	3.803	3.825	3.793	4.47
33) T	Tetrahydrofuran			1.517	1.442	1.411	1.445	1.421	3.36
34) T	1,2-dichloroeth			3.011	2.729	2.856	2.889	2.807	4.09
-----									
35) I	1,4-difluorobenzene	-----ISTD-----							
36) T	1,1,1-trichloro			0.965	0.932	0.867	0.850	0.880	4.95
37) T	Cyclohexane			0.447	0.521	0.424	0.446	0.444	7.93
38) T	Carbon tetrachl	1.015	1.032	1.008	0.916	0.883	0.885	0.932	6.53
39) T	Benzene			1.006	0.976	0.907	0.899	0.920	5.07
40) T	Methyl methacry			0.460	0.442	0.414	0.425	0.431	3.31
41) T	1,4-dioxane			0.212	0.220	0.173	0.171	0.190	9.21
42) T	2,2,4-trimethyl			1.489	1.363	1.366	1.380	1.383	3.42
43) T	Heptane			0.511	0.476	0.485	0.473	0.480	2.80
44) T	Trichloroethene	0.522	0.481	0.479	0.441	0.420	0.438	0.450	7.35
45) T	1,2-dichloropro			0.390	0.338	0.340	0.339	0.339	6.32
46) T	Bromodichlorome			0.935	0.878	0.868	0.876	0.873	3.18
47) T	cis-1,3-dichlor			0.623	0.564	0.542	0.535	0.559	4.98
48) T	trans-1,3-dichl			0.558	0.532	0.524	0.530	0.532	2.79
49) T	1,1,2-trichloro			0.427	0.425	0.407	0.399	0.407	3.32
-----									
50) I	Chlorobenzene-d5	-----ISTD-----							
51) T	Toluene			0.836	0.781	0.724	0.731	0.750	5.44

## Response Factor Report MSD #1

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:13:15 2017  
 Response via : Initial Calibration

## Calibration Files

0.04 =AO033112.D 0.10 =AO033111.D 0.15 =AO033110.D  
 0.30 =AO033109.D 0.50 =AO033108.D 0.75 =AO033107.D

	Compound	0.04	0.10	0.15	0.30	0.50	0.75	Avg	%RSD
52) T	Methyl Isobutyl			0.842	0.806	0.686	0.687	0.780	7.66
53) T	Dibromochlorome			0.996	0.909	0.905	0.897	0.917	3.54
54) T	Methyl Butyl Ke			0.769	0.697	0.572	0.590	0.702	11.13
55) T	1,2-dibromoetha			0.817	0.783	0.729	0.745	0.761	3.66
56) T	Tetrachloroethy		0.602	0.610	0.547	0.511	0.498	0.529	8.75
57) T	Chlorobenzene			1.150	1.055	1.013	1.034	1.044	4.31
58) T	Ethylbenzene			1.900	1.763	1.709	1.719	1.750	3.78
59) T	m&p-xylene			1.484	1.529	1.451	1.491	1.477	1.75
60) T	Nonane			0.859	0.833	0.787	0.794	0.818	2.84
61) T	Styrene			0.989	0.906	0.877	0.898	0.919	3.61
62) T	Bromoform			0.798	0.762	0.730	0.753	0.761	2.79
63) T	o-xylene			1.520	1.458	1.380	1.415	1.424	3.14
64) T	Cumene			2.030	1.935	1.787	1.817	1.851	4.65
65) S	Bromofluorobenz	0.658	0.652	0.688	0.674	0.670	0.685	0.682	3.13
66) T	1,1,2,2-tetrach			1.089	1.023	0.950	0.961	0.973	5.71
67) T	Propylbenzene			0.523	0.488	0.460	0.469	0.476	4.49
68) T	2-Chlorotoluene			0.467	0.480	0.433	0.424	0.440	4.85
69) T	4-ethyltoluene			1.807	1.724	1.685	1.688	1.729	2.22
70) T	1,3,5-trimethyl			1.560	1.598	1.507	1.507	1.538	2.14
71) T	1,2,4-trimethyl			1.437	1.474	1.391	1.418	1.434	1.91
72) T	1,3-dichloroben			0.840	0.891	0.842	0.922	0.902	4.60
73) T	benzyl chloride			0.688	0.683	0.675	0.714	0.755	10.12
74) T	1,4-dichloroben			0.809	0.797	0.798	0.837	0.834	4.08
75) T	1,2,3-trimethyl			1.517	1.458	1.401	1.409	1.443	2.58
76) T	1,2-dichloroben			0.890	0.863	0.835	0.860	0.869	2.82
77) T	1,2,4-trichloro			0.197	0.215	0.220	0.242	0.257	18.07
78) T	Naphthalene			0.504	0.604	0.546	0.603	0.661	18.21
79) T	Hexachloro-1,3~			0.652	0.675	0.607	0.641	0.644	3.53

Data File : C:\HPCHEM\1\DATA\AO033103.D  
 Acq On : 31 Mar 2017 3:35 pm  
 Sample : A1UG 2.0  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Mar 31 21:02:10 2017

Vial: 2  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Fri Mar 31 21:00:46 2017  
 Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D  
 DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	50190	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.93	114	230091	1.00	ppb	0.00
50) Chlorobenzene-d5	16.81	117	199884	1.00	ppb	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
65) Bromofluorobenzene	18.45	95	141811	1.02	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	102.00%

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Propylene	4.48	41	135733	1.89	ppb	86
3) Freon 12	4.55	85	771524	1.93	ppb	98
4) Chloromethane	4.71	50	104844m	1.87	ppb	
5) Freon 114	4.81	85	609896	1.97	ppb	97
6) Vinyl Chloride	4.94	62	156398	1.95	ppb	89
7) Butane	5.13	43	164187	2.02	ppb	91
8) 1,3-butadiene	5.07	39	103468m	1.80	ppb	
9) Bromomethane	5.35	94	217749	1.89	ppb	94
10) Chloroethane	5.52	64	67680	1.96	ppb	# 86
11) Ethanol	5.63	45	51068m	1.86	ppb	
12) Acrolein	5.99	56	55607m	1.85	ppb	
13) Vinyl Bromide	5.89	106	226013	1.92	ppb	93
14) Freon 11	6.31	101	757784	1.88	ppb	100
15) Acetone	6.13	58	57430	1.92	ppb	# 75
16) Pentane	6.68	42	120588	1.90	ppb	# 36
17) Isopropyl alcohol	6.37	45	202779	1.88	ppb	# 100
18) 1,1-dichloroethene	6.98	96	126324	2.06	ppb	90
19) Freon 113	7.38	101	301800	2.02	ppb	91
20) t-Butyl alcohol	7.01	59	316399	2.00	ppb	98
21) Methylene chloride	7.10	84	118066	1.91	ppb	# 77
22) Allyl chloride	7.23	41	156544	1.98	ppb	91
23) Carbon disulfide	7.44	76	383855	1.96	ppb	84
24) trans-1,2-dichloroethene	8.13	61	217211	2.07	ppb	90
25) methyl tert-butyl ether	8.43	73	407448	2.00	ppb	92
26) 1,1-dichloroethane	8.37	63	259676	1.95	ppb	94
27) Vinyl acetate	8.52	43	333912m	2.03	ppb	
28) Methyl Ethyl Ketone	8.82	72	65520	1.90	ppb	100
29) cis-1,2-dichloroethene	9.34	61	201214	1.99	ppb	94
30) Hexane	9.59	57	185188	1.95	ppb	88
31) Ethyl acetate	9.60	43	410515	2.00	ppb	97
32) Chloroform	9.69	83	367120	1.97	ppb	98
33) Tetrahydrofuran	10.20	42	140362	2.05	ppb	92
34) 1,2-dichloroethane	10.60	62	271231	1.98	ppb	98
36) 1,1,1-trichloroethane	10.93	97	396083	2.03	ppb	97
37) Cyclohexane	11.84	56	201108	2.20	ppb	98
38) Carbon tetrachloride	11.69	117	412399	2.05	ppb	99
39) Benzene	11.50	78	414323	2.08	ppb	92
40) Methyl methacrylate	13.00	41	194914	2.00	ppb	96
41) 1,4-dioxane	12.79	58	86895	2.08	ppb	# 56
42) 2,2,4-trimethylpentane	12.81	57	633637	2.08	ppb	92
43) Heptane	13.12	43	221392	2.06	ppb	96
44) Trichloroethene	12.77	130	198590	2.05	ppb	93
45) 1,2-dichloropropane	12.49	63	151921	2.04	ppb	98

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\AO033103.D

Vial: 2

Acq On : 31 Mar 2017 3:35 pm

Operator: RJP

Sample : A1UG\_2.0

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 31 21:02:10 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Fri Mar 31 21:00:46 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.72	83	401238	2.07	ppb	99
47) cis-1,3-dichloropropene	13.77	75	254736	2.05	ppb	97
48) trans-1,3-dichloropropene	14.36	75	247830	2.13	ppb	96
49) 1,1,2-trichloroethane	14.56	97	180605	2.01	ppb	99
51) Toluene	14.88	92	291345	1.94	ppb	# 85
52) Methyl Isobutyl Ketone	13.81	43	322762	2.00	ppb	95
53) Dibromochloromethane	15.35	129	364835	2.03	ppb	92
54) Methyl Butyl Ketone	15.16	43	303971	2.09	ppb	96
55) 1,2-dibromoethane	15.62	107	300368	1.96	ppb	97
56) Tetrachloroethylene	16.13	164	199230	1.98	ppb	98
57) Chlorobenzene	16.86	112	411684	1.99	ppb	95
58) Ethylbenzene	17.26	91	694504	2.04	ppb	98
59) m&p-xylene	17.45	91	1170790	4.00	ppb	99
60) Nonane	18.17	43	331734	2.03	ppb	92
61) Styrene	17.84	104	371528	2.06	ppb	100
62) Bromoform	17.54	173	311868	2.08	ppb	100
63) o-xylene	17.95	91	562995	2.01	ppb	93
64) Cumene	18.59	105	724745	1.99	ppb	99
66) 1,1,2,2-tetrachloroethane	17.95	83	373160	1.98	ppb	96
67) Propylbenzene	19.16	120	186824	1.98	ppb	# 1
68) 2-Chlorotoluene	19.13	126	170125	1.95	ppb	# 1
69) 4-ethyltoluene	19.32	105	687148	1.98	ppb	100
70) 1,3,5-trimethylbenzene	19.41	105	616942	1.98	ppb	97
71) 1,2,4-trimethylbenzene	19.88	105	578466	2.03	ppb	95
72) 1,3-dichlorobenzene	20.08	146	374300m	2.04	ppb	
73) benzyl chloride	20.05	91	346380	2.26	ppb	94
74) 1,4-dichlorobenzene	20.15	146	358330m	2.16	ppb	
75) 1,2,3-trimethylbenzene	20.40	105	575572	2.03	ppb	97
76) 1,2-dichlorobenzene	20.57	146	358230	2.11	ppb	98
77) 1,2,4-trichlorobenzene	22.75	180	131000	2.46	ppb	95
78) Naphthalene	22.89	128	336587	2.51	ppb	95
79) Hexachloro-1,3-butadiene	23.31	225	264113	2.05	ppb	99



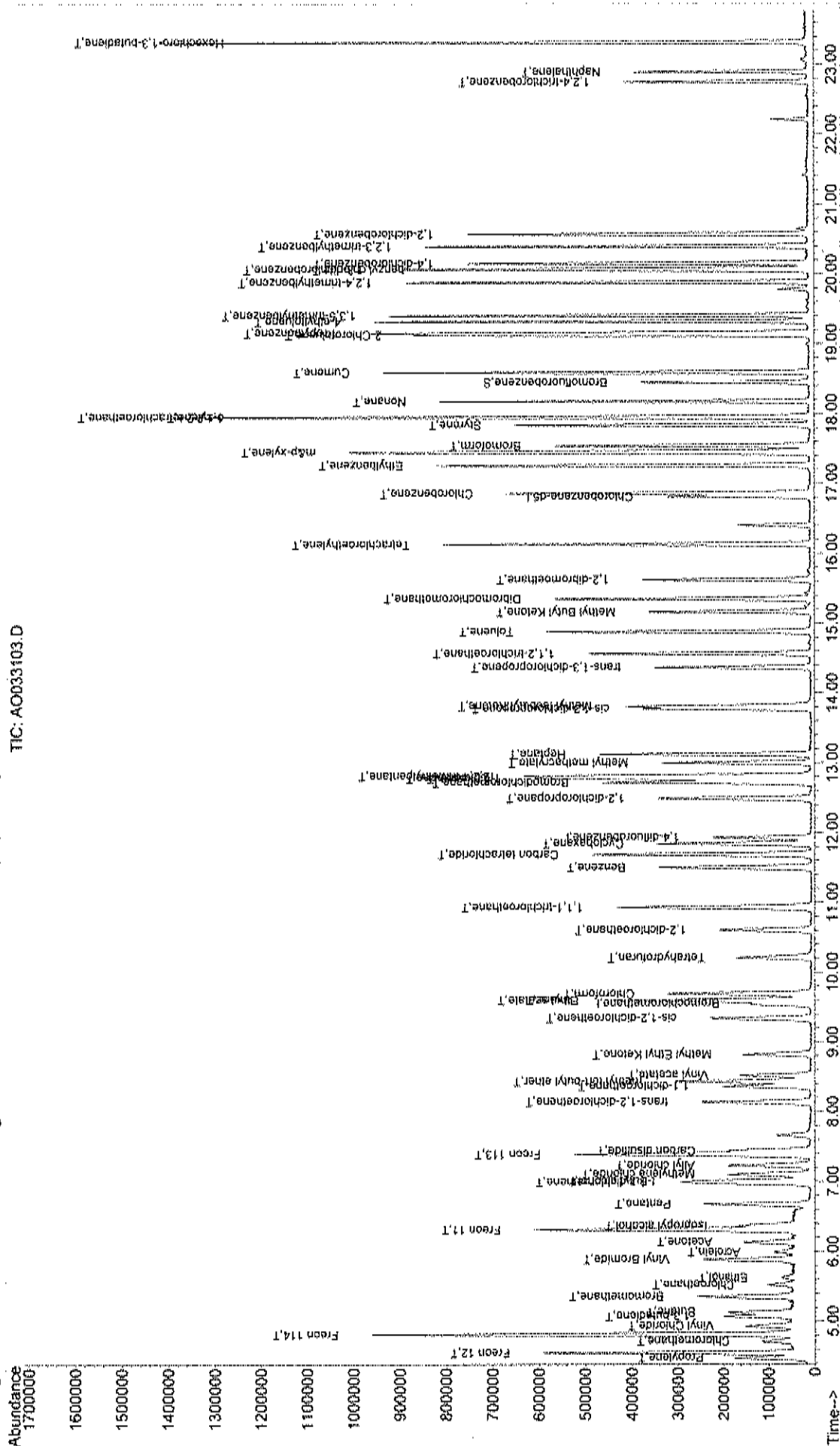
Data File : C:\HPCHEM\1\DATA\AO033103.D  
 Acq On : 31 Mar 2017 3:35 pm  
 Sample : ALUG 2.0  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 3 9:53 2017

Vial: 2  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:13:15 2017  
 Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

TIC: AO033103.D



Data File : C:\HPCHEM\1\DATA\AO033104.D

Vial: 3

Acq On : 31 Mar 2017 4:16 pm

Operator: RJP

Sample : A1UG\_1.5

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 31 21:01:43 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Fri Mar 31 21:00:46 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	49197	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.93	114	224035	1.00	ppb	0.00
50) Chlorobenzene-d5	16.81	117	194095	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.45	95	139602	1.04	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	104.00%

## Target Compounds

						Qvalue
2) Propylene	4.47	41	102672	1.45	ppb	89
3) Freon 12	4.55	85	561521	1.43	ppb	99
4) Chloromethane	4.71	50	81991m	1.49	ppb	
5) Freon 114	4.81	85	440539	1.46	ppb	95
6) Vinyl Chloride	4.93	62	113742	1.44	ppb	91
7) Butane	5.13	43	119938	1.50	ppb	95
8) 1,3-butadiene	5.07	39	70803m	1.25	ppb	
9) Bromomethane	5.35	94	161459	1.43	ppb	96
10) Chloroethane	5.52	64	51047	1.51	ppb	# 77
11) Ethanol	5.63	45	39277	1.46	ppb	86
12) Acrolein	5.99	56	37252	1.27	ppb	87
13) Vinyl Bromide	5.88	106	161219	1.40	ppb	92
14) Freon 11	6.30	101	563957	1.43	ppb	99
15) Acetone	6.13	58	44234	1.50	ppb	# 84
16) Pentane	6.68	42	87203	1.40	ppb	# 35
17) Isopropyl alcohol	6.38	45	150312	1.42	ppb	# 100
18) 1,1-dichloroethene	6.97	96	90564	1.51	ppb	# 88
19) Freon 113	7.38	101	223955	1.53	ppb	91
20) t-Butyl alcohol	7.02	59	226794	1.47	ppb	95
21) Methylene chloride	7.10	84	90177	1.49	ppb	84
22) Allyl chloride	7.23	41	114971	1.48	ppb	89
23) Carbon disulfide	7.44	76	279484	1.45	ppb	84
24) trans-1,2-dichloroethene	8.14	61	156541	1.52	ppb	92
25) methyl tert-butyl ether	8.44	73	296555	1.48	ppb	91
26) 1,1-dichloroethane	8.36	63	190247	1.46	ppb	94
27) Vinyl acetate	8.52	43	245094m	1.52	ppb	
28) Methyl Ethyl Ketone	8.82	72	48323	1.43	ppb	99
29) cis-1,2-dichloroethene	9.34	61	148832	1.50	ppb	91
30) Hexane	9.60	57	134826	1.45	ppb	88
31) Ethyl acetate	9.60	43	306467	1.52	ppb	96
32) Chloroform	9.69	83	263762	1.45	ppb	97
33) Tetrahydrofuran	10.21	42	101943	1.52	ppb	90
34) 1,2-dichloroethane	10.60	62	197805	1.47	ppb	96
36) 1,1,1-trichloroethane	10.93	97	288962	1.52	ppb	97
37) Cyclohexane	11.84	56	146732	1.65	ppb	98
38) Carbon tetrachloride	11.69	117	301856	1.54	ppb	100
39) Benzene	11.50	78	303620	1.57	ppb	91
40) Methyl methacrylate	13.00	41	145032	1.53	ppb	94
41) 1,4-dioxane	12.81	58	63132	1.55	ppb	# 51
42) 2,2,4-trimethylpentane	12.82	57	464131	1.57	ppb	92
43) Heptane	13.12	43	160113	1.53	ppb	96
44) Trichloroethene	12.78	130	144026	1.53	ppb	90
45) 1,2-dichloropropane	12.49	63	110806	1.53	ppb	98

(#)=qualifier out of range (m)=manual integration

AO033104.D A331\_1UG.M

Thu May 04 11:15:33 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO033104.D  
 Acq On : 31 Mar 2017 4:16 pm  
 Sample : A1UG\_1.5  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Mar 31 21:01:43 2017

Vial: 3  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

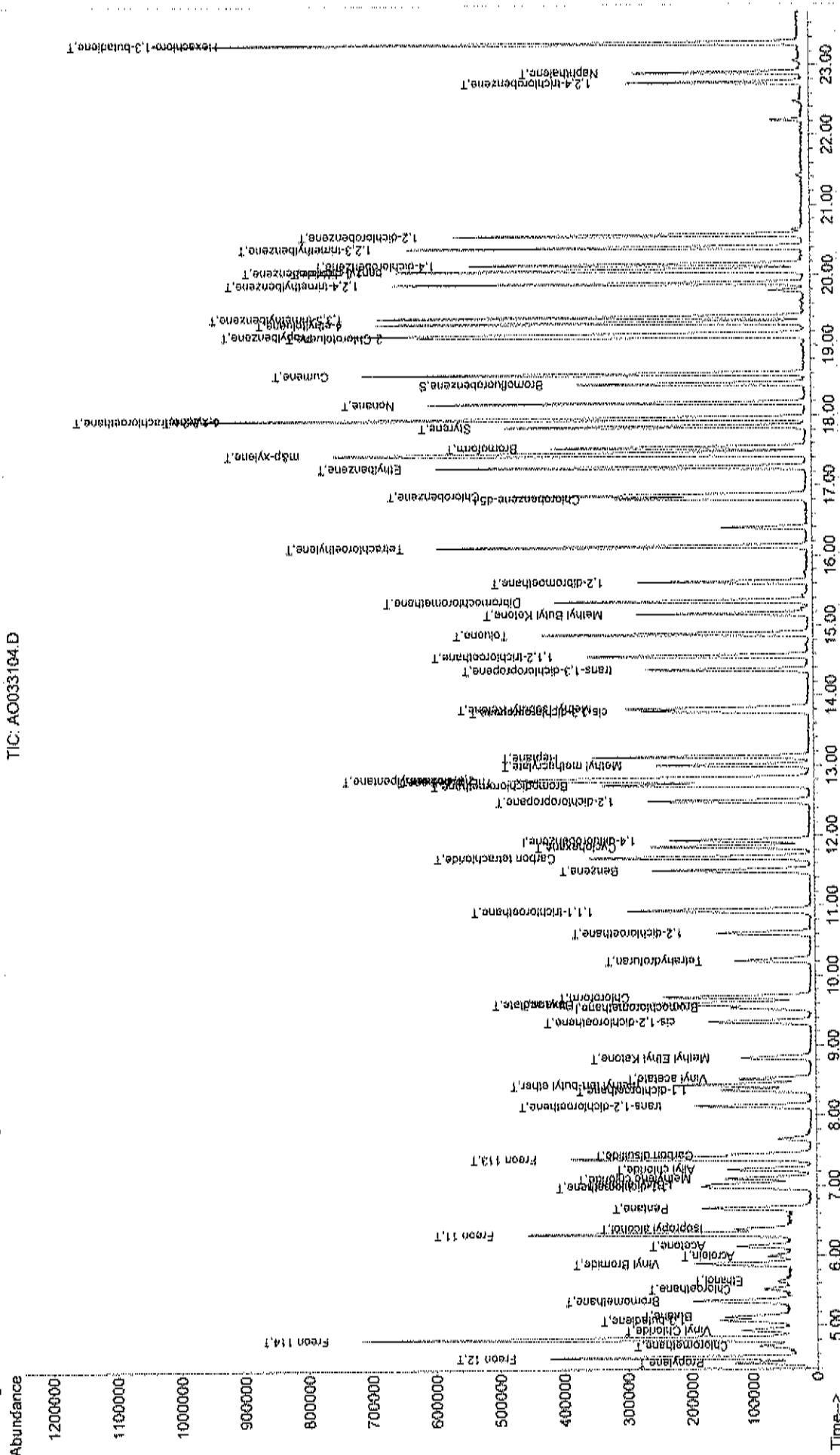
Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Fri Mar 31 21:00:46 2017  
 Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D  
 DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.72	83	289231	1.53	ppb	96
47) cis-1,3-dichloropropene	13.77	75	188783	1.56	ppb	98
48) trans-1,3-dichloropropene	14.37	75	180108	1.59	ppb	96
49) 1,1,2-trichloroethane	14.57	97	137201	1.57	ppb	97
51) Toluene	14.88	92	213554	1.46	ppb	# 82
52) Methyl Isobutyl Ketone	13.81	43	238170	1.52	ppb	94
53) Dibromochloromethane	15.35	129	265843	1.52	ppb	92
54) Methyl Butyl Ketone	15.17	43	222239	1.57	ppb	97
55) 1,2-dibromoethane	15.62	107	217690	1.46	ppb	95
56) Tetrachloroethylene	16.13	164	145863	1.49	ppb	97
57) Chlorobenzene	16.86	112	296747	1.47	ppb	94
58) Ethylbenzene	17.26	91	514456	1.56	ppb	99
59) m&p-xylene	17.45	91	863655	3.04	ppb	98
60) Nonane	18.17	43	238695	1.51	ppb	94
61) Styrene	17.84	104	271617	1.55	ppb	99
62) Bromoform	17.54	173	223130	1.53	ppb	99
63) o-xylene	17.96	91	413467	1.52	ppb	94
64) Cumene	18.59	105	525741	1.49	ppb	100
66) 1,1,2,2-tetrachloroethane	17.95	83	277397	1.52	ppb	97
67) Propylbenzene	19.16	120	134109	1.46	ppb	# 1
68) 2-Chlorotoluene	19.12	126	126693	1.50	ppb	# 1
69) 4-ethyltoluene	19.32	105	509228	1.51	ppb	99
70) 1,3,5-trimethylbenzene	19.40	105	443048	1.47	ppb	98
71) 1,2,4-trimethylbenzene	19.88	105	426277	1.54	ppb	97
72) 1,3-dichlorobenzene	20.08	146	276528	1.55	ppb	98
73) benzyl chloride	20.06	91	246509	1.66	ppb	94
74) 1,4-dichlorobenzene	20.16	146	249352	1.55	ppb	96
75) 1,2,3-trimethylbenzene	20.40	105	426217	1.55	ppb	98
76) 1,2-dichlorobenzene	20.57	146	262645	1.60	ppb	98
77) 1,2,4-trichlorobenzene	22.75	180	90071	1.74	ppb	95
78) Naphthalene	22.89	128	235547	1.81	ppb	94
79) Hexachloro-1,3-butadiene	23.31	225	191216	1.53	ppb	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed  
 AO033104.D A331\_1UG.M Thu May 04 11:15:34 2017 MSD1

TIC: A0033104.D



Data File : C:\HPCHEM\1\DATA\AO033105.D  
 Acq On : 31 Mar 2017 4:57 pm  
 Sample : A1UG\_1.25  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Mar 31 21:01:19 2017

Vial: 4  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Fri Mar 31 21:00:46 2017  
 Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D  
 DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	46123	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.94	114	218725	1.00	ppb	0.00
50) Chlorobenzene-d5	16.81	117	193903	1.00	ppb	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
65) Bromofluorobenzene	18.45	95	130135	0.97	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	97.00%

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Propylene	4.48	41	78567	1.19	ppb	84
3) Freon 12	4.55	85	467760	1.27	ppb	98
4) Chloromethane	4.72	50	65723m A	1.28	ppb	
5) Freon 114	4.82	85	357396	1.26	ppb	94
6) Vinyl Chloride	4.94	62	96877	1.31	ppb	90
7) Butane	5.14	43	100777	1.35	ppb	95
8) 1,3-butadiene	5.07	39	68213m P	1.29	ppb	
9) Bromomethane	5.36	94	131929	1.25	ppb	93
10) Chloroethane	5.53	64	41303	1.30	ppb	# 84
11) Ethanol	5.63	45	31486	1.25	ppb	88
12) Acrolein	5.99	56	32448	1.18	ppb	89
13) Vinyl Bromide	5.89	106	136500	1.26	ppb	95
14) Freon 11	6.31	101	468764	1.27	ppb	99
15) Acetone	6.14	58	36412	1.32	ppb	85
16) Pentane	6.68	42	71803	1.23	ppb	# 31
17) Isopropyl alcohol	6.37	45	118797	1.20	ppb	# 100
18) 1,1-dichloroethene	6.98	96	74889	1.33	ppb	# 89
19) Freon 113	7.39	101	179185	1.30	ppb	91
20) t-Butyl alcohol	7.02	59	183128	1.26	ppb	96
21) Methylene chloride	7.10	84	72610	1.28	ppb	# 83
22) Allyl chloride	7.23	41	91853	1.26	ppb	95
23) Carbon disulfide	7.45	76	232308	1.29	ppb	84
24) trans-1,2-dichloroethene	8.14	61	126965	1.32	ppb	93
25) methyl tert-butyl ether	8.44	73	240569	1.28	ppb	91
26) 1,1-dichloroethane	8.37	63	154227	1.26	ppb	96
27) Vinyl acetate	8.52	43	201074m A	1.33	ppb	
28) Methyl Ethyl Ketone	8.83	72	39068	1.23	ppb	97
29) cis-1,2-dichloroethene	9.35	61	120450	1.30	ppb	94
30) Hexane	9.60	57	110322	1.27	ppb	90
31) Ethyl acetate	9.60	43	246897	1.31	ppb	97
32) Chloroform	9.69	83	222198	1.30	ppb	99
33) Tetrahydrofuran	10.21	42	81151	1.29	ppb	89
34) 1,2-dichloroethane	10.60	62	164628	1.31	ppb	97
36) 1,1,1-trichloroethane	10.93	97	235511	1.27	ppb	100
37) Cyclohexane	11.85	56	121625	1.40	ppb	98
38) Carbon tetrachloride	11.69	117	248909	1.30	ppb	99
39) Benzene	11.50	78	246604	1.30	ppb	92
40) Methyl methacrylate	13.01	41	117282	1.27	ppb	97
41) 1,4-dioxane	12.80	58	50550	1.27	ppb	# 50
42) 2,2,4-trimethylpentane	12.82	57	379009	1.31	ppb	91
43) Heptane	13.13	43	129459	1.27	ppb	96
44) Trichloroethene	12.78	130	119946	1.30	ppb	93
45) 1,2-dichloropropane	12.49	63	88102	1.24	ppb	95

(#) = qualifier out of range (m) = manual integration  
 AO033105.D A331\_1UG.M Thu May 04 11:15:37 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO033105.D

Vial: 4

Acq On : 31 Mar 2017 4:57 pm

Operator: RJP

Sample : A1UG\_1.25

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 31 21:01:19 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Fri Mar 31 21:00:46 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.72	83	233632	1.27	ppb	98
47) cis-1,3-dichloropropene	13.78	75	150325	1.27	ppb	98
48) trans-1,3-dichloropropene	14.37	75	145587	1.32	ppb	98
49) 1,1,2-trichloroethane	14.57	97	111276	1.30	ppb	98
51) Toluene	14.88	92	172202	1.18	ppb	# 83
52) Methyl Isobutyl Ketone	13.82	43	190568	1.22	ppb	94
53) Dibromochloromethane	15.35	129	219080	1.26	ppb	95
54) Methyl Butyl Ketone	15.17	43	177195	1.25	ppb	96
55) 1,2-dibromoethane	15.62	107	180999	1.22	ppb	98
56) Tetrachloroethylene	16.13	164	119394	1.22	ppb	97
57) Chlorobenzene	16.86	112	246201	1.22	ppb	95
58) Ethylbenzene	17.26	91	412665	1.25	ppb	100
59) m&p-xylene	17.45	91	701907	2.47	ppb	98
60) Nonane	18.18	43	195655	1.24	ppb	94
61) Styrene	17.84	104	222122	1.27	ppb	88
62) Bromoform	17.54	173	180532	1.24	ppb	98
63) o-xylene	17.95	91	337985	1.24	ppb	93
64) Cumene	18.59	105	435179	1.23	ppb	100
66) 1,1,2,2-tetrachloroethane	17.95	83	225167	1.23	ppb	97
67) Propylbenzene	19.16	120	112167	1.22	ppb	# 1
68) 2-Chlorotoluene	19.13	126	102454	1.21	ppb	# 1
69) 4-ethyltoluene	19.32	105	417833	1.24	ppb	99
70) 1,3,5-trimethylbenzene	19.41	105	365290	1.21	ppb	100
71) 1,2,4-trimethylbenzene	19.89	105	343051	1.24	ppb	94
72) 1,3-dichlorobenzene	20.08	146	222984m	1.25	ppb	
73) benzyl chloride	20.06	91	194194	1.31	ppb	94
74) 1,4-dichlorobenzene	20.16	146	206810m	1.29	ppb	
75) 1,2,3-trimethylbenzene	20.40	105	349260	1.27	ppb	97
76) 1,2-dichlorobenzene	20.57	146	206999	1.26	ppb	99
77) 1,2,4-trichlorobenzene	22.75	180	66634	1.29	ppb	93
78) Naphthalene	22.89	128	171519	1.32	ppb	93
79) Hexachloro-1,3-butadiene	23.32	225	149165	1.19	ppb	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed  
 AO033105.D A331\_1UG.M Thu May 04 11:15:37 2017 MSD1



[illegible]

Data File : C:\HPCHEM\1\DATA\AO033106.D

Vial: 5

Acq On : 31 Mar 2017 5:36 pm

Operator: RJP

Sample : A1UG\_1.0

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 31 21:01:04 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Fri Mar 31 21:00:46 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	46112	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.93	114	220001	1.00	ppb	0.00
50) Chlorobenzene-d5	16.82	117	186025	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.45	95	128961	1.00	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	= 100.00%	

## Target Compounds

						Qvalue
2) Propylene	4.49	41	66145	1.00	ppb	96
3) Freon 12	4.56	85	368126	1.00	ppb	98
4) Chloromethane	4.72	50	51707m	1.01	ppb	
5) Freon 114	4.82	85	283775	1.00	ppb	95
6) Vinyl Chloride	4.95	62	73850	1.00	ppb	89
7) Butane	5.14	43	74701	1.00	ppb	92
8) 1,3-butadiene	5.09	39	49285m	0.93	ppb	
9) Bromomethane	5.36	94	105576	1.00	ppb	92
10) Chloroethane	5.54	64	31781	1.00	ppb	# 75
11) Ethanol	5.63	45	25232	1.00	ppb	83
12) Acrolein	6.00	56	27310m	0.99	ppb	
13) Vinyl Bromide	5.89	106	108238	1.00	ppb	93
14) Freon 11	6.31	101	370422	1.00	ppb	99
15) Acetone	6.14	58	27552	1.00	ppb	# 76
16) Pentane	6.68	42	58367	1.00	ppb	# 40
17) Isopropyl alcohol	6.38	45	99005	1.00	ppb	# 100
18) 1,1-dichloroethene	6.99	96	56363	1.00	ppb	# 87
19) Freon 113	7.39	101	137328	1.00	ppb	91
20) t-Butyl alcohol	7.02	59	145442	1.00	ppb	95
21) Methylene chloride	7.10	84	56713	1.00	ppb	# 77
22) Allyl chloride	7.24	41	72786	1.00	ppb	93
23) Carbon disulfide	7.45	76	180433	1.00	ppb	83
24) trans-1,2-dichloroethene	8.14	61	96246	1.00	ppb	97
25) methyl tert-butyl ether	8.44	73	187288	1.00	ppb	90
26) 1,1-dichloroethane	8.38	63	122504	1.00	ppb	95
27) Vinyl acetate	8.53	43	148055m	0.98	ppb	
28) Methyl Ethyl Ketone	8.83	72	31753	1.00	ppb	96
29) cis-1,2-dichloroethene	9.35	61	92886	1.00	ppb	92
30) Hexane	9.60	57	87106	1.00	ppb	89
31) Ethyl acetate	9.61	43	188952	1.00	ppb	96
32) Chloroform	9.70	83	171011	1.00	ppb	100
33) Tetrahydrofuran	10.22	42	62907	1.00	ppb	88
34) 1,2-dichloroethane	10.61	62	125908	1.00	ppb	95
36) 1,1,1-trichloroethane	10.93	97	186177	1.00	ppb	98
37) Cyclohexane	11.85	56	87220	1.00	ppb	93
38) Carbon tetrachloride	11.69	117	192761	1.00	ppb	100
39) Benzene	11.51	78	190289	1.00	ppb	91
40) Methyl methacrylate	13.01	41	93083	1.00	ppb	97
41) 1,4-dioxane	12.81	58	40035	1.00	ppb	# 56
42) 2,2,4-trimethylpentane	12.81	57	291037	1.00	ppb	90
43) Heptane	13.13	43	102640	1.00	ppb	99
44) Trichloroethene	12.78	130	92651	1.00	ppb	93
45) 1,2-dichloropropane	12.49	63	71331	1.00	ppb	99

(#)=qualifier out of range (m)=manual integration

AO033106.D A331\_1UG.M

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MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO033106.D

Vial: 5

Acq On : 31 Mar 2017 5:36 pm

Operator: RJP

Sample : A1UG\_1.0

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 31 21:01:04 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Fri Mar 31 21:00:46 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.72	83	185244	1.00	ppb	98
47) cis-1,3-dichloropropene	13.78	75	118799	1.00	ppb	96
48) trans-1,3-dichloropropene	14.37	75	111119	1.00	ppb	95
49) 1,1,2-trichloroethane	14.57	97	85955	1.00	ppb	99
51) Toluene	14.88	92	139995	1.00	ppb	88
52) Methyl Isobutyl Ketone	13.82	43	149970	1.00	ppb	94
53) Dibromochloromethane	15.35	129	167207	1.00	ppb	89
54) Methyl Butyl Ketone	15.17	43	135647	1.00	ppb	96
55) 1,2-dibromoethane	15.62	107	142683	1.00	ppb	97
56) Tetrachloroethylene	16.13	164	93558	1.00	ppb	100
57) Chlorobenzene	16.86	112	192859	1.00	ppb	96
58) Ethylbenzene	17.26	91	316266	1.00	ppb	100
59) m&p-xylene	17.46	91	546554	2.01	ppb	98
60) Nonane	18.18	43	151897	1.00	ppb	93
61) Styrene	17.84	104	168155	1.00	ppb	99
62) Bromoform	17.54	173	139588	1.00	ppb	98
63) o-xylene	17.96	91	260479	1.00	ppb	93
64) Cumene	18.59	105	339221	1.00	ppb	100
66) 1,1,2,2-tetrachloroethane	17.95	83	175288	1.00	ppb	96
67) Propylbenzene	19.16	120	87986	1.00	ppb	# 1
68) 2-Chlorotoluene	19.12	126	81024	1.00	ppb	# 1
69) 4-ethyltoluene	19.32	105	322806	1.00	ppb	99
70) 1,3,5-trimethylbenzene	19.41	105	289695	1.00	ppb	98
71) 1,2,4-trimethylbenzene	19.89	105	264743	1.00	ppb	94
72) 1,3-dichlorobenzene	20.08	146	170760	1.00	ppb	99
73) benzyl chloride	20.05	91	142227	1.00	ppb	92
74) 1,4-dichlorobenzene	20.16	146	154100	1.00	ppb	98
75) 1,2,3-trimethylbenzene	20.40	105	263431	1.00	ppb	94
76) 1,2-dichlorobenzene	20.58	146	157732	1.00	ppb	97
77) 1,2,4-trichlorobenzene	22.75	180	49634	1.00	ppb	96
78) Naphthalene	22.90	128	124780	1.00	ppb	94
79) Hexachloro-1,3-butadiene	23.31	225	119753	1.00	ppb	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed  
 AO033106.D A331\_1UG.M Thu May 04 11:15:41 2017 MSD1

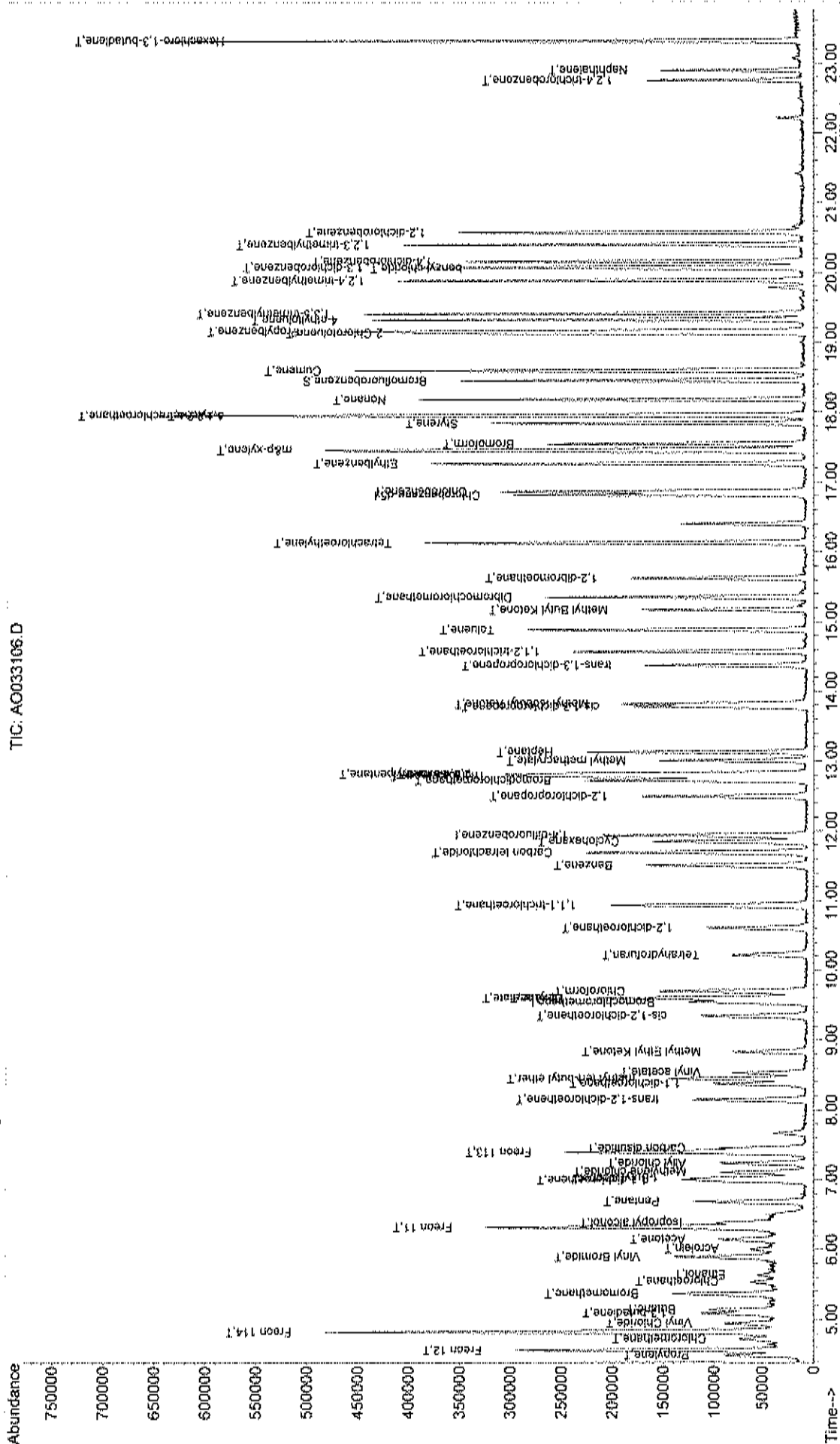
Data File : C:\HPCHEM\1\DATA\AO033106.D  
Acq On : 31 Mar 2017 5:36 pm  
Sample : ALUG\_1.0  
Misc : A331\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Apr 3 10:01 2017

Vial: 5  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Thu May 04 11:13:15 2017  
Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

TIC: AO033106.D



Data File : C:\HPCHEM\1\DATA\AO033107.D  
 Acq On : 31 Mar 2017 6:15 pm  
 Sample : A1UG\_0.75  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Mar 31 21:02:35 2017

Vial: 6  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Fri Mar 31 21:00:46 2017  
 Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D  
 DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	44024	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.93	114	210749	1.00	ppb	0.00
50) Chlorobenzene-d5	16.82	117	183483	1.00	ppb	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
65) Bromofluorobenzene	18.45	95	125665	0.99	ppb	0.00
Spiked Amount 1.000	Range 70 - 130		Recovery	=	99.00%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Propylene	4.49	41	45360	0.72	ppb	80
3) Freon 12	4.56	85	277878	0.79	ppb	97
4) Chloromethane	4.73	50	37635	0.77	ppb	89
5) Freon 114	4.82	85	215987	0.80	ppb	96
6) Vinyl Chloride	4.94	62	56874	0.81	ppb	90
7) Butane	5.14	43	58073	0.81	ppb	92
8) 1,3-butadiene	5.08	39	39119m	0.77	ppb	
9) Bromomethane	5.36	94	77287	0.77	ppb	93
10) Chloroethane	5.54	64	23478	0.77	ppb	# 63
11) Ethanol	5.64	45	18822m	0.78	ppb	
12) Acrolein	6.01	56	19805	0.75	ppb	92
13) Vinyl Bromide	5.89	106	79505	0.77	ppb	94
14) Freon 11	6.31	101	275669	0.78	ppb	98
15) Acetone	6.15	58	20215	0.77	ppb	# 67
16) Pentane	6.68	42	42253	0.76	ppb	# 39
17) Isopropyl alcohol	6.39	45	72401	0.77	ppb	# 100
18) 1,1-dichloroethene	6.98	96	42805	0.80	ppb	# 88
19) Freon 113	7.39	101	102624	0.78	ppb	92
20) t-Butyl alcohol	7.03	59	101096	0.73	ppb	# 94
21) Methylene chloride	7.11	84	40588	0.75	ppb	# 76
22) Allyl chloride	7.24	41	54748	0.79	ppb	89
23) Carbon disulfide	7.45	76	136144	0.79	ppb	83
24) trans-1,2-dichloroethene	8.14	61	71552	0.78	ppb	93
25) methyl tert-butyl ether	8.45	73	140829	0.79	ppb	90
26) 1,1-dichloroethane	8.37	63	89176	0.76	ppb	93
27) Vinyl acetate	8.53	43	113620m	0.79	ppb	
28) Methyl Ethyl Ketone	8.83	72	21826	0.72	ppb	94
29) cis-1,2-dichloroethene	9.34	61	69108	0.78	ppb	91
30) Hexane	9.60	57	65405	0.79	ppb	90
31) Ethyl acetate	9.61	43	140819	0.78	ppb	98
32) Chloroform	9.70	83	126305	0.77	ppb	99
33) Tetrahydrofuran	10.22	42	47717	0.79	ppb	88
34) 1,2-dichloroethane	10.62	62	95385	0.79	ppb	99
36) 1,1,1-trichloroethane	10.93	97	134349	0.75	ppb	98
37) Cyclohexane	11.85	56	70550	0.84	ppb	96
38) Carbon tetrachloride	11.69	117	139907	0.76	ppb	98
39) Benzene	11.50	78	142039	0.78	ppb	91
40) Methyl methacrylate	13.00	41	67157m	0.75	ppb	
41) 1,4-dioxane	12.81	58	26970	0.70	ppb	# 56
42) 2,2,4-trimethylpentane	12.82	57	218109	0.78	ppb	91
43) Heptane	13.13	43	74753	0.76	ppb	99
44) Trichloroethene	12.78	130	69266	0.78	ppb	94
45) 1,2-dichloropropane	12.49	63	53577	0.78	ppb	99

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\AO033107.D

Vial: 6

Acq On : 31 Mar 2017 6:15 pm

Operator: RJP

Sample : A1UG\_0.75

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 31 21:02:35 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Fri Mar 31 21:00:46 2017

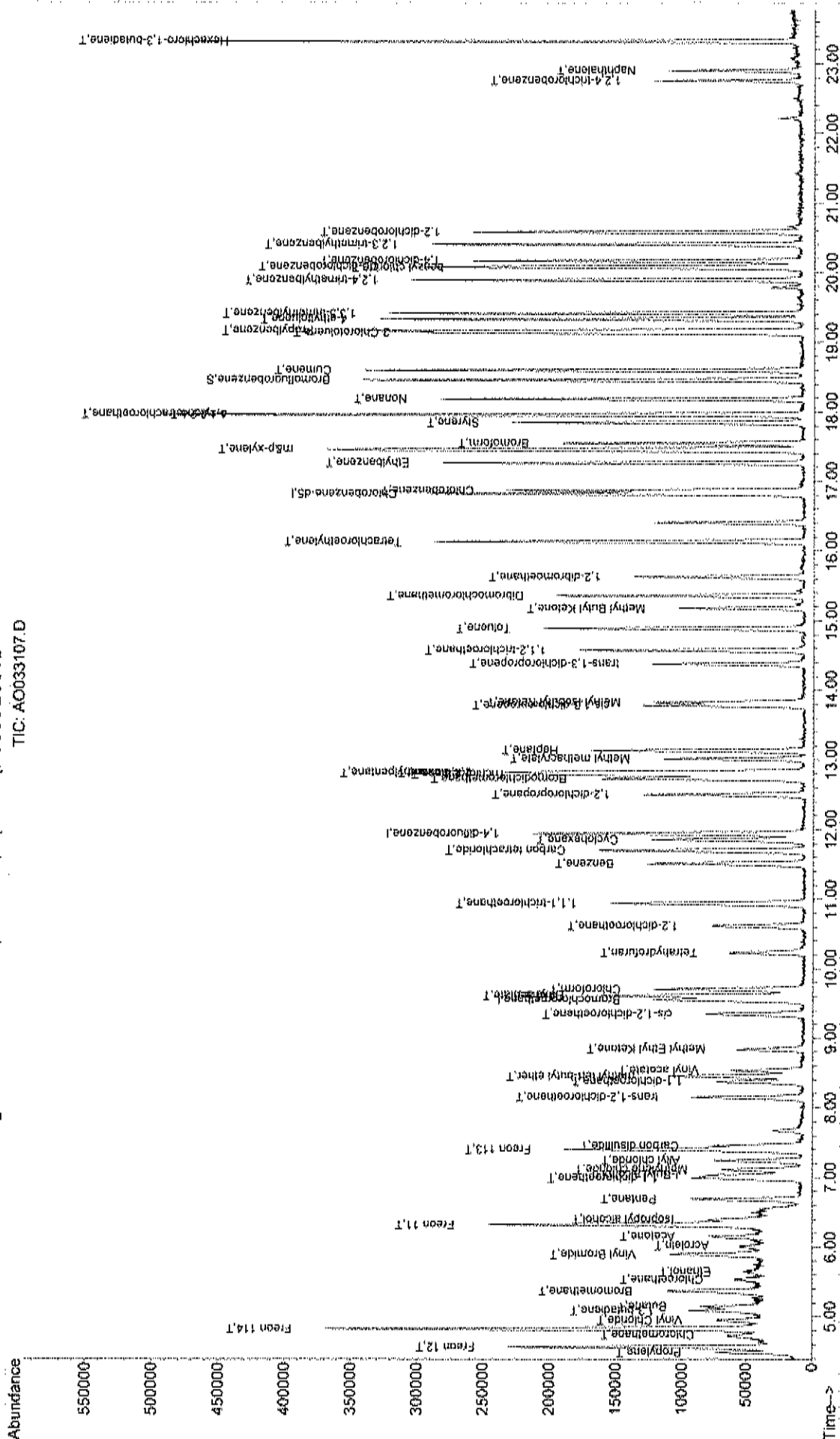
Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.72	83	138457	0.78	ppb	100
47) cis-1,3-dichloropropene	13.78	75	84571	0.74	ppb	97
48) trans-1,3-dichloropropene	14.38	75	83731	0.79	ppb	97
49) 1,1,2-trichloroethane	14.57	97	63090	0.77	ppb	97
51) Toluene	14.89	92	100613	0.73	ppb #	83
52) Methyl Isobutyl Ketone	13.82	43	94497	0.64	ppb	95
53) Dibromochloromethane	15.36	129	123463	0.75	ppb	95
54) Methyl Butyl Ketone	15.18	43	81165	0.61	ppb	97
55) 1,2-dibromoethane	15.62	107	102493	0.73	ppb	97
56) Tetrachloroethylene	16.13	164	68551	0.74	ppb	99
57) Chlorobenzene	16.86	112	142252	0.75	ppb	94
58) Ethylbenzene	17.26	91	236565	0.76	ppb	100
59) m&p-xylene	17.45	91	410271	1.53	ppb	99
60) Nonane	18.18	43	109229	0.73	ppb	94
61) Styrene	17.84	104	123584	0.75	ppb	98
62) Bromoform	17.55	173	103675	0.75	ppb	98
63) o-xylene	17.96	91	194699	0.76	ppb	93
64) Cumene	18.59	105	250016	0.75	ppb	99
66) 1,1,2,2-tetrachloroethane	17.95	83	132210	0.76	ppb	97
67) Propylbenzene	19.17	120	64596	0.74	ppb #	1
68) 2-Chlorotoluene	19.13	126	58331	0.73	ppb #	1
69) 4-ethyltoluene	19.33	105	232250	0.73	ppb	100
70) 1,3,5-trimethylbenzene	19.41	105	207423	0.73	ppb	98
71) 1,2,4-trimethylbenzene	19.89	105	195195	0.75	ppb	95
72) 1,3-dichlorobenzene	20.08	146	126865	0.75	ppb #	55
73) benzyl chloride	20.06	91	98302	0.70	ppb	93
74) 1,4-dichlorobenzene	20.16	146	115236	0.76	ppb #	52
75) 1,2,3-trimethylbenzene	20.40	105	193918	0.75	ppb	93
76) 1,2-dichlorobenzene	20.57	146	118355	0.76	ppb	99
77) 1,2,4-trichlorobenzene	22.75	180	33339	0.68	ppb	92
78) Naphthalene	22.90	128	82942	0.67	ppb	96
79) Hexachloro-1,3-butadiene	23.32	225	88256	0.75	ppb	98

-----  
 (#) = qualifier out of range (m) = manual integration (+) = signals summed  
 AO033107.D A331\_1UG.M Thu May 04 11:15:45 2017 MSD1

TIC: A0033107.D





Data File : C:\HPCHEM\1\DATA\AO033108.D

Vial: 7

Acq On : 31 Mar 2017 6:53 pm

Operator: RJP

Sample : A1UG\_0.50

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 31 21:03:17 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Fri Mar 31 21:00:46 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.55	128	43920	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.93	114	204815	1.00	ppb	0.00
50) Chlorobenzene-d5	16.82	117	181136	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.45	95	121444	0.97	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	97.00%

## Target Compounds

						Qvalue
2) Propylene	4.48	41	31388	0.50	ppb	79
3) Freon 12	4.56	85	182600	0.52	ppb	98
4) Chloromethane	4.72	50	24920	0.51	ppb	91
5) Freon 114	4.82	85	143482	0.53	ppb	96
6) Vinyl Chloride	4.95	62	37274	0.53	ppb	87
7) Butane	5.14	43	41459	0.58	ppb	92
8) 1,3-butadiene	5.08	39	25730m	0.51	ppb	
9) Bromomethane	5.35	94	51121	0.51	ppb	96
10) Chloroethane	5.53	64	16560	0.55	ppb	# 69
11) Ethanol	5.64	45	11702m	0.49	ppb	
12) Acrolein	6.00	56	14036	0.53	ppb	90
13) Vinyl Bromide	5.89	106	50717	0.49	ppb	91
14) Freon 11	6.31	101	185391	0.53	ppb	99
15) Acetone	6.15	58	14318	0.55	ppb	# 84
16) Pentane	6.68	42	27157	0.49	ppb	# 31
17) Isopropyl alcohol	6.38	45	44936	0.48	ppb	# 100
18) 1,1-dichloroethene	6.98	96	29783	0.55	ppb	93
19) Freon 113	7.40	101	68337	0.52	ppb	91
20) t-Butyl alcohol	7.03	59	65506	0.47	ppb	94
21) Methylene chloride	7.11	84	28901	0.54	ppb	# 82
22) Allyl chloride	7.23	41	33667	0.49	ppb	93
23) Carbon disulfide	7.45	76	89263	0.52	ppb	84
24) trans-1,2-dichloroethene	8.15	61	47197	0.51	ppb	95
25) methyl tert-butyl ether	8.45	73	88167	0.49	ppb	89
26) 1,1-dichloroethane	8.37	63	60989	0.52	ppb	96
27) Vinyl acetate	8.52	43	72472m	0.50	ppb	
28) Methyl Ethyl Ketone	8.84	72	14916	0.49	ppb	96
29) cis-1,2-dichloroethene	9.35	61	45517	0.51	ppb	92
30) Hexane	9.59	57	41945	0.51	ppb	88
31) Ethyl acetate	9.61	43	93257	0.52	ppb	96
32) Chloroform	9.70	83	83520	0.51	ppb	98
33) Tetrahydrofuran	10.23	42	30981	0.52	ppb	91
34) 1,2-dichloroethane	10.61	62	62717	0.52	ppb	93
36) 1,1,1-trichloroethane	10.94	97	88752	0.51	ppb	99
37) Cyclohexane	11.85	56	43436	0.53	ppb	98
38) Carbon tetrachloride	11.69	117	90406	0.50	ppb	95
39) Benzene	11.51	78	92910	0.52	ppb	90
40) Methyl methacrylate	13.01	41	42361	0.49	ppb	98
41) 1,4-dioxane	12.82	58	17766	0.48	ppb	# 46
42) 2,2,4-trimethylpentane	12.82	57	139898	0.52	ppb	89
43) Heptane	13.13	43	49667	0.52	ppb	98
44) Trichloroethene	12.78	130	43044	0.50	ppb	86
45) 1,2-dichloropropane	12.50	63	34841	0.52	ppb	99

(#)=qualifier out of range (m)=manual integration

AO033108.D A331\_1UG.M

Thu May 04 11:15:49 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO033108.D

Vial: 7

Acq On : 31 Mar 2017 6:53 pm

Operator: RJP

Sample : A1UG\_0.50

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 31 21:03:17 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Fri Mar 31 21:00:46 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

DataAcq Meth : 1UG\_RUN

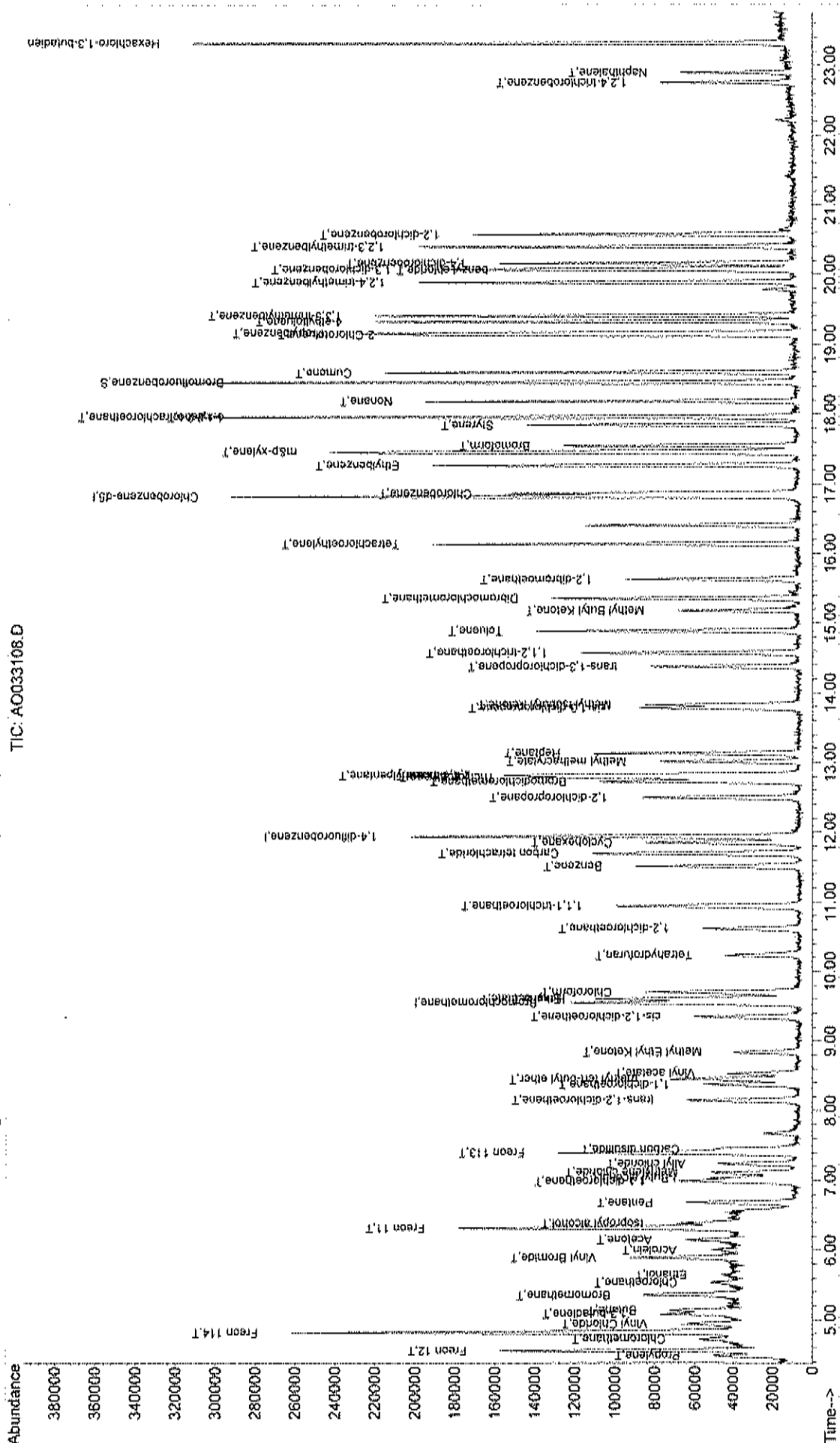
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.72	83	88899	0.52	ppb	99
47) cis-1,3-dichloropropene	13.78	75	55495	0.50	ppb	98
48) trans-1,3-dichloropropene	14.37	75	53706	0.52	ppb	91
49) 1,1,2-trichloroethane	14.57	97	41634	0.52	ppb	99
51) Toluene	14.88	92	65569	0.48	ppb	# 82
52) Methyl Isobutyl Ketone	13.83	43	62120	0.43	ppb	94
53) Dibromochloromethane	15.36	129	81972	0.50	ppb	94
54) Methyl Butyl Ketone	15.17	43	51833	0.39	ppb	95
55) 1,2-dibromoethane	15.62	107	66022	0.48	ppb	97
56) Tetrachloroethylene	16.14	164	46315	0.51	ppb	98
57) Chlorobenzene	16.87	112	91788	0.49	ppb	93
58) Ethylbenzene	17.26	91	154803	0.50	ppb	100
59) m&p-xylene	17.46	91	262800	0.99	ppb	100
60) Nonane	18.18	43	71262	0.48	ppb	96
61) Styrene	17.84	104	79431	0.49	ppb	97
62) Bromoform	17.55	173	66152	0.49	ppb	98
63) o-xylene	17.96	91	124998	0.49	ppb	94
64) Cumene	18.60	105	161847	0.49	ppb	99
66) 1,1,2,2-tetrachloroethane	17.95	83	86064	0.50	ppb	97
67) Propylbenzene	19.16	120	41634	0.49	ppb	# 1
68) 2-Chlorotoluene	19.13	126	39209	0.50	ppb	# 1
69) 4-ethyltoluene	19.32	105	152603	0.49	ppb	99
70) 1,3,5-trimethylbenzene	19.41	105	136526	0.48	ppb	99
71) 1,2,4-trimethylbenzene	19.89	105	126003	0.49	ppb	94
72) 1,3-dichlorobenzene	20.08	146	76215m	0.46	ppb	
73) benzyl chloride	20.05	91	61148	0.44	ppb	93
74) 1,4-dichlorobenzene	20.16	146	72236m	0.48	ppb	
75) 1,2,3-trimethylbenzene	20.40	105	126884	0.49	ppb	97
76) 1,2-dichlorobenzene	20.58	146	75655	0.49	ppb	97
77) 1,2,4-trichlorobenzene	22.76	180	19912	0.41	ppb	93
78) Naphthalene	22.90	128	49444	0.41	ppb	96
79) Hexachloro-1,3-butadiene	23.32	225	55006	0.47	ppb	96

Data File : C:\HPCHEM\1\DATA\AO033108.D  
 Acq On : 31 Mar 2017 6:53 pm  
 Sample : A1UG 0.50  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 3 10:05 2017

Vial: 7  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00  
 Quant Results File: A331\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:13:15 2017  
 Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

TIC: AO033108.D



Data File : C:\HPCHEM\1\DATA\AO033109.D

Vial: 8

Acq On : 31 Mar 2017 7:30 pm

Operator: RJP

Sample : A1UG\_0.30

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 31 21:03:49 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Fri Mar 31 21:00:46 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	44360	1.00	ppb	0.01
35) 1,4-difluorobenzene	11.94	114	201658	1.00	ppb	0.00
50) Chlorobenzene-d5	16.82	117	174946	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	117955	0.97	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	97.00%

## Target Compounds

						Qvalue
2) Propylene	4.49	41	20130	0.32	ppb	# 57
3) Freon 12	4.55	85	111633	0.32	ppb	93
4) Chloromethane	4.72	50	16029	0.32	ppb	67
5) Freon 114	4.81	85	88039	0.32	ppb	96
6) Vinyl Chloride	4.94	62	23880	0.34	ppb	89
7) Butane	5.13	43	24791	0.34	ppb	92
8) 1,3-butadiene	5.09	39	17729m	0.35	ppb	
9) Bromomethane	5.35	94	36947	0.36	ppb	81
10) Chloroethane	5.53	64	10936	0.36	ppb	# 38
11) Ethanol	5.65	45	9308	0.38	ppb	# 67
12) Acrolein	6.01	56	7336m	0.28	ppb	
13) Vinyl Bromide	5.89	106	32409	0.31	ppb	95
14) Freon 11	6.32	101	112125	0.31	ppb	99
15) Acetone	6.14	58	9097	0.34	ppb	# 77
16) Pentane	6.69	42	16757	0.30	ppb	# 39
17) Isopropyl alcohol	6.39	45	31119	0.33	ppb	# 100
18) 1,1-dichloroethene	6.99	96	17110	0.32	ppb	92
19) Freon 113	7.39	101	40494	0.31	ppb	92
20) t-Butyl alcohol	7.03	59	43837	0.31	ppb	# 94
21) Methylene chloride	7.11	84	17210	0.32	ppb	# 76
22) Allyl chloride	7.23	41	21572	0.31	ppb	90
23) Carbon disulfide	7.45	76	55879	0.32	ppb	84
24) trans-1,2-dichloroethene	8.15	61	29814	0.32	ppb	91
25) methyl tert-butyl ether	8.45	73	56728	0.31	ppb	88
26) 1,1-dichloroethane	8.37	63	36458	0.31	ppb	98
27) Vinyl acetate	8.52	43	45130m	0.31	ppb	
28) Methyl Ethyl Ketone	8.84	72	8808	0.29	ppb	# 91
29) cis-1,2-dichloroethene	9.35	61	27589	0.31	ppb	95
30) Hexane	9.59	57	25849	0.31	ppb	89
31) Ethyl acetate	9.62	43	58527	0.32	ppb	97
32) Chloroform	9.70	83	50295	0.31	ppb	96
33) Tetrahydrofuran	10.23	42	19194	0.32	ppb	92
34) 1,2-dichloroethane	10.61	62	36324	0.30	ppb	93
36) 1,1,1-trichloroethane	10.94	97	56401	0.33	ppb	94
37) Cyclohexane	11.85	56	31511	0.39	ppb	87
38) Carbon tetrachloride	11.69	117	55427	0.31	ppb	96
39) Benzene	11.51	78	59047	0.34	ppb	93
40) Methyl methacrylate	13.02	41	26739	0.31	ppb	97
41) 1,4-dioxane	12.81	58	13328	0.36	ppb	# 47
42) 2,2,4-trimethylpentane	12.82	57	82460	0.31	ppb	87
43) Heptane	13.13	43	28810	0.31	ppb	97
44) Trichloroethene	12.78	130	26660	0.31	ppb	89
45) 1,2-dichloropropane	12.51	63	20460	0.31	ppb	96

(# ) = qualifier out of range (m) = manual integration

AO033109.D A331\_1UG.M

Thu May 04 11:15:53 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO033109.D

Vial: 8

Acq On : 31 Mar 2017 7:30 pm

Operator: RJP

Sample : A1UG\_0.30

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 31 21:03:49 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Fri Mar 31 21:00:46 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.73	83	53117	0.31	ppb	95
47) cis-1,3-dichloropropene	13.78	75	34145	0.31	ppb	95
48) trans-1,3-dichloropropene	14.37	75	32186	0.32	ppb	94
49) 1,1,2-trichloroethane	14.57	97	25707	0.33	ppb	96
51) Toluene	14.89	92	40986	0.31	ppb	# 83
52) Methyl Isobutyl Ketone	13.82	43	42286	0.30	ppb	95
53) Dibromochloromethane	15.35	129	47684	0.30	ppb	91
54) Methyl Butyl Ketone	15.18	43	36598	0.29	ppb	96
55) 1,2-dibromoethane	15.62	107	41069	0.31	ppb	97
56) Tetrachloroethylene	16.13	164	28697	0.33	ppb	95
57) Chlorobenzene	16.86	112	55379	0.31	ppb	88
58) Ethylbenzene	17.27	91	92529	0.31	ppb	97
59) m&p-xylene	17.45	91	160464	0.63	ppb	100
60) Nonane	18.18	43	43730	0.31	ppb	91
61) Styrene	17.85	104	47576	0.30	ppb	98
62) Bromoform	17.54	173	40010	0.30	ppb	97
63) o-xylene	17.96	91	76534	0.31	ppb	93
64) Cumene	18.60	105	101572	0.32	ppb	99
66) 1,1,2,2-tetrachloroethane	17.96	83	53717	0.33	ppb	98
67) Propylbenzene	19.16	120	25611	0.31	ppb	# 1
68) 2-Chlorotoluene	19.13	126	25192	0.33	ppb	# 1
69) 4-ethyltoluene	19.32	105	90497m	0.30	ppb	
70) 1,3,5-trimethylbenzene	19.41	105	83853m	0.31	ppb	
71) 1,2,4-trimethylbenzene	19.89	105	77387	0.31	ppb	94
72) 1,3-dichlorobenzene	20.08	146	46758	0.29	ppb	99
73) benzyl chloride	20.06	91	35829	0.27	ppb	91
74) 1,4-dichlorobenzene	20.16	146	41836	0.29	ppb	99
75) 1,2,3-trimethylbenzene	20.41	105	76513	0.31	ppb	95
76) 1,2-dichlorobenzene	20.58	146	45300	0.31	ppb	97
77) 1,2,4-trichlorobenzene	22.76	180	11306	0.24	ppb	87
78) Naphthalene	22.89	128	31679	0.27	ppb	95
79) Hexachloro-1,3-butadiene	23.32	225	35411	0.31	ppb	97

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

AO033109.D A331\_1UG.M

Thu May 04 11:15:53 2017

MSD1

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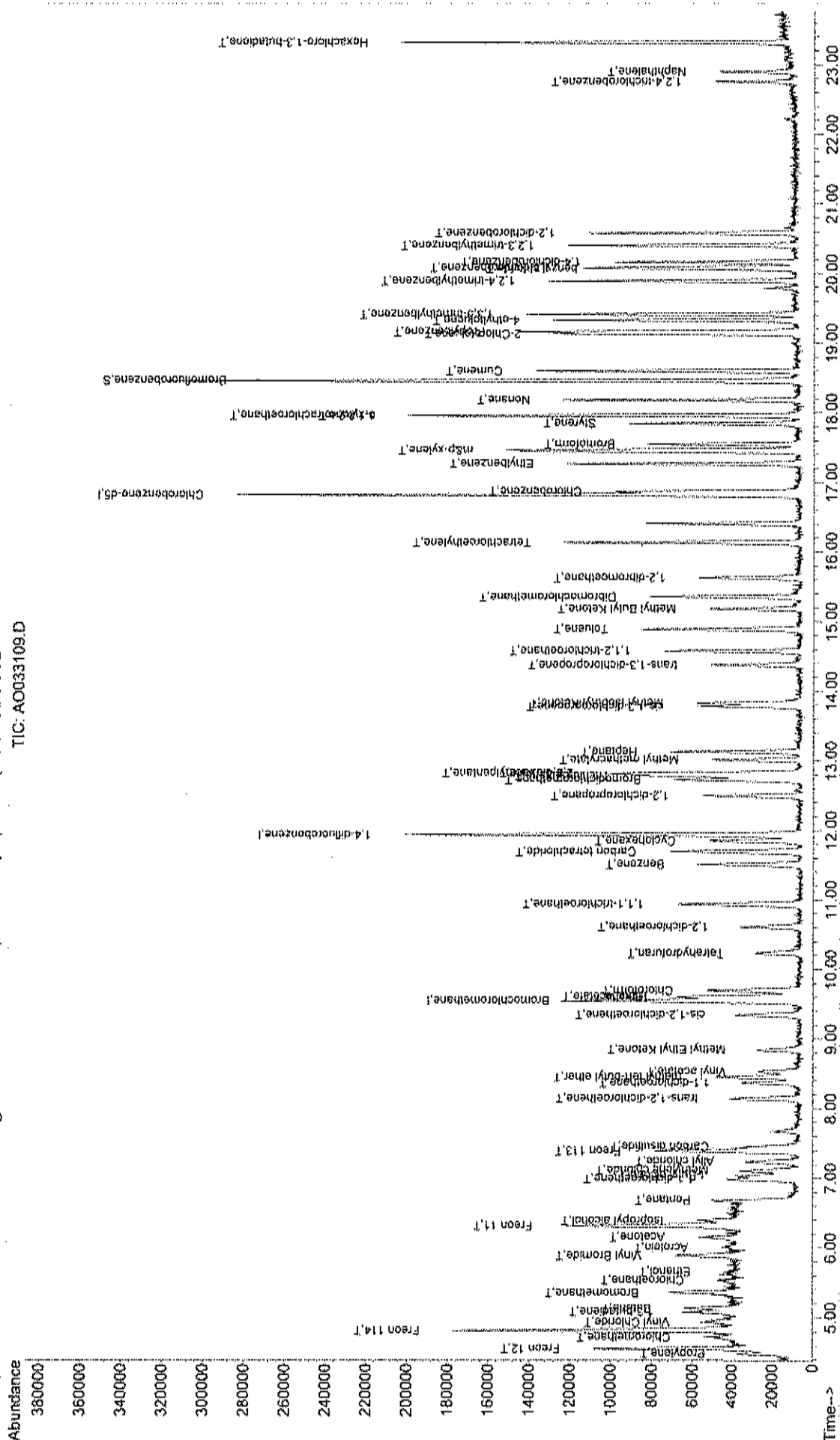
Data File : C:\HPCHEM\1\DATA\AO033109.D  
 Acq On : 31 Mar 2017 7:30 pm  
 Sample : A1UG\_0.30  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 3 10:06 2017

Vial: 8  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:13:15 2017  
 Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

TIC: AO033109.D



Data File : C:\HPCHEM\1\DATA\AO033110.D

Vial: 9

Acq On : 31 Mar 2017 8:07 pm

Operator: RJP

Sample : A1UG\_0.15

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 31 21:04:16 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Fri Mar 31 21:00:46 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	42792	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.94	114	193548	1.00	ppb	0.00
50) Chlorobenzene-d5	16.82	117	165853	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	114097	0.99	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	99.00%

## Target Compounds

						Qvalue
2) Propylene	4.49	41	9911	0.16	ppb	# 62
3) Freon 12	4.55	85	59977	0.18	ppb	99
4) Chloromethane	4.72	50	9041m	0.19	ppb	
5) Freon 114	4.82	85	46015	0.17	ppb	94
6) Vinyl Chloride	4.95	62	12362	0.18	ppb	85
7) Butane	5.14	43	14396	0.21	ppb	88
8) 1,3-butadiene	5.09	39	8893m	0.18	ppb	
9) Bromomethane	5.36	94	16794	0.17	ppb	95
10) Chloroethane	5.54	64	5195m	0.18	ppb	
11) Ethanol	5.65	45	4334m	0.19	ppb	
12) Acrolein	6.01	56	5298	0.21	ppb	88
13) Vinyl Bromide	5.89	106	17806	0.18	ppb	100
14) Freon 11	6.32	101	59840	0.17	ppb	99
15) Acetone	6.15	58	5209	0.20	ppb	# 67
16) Pentane	6.68	42	9149	0.17	ppb	# 29
17) Isopropyl alcohol	6.40	45	14755m	0.16	ppb	
18) 1,1-dichloroethene	7.00	96	9344	0.18	ppb	93
19) Freon 113	7.39	101	22230	0.17	ppb	89
20) t-Butyl alcohol	7.04	59	22290	0.17	ppb	# 91
21) Methylene chloride	7.11	84	9103	0.17	ppb	# 68
22) Allyl chloride	7.24	41	11651	0.17	ppb	87
23) Carbon disulfide	7.44	76	29034	0.17	ppb	90
24) trans-1,2-dichloroethene	8.15	61	15274	0.17	ppb	92
25) methyl tert-butyl ether	8.46	73	27786	0.16	ppb	89
26) 1,1-dichloroethane	8.38	63	19013	0.17	ppb	91
27) Vinyl acetate	8.53	43	28787	0.20	ppb	88
28) Methyl Ethyl Ketone	8.85	72	4136	0.14	ppb	# 80
29) cis-1,2-dichloroethene	9.35	61	14449	0.17	ppb	89
30) Hexane	9.60	57	12497	0.15	ppb	89
31) Ethyl acetate	9.62	43	28876	0.16	ppb	95
32) Chloroform	9.71	83	26592	0.17	ppb	96
33) Tetrahydrofuran	10.24	42	9740	0.17	ppb	92
34) 1,2-dichloroethane	10.61	62	19328	0.17	ppb	96
36) 1,1,1-trichloroethane	10.94	97	28019	0.17	ppb	99
37) Cyclohexane	11.85	56	12963	0.17	ppb	89
38) Carbon tetrachloride	11.70	117	29251	0.17	ppb	96
39) Benzene	11.51	78	29211	0.17	ppb	86
40) Methyl methacrylate	13.02	41	13355	0.16	ppb	99
41) 1,4-dioxane	12.81	58	6152	0.17	ppb	# 52
42) 2,2,4-trimethylpentane	12.82	57	43239	0.17	ppb	89
43) Heptane	13.14	43	14828	0.16	ppb	96
44) Trichloroethene	12.78	130	13911	0.17	ppb	87
45) 1,2-dichloropropane	12.51	63	11309	0.18	ppb	99

(#)= qualifier out of range (m)= manual integration

AO033110.D A331\_1UG.M

Thu May 04 11:15:57 2017

MSD1

Page 1



Data File : C:\HPCHEM\1\DATA\AO033110.D

Vial: 9

Acq On : 31 Mar 2017 8:07 pm

Operator: RJP

Sample : A1UG\_0.15

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 31 21:04:16 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Fri Mar 31 21:00:46 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.73	83	27155	0.17	ppb	94
47) cis-1,3-dichloropropene	13.78	75	18076	0.17	ppb	99
48) trans-1,3-dichloropropene	14.37	75	16213	0.17	ppb	96
49) 1,1,2-trichloroethane	14.57	97	12408	0.16	ppb	99
51) Toluene	14.88	92	20793	0.17	ppb	85
52) Methyl Isobutyl Ketone	13.83	43	20938	0.16	ppb	96
53) Dibromochloromethane	15.36	129	24776	0.17	ppb	88
54) Methyl Butyl Ketone	15.18	43	19129	0.16	ppb	94
55) 1,2-dibromoethane	15.62	107	20335	0.16	ppb	94
56) Tetrachloroethylene	16.13	164	15170	0.18	ppb	99
57) Chlorobenzene	16.86	112	28620	0.17	ppb	89
58) Ethylbenzene	17.27	91	47280	0.17	ppb	100
59) m&p-xylene	17.45	91	73840	0.30	ppb	95
60) Nonane	18.18	43	21381	0.16	ppb	92
61) Styrene	17.85	104	24592	0.16	ppb	93
62) Bromoform	17.55	173	19862	0.16	ppb	96
63) o-xylene	17.96	91	37803	0.16	ppb	96
64) Cumene	18.60	105	50505	0.17	ppb	98
66) 1,1,2,2-tetrachloroethane	17.96	83	27102	0.17	ppb	96
67) Propylbenzene	19.16	120	13021	0.17	ppb	# 1
68) 2-Chlorotoluene	19.13	126	11618	0.16	ppb	# 1
69) 4-ethyltoluene	19.33	105	44953	0.16	ppb	98
70) 1,3,5-trimethylbenzene	19.41	105	38812	0.15	ppb	98
71) 1,2,4-trimethylbenzene	19.89	105	35741	0.15	ppb	88
72) 1,3-dichlorobenzene	20.08	146	20897	0.14	ppb	97
73) benzyl chloride	20.06	91	17111	0.13	ppb	95
74) 1,4-dichlorobenzene	20.16	146	20124	0.15	ppb	94
75) 1,2,3-trimethylbenzene	20.41	105	37732	0.16	ppb	97
76) 1,2-dichlorobenzene	20.58	146	22151	0.16	ppb	96
77) 1,2,4-trichlorobenzene	22.76	180	4904	0.11	ppb	91
78) Naphthalene	22.90	128	12548	0.11	ppb	90
79) Hexachloro-1,3-butadiene	23.33	225	16218	0.15	ppb	98

Time (h)	10.00	9.00	8.00	7.00	6.00	5.00
23.00						
22.00						
21.00						
20.00						
19.00						
18.00						
17.00						
16.00						
15.00						
14.00						
13.00						
12.00						
11.00						
10.00						

Data File : C:\HPCHEM\1\DATA\AO033111.D

Vial: 10

Acq On : 31 Mar 2017 8:44 pm

Operator: RJP

Sample : A1UG\_0.10

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 31 22:13:02 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Fri Mar 31 21:00:46 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.55	128	41392	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.94	114	192052	1.00	ppb	0.00
50) Chlorobenzene-d5	16.82	117	166270	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	108405	0.94	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	94.00%

## Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
6) Vinyl Chloride	4.94	62	8904	0.13	ppb	66
38) Carbon tetrachloride	11.69	117	19812	0.12	ppb	96
44) Trichloroethene	12.78	130	9232	0.11	ppb	89
56) Tetrachloroethylene	16.14	164	10008	0.12	ppb	93

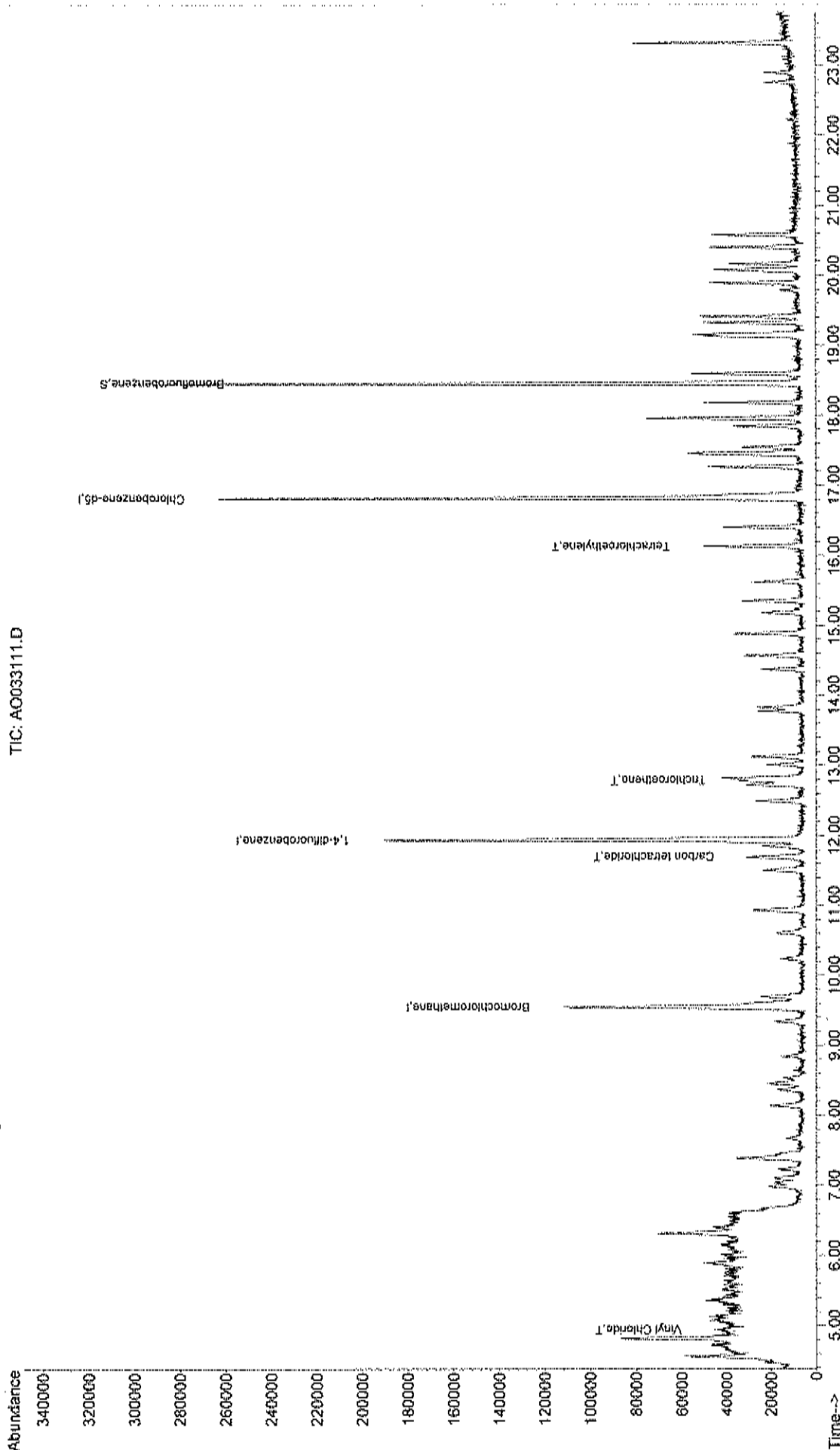
Data File : C:\HPCHEM\1\DATA\AO033111.D  
 Acq On : 31 Mar 2017 8:44 pm  
 Sample : A1UG 0.10  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 3 10:09 2017

Vial: 10  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:13:15 2017  
 Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

TIC: AO033111.D



Data File : C:\HPCHEM\1\DATA\AO033112.D

Vial: 11

Acq On : 31 Mar 2017 9:20 pm

Operator: RJP

Sample : A1UG\_0.04

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 31 22:13:36 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Fri Mar 31 21:00:46 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.55	128	39334	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.94	114	186827	1.00	ppb	0.00
50) Chlorobenzene-d5	16.82	117	162286	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.45	95	106786	0.95	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	95.00%

## Target Compounds

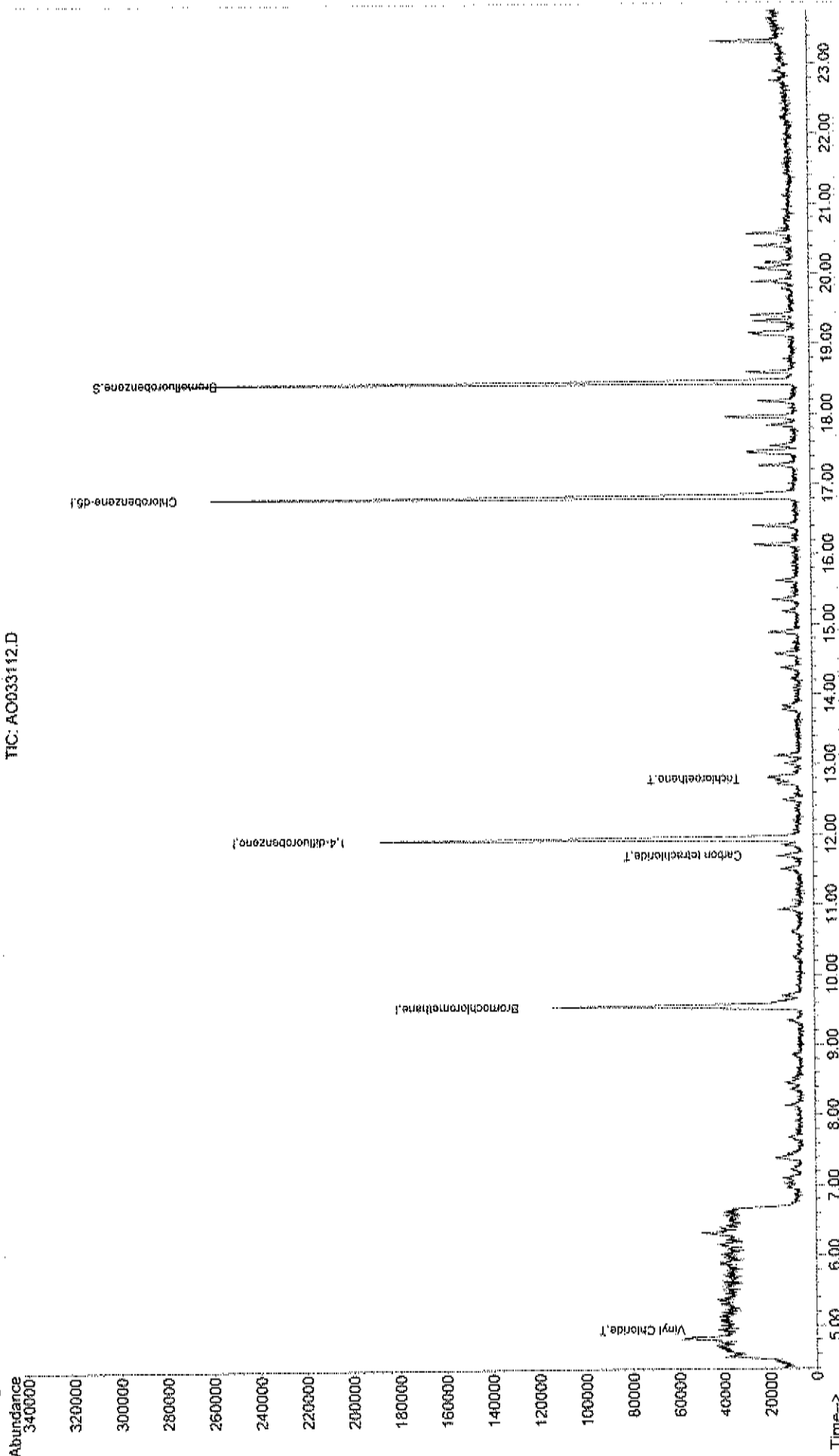
	R.T.	QIon	Response	Conc	Units	Qvalue
6) Vinyl Chloride	4.93	62	3087	0.05	ppb	89
38) Carbon tetrachloride	11.70	117	7584	0.05	ppb	98
44) Trichloroethene	12.79	130	3902	0.05	ppb	# 23

Data File : C:\HPCHEM\1\DATA\AO033112.D  
 Acq On : 31 Mar 2017 9:20 pm  
 Sample : A1UG 0.04  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 3 10:11 2017

Vial: 11  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:13:15 2017  
 Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO033106.D  
 TIC: AO033112.D



**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**CALIBRATION VERIFICATION**



## Evaluate Continuing Calibration Report

Data File : C:\HPCHEM\1\DATA\AO040602.D

Vial: 2

Acq On : 6 Apr 2017 9:31 am

Operator: RJP

Sample : A1UG\_1.0

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Thu May 04 11:27:28 2017

Response via : Multiple Level Calibration

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

Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane	1.000	1.000	0.0	58	0.00
2 T	Propylene	1.425	1.538	-7.9	62	0.00
3 T	Freon 12	8.232	7.850	4.6	57	0.00
4 T	Chloromethane	1.163	1.228	-5.6	63	0.00
5 T	Freon 114	6.407	6.744	-5.3	63	0.00
6 T	Vinyl Chloride	1.763	1.645	6.7	60	0.00
7 T	Butane	1.798	1.735	3.5	62	0.00
8 T	1,3-butadiene	1.165	1.240	-6.4	67	0.00
9 T	Bromomethane	2.375	2.252	5.2	57	0.00
10 T	Chloroethane	0.733	0.732	0.1	62	0.00
11 T	Ethanol	0.576	0.494	14.2	52	0.00
12 T	Acrolein	0.604	0.474	21.5	46#	0.00
13 T	Vinyl Bromide	2.385	2.040	14.5	50	0.00
14 T	Freon 11	8.237	10.020	-21.6	72	0.00
15 T	Acetone	0.645	0.664	-2.9	64	0.00
16 T	Pentane	1.262	1.246	1.3	57	0.00
17 T	Isopropyl alcohol	2.143	2.170	-1.3	59	0.00
18 T	1,1-dichloroethene	1.300	1.185	8.8	56	0.00
19 T	Freon 113	3.107	3.062	1.4	60	0.00
20 t	t-Butyl alcohol	3.171	2.505	21.0	46#	0.00
21 T	Methylene chloride	1.268	1.230	3.0	58	0.00
22 T	Allyl chloride	1.615	1.358	15.9	50#	0.00
23 T	Carbon disulfide	4.058	3.972	2.1	59	0.00
24 T	trans-1,2-dichloroethene	2.189	2.057	6.0	57	0.00
25 T	methyl tert-butyl ether	4.148	3.676	11.4	52	0.00
26 T	1,1-dichloroethane	2.710	2.632	2.9	57	0.00
27 T	Vinyl acetate	3.495	3.889	-11.3	70	0.00
28 T	Methyl Ethyl Ketone	0.665	0.614	7.7	52	0.00
29 T	cis-1,2-dichloroethene	2.077	1.918	7.7	55	0.00
30 T	Hexane	1.907	1.788	6.2	55	0.00
31 T	Ethyl acetate	4.254	3.890	8.6	55	0.00
32 T	Chloroform	3.793	3.590	5.4	56	0.00
33 T	Tetrahydrofuran	1.421	1.284	9.6	55	0.00
34 T	1,2-dichloroethane	2.807	2.513	10.5	53	0.00
35 I	1,4-difluorobenzene	1.000	1.000	0.0	55	0.00
36 T	1,1,1-trichloroethane	0.880	0.890	-1.1	58	0.00
37 T	Cyclohexane	0.444	0.426	4.1	59	0.00
38 T	Carbon tetrachloride	0.932	0.926	0.6	58	0.00
39 T	Benzene	0.920	0.911	1.0	58	0.00
40 T	Methyl methacrylate	0.431	0.389	9.7	51	0.00
41 T	1,4-dioxane	0.190	0.162	14.7	49#	0.00
42 T	2,2,4-trimethylpentane	1.383	1.364	1.4	57	0.00
43 T	Heptane	0.480	0.462	3.7	55	0.00
44 T	Trichloroethene	0.450	0.443	1.6	58	0.00
45 T	1,2-dichloropropane	0.339	0.348	-2.7	59	0.00
46 T	Bromodichloromethane	0.873	0.877	-0.5	58	0.00
47 T	cis-1,3-dichloropropene	0.559	0.551	1.4	56	0.00
48 T	trans-1,3-dichloropropene	0.532	0.524	1.5	57	0.00
49 T	1,1,2-trichloroethane	0.407	0.427	-4.9	60	0.00

-----  
(#) = Out of Range

AO040602.D A331\_1UG.M

Thu May 04 11:59:46 2017

MSD1

Page 1

## Evaluate Continuing Calibration Report

Data File : C:\HPCHEM\1\DATA\AO040602.D

Vial: 2

Acq On : 6 Apr 2017 9:31 am

Operator: RJP

Sample : A1UG\_1.0

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Thu May 04 11:27:28 2017

Response via : Multiple Level Calibration

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

Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
51 T	Toluene	0.750	0.686	8.5	55	0.00
52 T	Methyl Isobutyl Ketone	0.780	0.558	28.5	42#	0.00
53 T	Dibromochloromethane	0.917	0.883	3.7	59	0.00
54 T	Methyl Butyl Ketone	0.702	0.528	24.8	44#	0.00
55 T	1,2-dibromoethane	0.761	0.740	2.8	58	0.00
56 T	Tetrachloroethylene	0.529	0.507	4.2	61	0.00
57 T	Chlorobenzene	1.044	0.978	6.3	57	0.00
58 T	Ethylbenzene	1.750	1.612	7.9	57	0.00
59 T	m&p-xylene	1.477	1.367	7.4	56	0.00
60 T	Nonane	0.818	0.751	8.2	55	0.00
61 T	Styrene	0.919	0.821	10.7	55	0.00
62 T	Bromoform	0.761	0.735	3.4	59	0.00
63 T	o-xylene	1.424	1.330	6.6	57	0.00
64 T	Cumene	1.851	1.691	8.6	56	0.00
65 S	Bromofluorobenzene	0.682	0.680	0.3	59	0.00
66 T	1,1,2,2-tetrachloroethane	0.973	0.966	0.7	62	0.00
67 T	Propylbenzene	0.476	0.460	3.4	58	0.00
68 T	2-Chlorotoluene	0.440	0.428	2.7	59	0.00
69 T	4-ethyltoluene	1.729	1.571	9.1	54	0.00
70 T	1,3,5-trimethylbenzene	1.538	1.445	6.0	56	0.00
71 T	1,2,4-trimethylbenzene	1.434	1.300	9.3	55	0.00
72 T	1,3-dichlorobenzene	0.902	0.907	-0.6	59	0.00
73 T	benzyl chloride	0.755	0.754	0.1	59	0.00
74 T	1,4-dichlorobenzene	0.834	0.822	1.4	60	0.00
75 T	1,2,3-trimethylbenzene	1.443	1.318	8.7	56	0.00
76 T	1,2-dichlorobenzene	0.869	0.850	2.2	60	0.00
77 T	1,2,4-trichlorobenzene	0.257	0.268	-4.3	60	0.00
78 T	Naphthalene	0.661	0.565	14.5	51	0.00
79 T	Hexachloro-1,3-butadiene	0.644	0.635	1.4	59	0.00

Data File : C:\HPCHEM\1\DATA\A0040602.D

Vial: 2

Acq On : 6 Apr 2017 9:31 am

Operator: RJP

Sample : A1UG\_1.0

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 06 11:05:43 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.57	128	26713	1.00	ppb	0.03
35) 1,4-difluorobenzene	11.95	114	121699	1.00	ppb	0.02
50) Chlorobenzene-d5	16.83	117	111761	1.00	ppb	0.01

## System Monitoring Compounds

65) Bromofluorobenzene	18.47	95	75986	1.00	ppb	0.02
Spiked Amount	1.000	Range	70 - 130	Recovery	=	100.00%

## Target Compounds

						Qvalue
2) Propylene	4.49	41	41087	1.08	ppb	85
3) Freon 12	4.57	85	209693	0.95	ppb	98
4) Chloromethane	4.73	50	32806	1.06	ppb	87
5) Freon 114	4.83	85	180160	1.05	ppb	94
6) Vinyl Chloride	4.97	62	43942	0.93	ppb	93
7) Butane	5.16	43	46357	0.97	ppb	87
8) 1,3-butadiene	5.09	39	33128m	1.06	ppb	
9) Bromomethane	5.38	94	60166	0.95	ppb	89
10) Chloroethane	5.56	64	19566	1.00	ppb	91
11) Ethanol	5.68	45	13202	0.86	ppb	85
12) Acrolein	6.02	56	12651	0.78	ppb	99
13) Vinyl Bromide	5.91	106	54506	0.86	ppb	95
14) Freon 11	6.33	101	267654	1.22	ppb	99
15) Acetone	6.17	58	17734	1.03	ppb	# 78
16) Pentane	6.69	42	33284	0.99	ppb	# 45
17) Isopropyl alcohol	6.41	45	57975	1.01	ppb	# 100
18) 1,1-dichloroethene	7.00	96	31668	0.91	ppb	89
19) Freon 113	7.41	101	81782	0.99	ppb	91
20) t-Butyl alcohol	7.05	59	66928	0.79	ppb	# 92
21) Methylene chloride	7.12	84	32863	0.97	ppb	85
22) Allyl chloride	7.25	41	36285	0.84	ppb	93
23) Carbon disulfide	7.47	76	106097	0.98	ppb	84
24) trans-1,2-dichloroethene	8.17	61	54946	0.94	ppb	93
25) methyl tert-butyl ether	8.47	73	98209	0.89	ppb	87
26) 1,1-dichloroethane	8.40	63	70315	0.97	ppb	98
27) Vinyl acetate	8.55	43	103887	1.11	ppb	91
28) Methyl Ethyl Ketone	8.87	72	16394	0.92	ppb	98
29) cis-1,2-dichloroethene	9.37	61	51232	0.92	ppb	95
30) Hexane	9.63	57	47775	0.94	ppb	89
31) Ethyl acetate	9.63	43	103920	0.91	ppb	98
32) Chloroform	9.72	83	95910	0.95	ppb	96
33) Tetrahydrofuran	10.25	42	34290	0.90	ppb	87
34) 1,2-dichloroethane	10.63	62	67117	0.90	ppb	95
36) 1,1,1-trichloroethane	10.96	97	108277	1.01	ppb	95
37) Cyclohexane	11.87	56	51864	0.96	ppb	98
38) Carbon tetrachloride	11.71	117	112693	0.99	ppb	98
39) Benzene	11.52	78	110859	0.99	ppb	91
40) Methyl methacrylate	13.03	41	47367	0.90	ppb	96
41) 1,4-dioxane	12.83	58	19760	0.85	ppb	# 61
42) 2,2,4-trimethylpentane	12.84	57	165964	0.99	ppb	90
43) Heptane	13.14	43	56271	0.96	ppb	95
44) Trichloroethene	12.80	130	53897	0.98	ppb	93
45) 1,2-dichloropropane	12.51	63	42378	1.03	ppb	99

(#)= qualifier out of range (m) = manual integration

A0040602.D A331\_1UG.M

Thu May 04 11:59:53 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO040602.D

Vial: 2

Acq On : 6 Apr 2017 9:31 am

Operator: RJP

Sample : A1UG\_1.0

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 06 11:05:43 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.74	83	106680	1.00	ppb	96
47) cis-1,3-dichloropropene	13.80	75	67019	0.99	ppb	97
48) trans-1,3-dichloropropene	14.38	75	63822	0.99	ppb	95
49) 1,1,2-trichloroethane	14.58	97	51916	1.05	ppb	97
51) Toluene	14.90	92	76639	0.91	ppb	# 83
52) Methyl Isobutyl Ketone	13.84	43	62379m	0.72	ppb	
53) Dibromochloromethane	15.37	129	98703	0.96	ppb	93
54) Methyl Butyl Ketone	15.20	43	59023m	0.75	ppb	
55) 1,2-dibromoethane	15.64	107	82733	0.97	ppb	95
56) Tetrachloroethylene	16.15	164	56610	0.96	ppb	99
57) Chlorobenzene	16.87	112	109355	0.94	ppb	95
58) Ethylbenzene	17.27	91	180188	0.92	ppb	97
59) m&p-xylene	17.46	91	305465	1.85	ppb	100
60) Nonane	18.19	43	83926	0.92	ppb	93
61) Styrene	17.85	104	91731	0.89	ppb	98
62) Bromoform	17.56	173	82114	0.97	ppb	100
63) o-xylene	17.97	91	148687	0.93	ppb	93
64) Cumene	18.61	105	189034	0.91	ppb	98
66) 1,1,2,2-tetrachloroethane	17.96	83	107989	0.99	ppb	97
67) Propylbenzene	19.18	120	51380	0.97	ppb	# 1
68) 2-Chlorotoluene	19.14	126	47884	0.97	ppb	# 1
69) 4-ethyltoluene	19.34	105	175572	0.91	ppb	100
70) 1,3,5-trimethylbenzene	19.42	105	161521	0.94	ppb	99
71) 1,2,4-trimethylbenzene	19.90	105	145281	0.91	ppb	93
72) 1,3-dichlorobenzene	20.09	146	101408	1.01	ppb	97
73) benzyl chloride	20.06	91	84267	1.00	ppb	94
74) 1,4-dichlorobenzene	20.17	146	91864	0.99	ppb	95
75) 1,2,3-trimethylbenzene	20.41	105	147288	0.91	ppb	93
76) 1,2-dichlorobenzene	20.59	146	94991	0.98	ppb	96
77) 1,2,4-trichlorobenzene	22.76	180	29991	1.05	ppb	99
78) Naphthalene	22.91	128	63149	0.86	ppb	94
79) Hexachloro-1,3-butadiene	23.33	225	70991	0.99	ppb	97

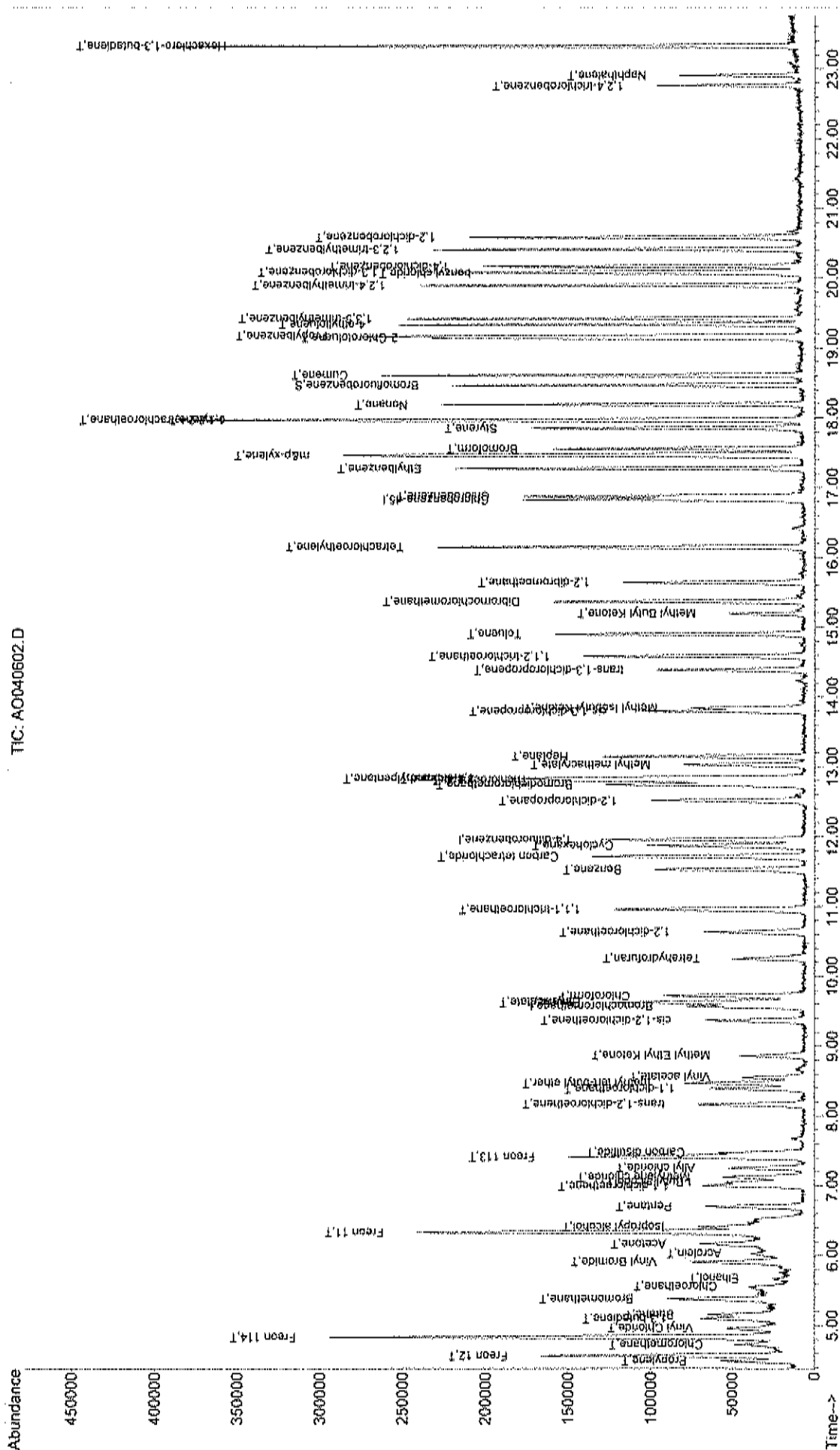
-----  
 (#) = qualifier out of range (m) = manual integration (+) = signals summed  
 AO040602.D A331\_1UG.M Thu May 04 11:59:54 2017 MSD1

Data File : C:\HPCHEM\1\DATA\A0040602.D  
Acq On : 6 Apr 2017 9:31 am  
Sample : A1UG\_1.0  
Misc : A331\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Apr 7 10:14 2017

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Thu May 04 11:27:28 2017  
Response via : Initial Calibration

Vial: 2  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A331\_1UG.RES



## Evaluate Continuing Calibration Report

Data File : C:\HPCHEM\1\DATA\A0040702.D  
 Acq On : 7 Apr 2017 11:20 am  
 Sample : A1UG\_1.0  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P

Vial: 2  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min  
 Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane	1.000	1.000	0.0	50	-0.02
2 T	Propylene	1.425	1.469	-3.1	52	0.00
3 T	Freon 12	8.232	9.830	-19.4	62	-0.01
4 T	Chloromethane	1.163	1.259	-8.3	56	-0.02
5 T	Freon 114	6.407	7.985	-24.6	65	-0.02
6 T	Vinyl Chloride	1.763	1.981	-12.4	62	-0.02
7 T	Butane	1.798	2.051	-14.1	64	-0.02
8 T	1,3-butadiene	1.165	1.307	-12.2	62	-0.01
9 T	Bromomethane	2.375	2.910	-22.5	64	-0.01
10 T	Chloroethane	0.733	0.845	-15.3	62	-0.03
11 T	Ethanol	0.576	0.568	1.4	52	-0.04
12 T	Acrolein	0.604	0.663	-9.8	56	-0.02
13 T	Vinyl Bromide	2.385	2.836	-18.9	61	-0.02
14 T	Freon 11	8.237	10.475	-27.2	66	0.00
15 T	Acetone	0.645	0.792	-22.8	67	-0.02
16 T	Pentane	1.262	1.188	5.9	47#	0.00
17 T	Isopropyl alcohol	2.143	2.346	-9.5	55	-0.03
18 T	1,1-dichloroethene	1.300	1.339	-3.0	55	-0.02
19 T	Freon 113	3.107	3.265	-5.1	55	-0.01
20 T	t-Butyl alcohol	3.171	2.625	17.2	42#	-0.02
21 T	Methylene chloride	1.268	1.314	-3.6	54	-0.01
22 T	Allyl chloride	1.615	1.415	12.4	45#	0.00
23 T	Carbon disulfide	4.058	4.135	-1.9	53	-0.01
24 T	trans-1,2-dichloroethene	2.189	2.150	1.8	52	-0.02
25 T	methyl tert-butyl ether	4.148	3.833	7.6	48#	-0.02
26 T	1,1-dichloroethane	2.710	2.713	-0.1	51	-0.02
27 T	Vinyl acetate	3.495	3.170	9.3	50#	-0.02
28 T	Methyl Ethyl Ketone	0.665	0.594	10.7	43#	-0.03
29 T	cis-1,2-dichloroethene	2.077	2.006	3.4	50	-0.02
30 T	Hexane	1.907	1.835	3.8	49#	-0.02
31 T	Ethyl acetate	4.254	3.961	6.9	49#	-0.02
32 T	Chloroform	3.793	3.922	-3.4	53	0.00
33 T	Tetrahydrofuran	1.421	1.277	10.1	47#	-0.02
34 T	1,2-dichloroethane	2.807	2.738	2.5	50	-0.01
35 I	1,4-difluorobenzene	1.000	1.000	0.0	45#	0.00
36 T	1,1,1-trichloroethane	0.880	1.005	-14.2	54	-0.01
37 T	Cyclohexane	0.444	0.448	-0.9	51	0.00
38 T	Carbon tetrachloride	0.932	1.012	-8.6	52	0.00
39 T	Benzene	0.920	1.006	-9.3	52	0.00
40 T	Methyl methacrylate	0.431	0.421	2.3	45#	-0.01
41 T	1,4-dioxane	0.190	0.186	2.1	46#	0.00
42 T	2,2,4-trimethylpentane	1.383	1.522	-10.1	52	0.00
43 T	Heptane	0.480	0.531	-10.6	51	0.00
44 T	Trichloroethene	0.450	0.502	-11.6	54	0.00
45 T	1,2-dichloropropane	0.339	0.391	-15.3	54	0.00
46 T	Bromodichloromethane	0.873	0.977	-11.9	52	0.00
47 T	cis-1,3-dichloropropene	0.559	0.609	-8.9	51	-0.01
48 T	trans-1,3-dichloropropene	0.532	0.577	-8.5	52	0.00
49 T	1,1,2-trichloroethane	0.407	0.462	-13.5	53	0.00

(#) = Out of Range

A0040702.D A331\_1UG.M

Thu May 04 12:01:41 2017

MSD1

Page 1

## Evaluate Continuing Calibration Report

Data File : C:\HPCHEM\1\DATA\AO040702.D

Vial: 2

Acq On : 7 Apr 2017 11:20 am

Operator: RJP

Sample : A1UG\_1.0

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Thu May 04 11:27:28 2017

Response via : Multiple Level Calibration

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

Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
51 T	Toluene	0.750	0.788	-5.1	50#	0.00
52 T	Methyl Isobutyl Ketone	0.780	0.724	7.2	43#	-0.02
53 T	Dibromochloromethane	0.917	1.003	-9.4	53	0.00
54 T	Methyl Butyl Ketone	0.702	0.499	28.9	33#	-0.02
55 T	1,2-dibromoethane	0.761	0.829	-8.9	52	0.00
56 T	Tetrachloroethylene	0.529	0.570	-7.8	54	0.00
57 T	Chlorobenzene	1.044	1.168	-11.9	54	0.00
58 T	Ethylbenzene	1.750	1.875	-7.1	53	0.00
59 T	m&p-xylene	1.477	1.572	-6.4	51	0.00
60 T	Nonane	0.818	0.866	-5.9	51	0.00
61 T	Styrene	0.919	0.969	-5.4	51	0.00
62 T	Bromoform	0.761	0.848	-11.4	54	0.00
63 T	o-xylene	1.424	1.535	-7.8	52	0.00
64 T	Cumene	1.851	1.947	-5.2	51	0.00
65 S	Bromofluorobenzene	0.682	0.628	7.9	43#	0.00
66 T	1,1,2,2-tetrachloroethane	0.973	1.144	-17.6	58	0.00
67 T	Propylbenzene	0.476	0.524	-10.1	53	0.00
68 T	2-Chlorotoluene	0.440	0.474	-7.7	52	0.00
69 T	4-ethyltoluene	1.729	1.869	-8.1	51	0.00
70 T	1,3,5-trimethylbenzene	1.538	1.571	-2.1	48#	0.00
71 T	1,2,4-trimethylbenzene	1.434	1.465	-2.2	49#	0.00
72 T	1,3-dichlorobenzene	0.902	1.068	-18.4	56	0.00
73 T	benzyl chloride	0.755	0.804	-6.5	50	0.00
74 T	1,4-dichlorobenzene	0.834	0.944	-13.2	54	0.00
75 T	1,2,3-trimethylbenzene	1.443	1.458	-1.0	49#	0.00
76 T	1,2-dichlorobenzene	0.869	0.976	-12.3	55	0.00
77 T	1,2,4-trichlorobenzene	0.257	0.273	-6.2	49#	0.00
78 T	Naphthalene	0.661	0.561	15.1	40#	0.00
79 T	Hexachloro-1,3-butadiene	0.644	0.689	-7.0	51	0.00



Data File : C:\HPCHEM\1\DATA\A0040702.D

Acq On : 7 Apr 2017 11:20 am

Sample : A1UG\_1.0

Misc : A331\_1UG

MS Integration Params: RTEINT.P

Quant Time: Apr 07 12:17:46 2017

Vial: 2

Operator: RJP

Inst : MSD #1

Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	9.55	128	23210m	1.00	ppb	0.01
35) 1,4-difluorobenzene	11.95	114	99215	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	88822	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	55801	0.92	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	92.00%

## Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Propylene	4.48	41	34098	1.03	ppb	77
3) Freon 12	4.56	85	228157	1.19	ppb	99
4) Chloromethane	4.71	50	29211m	1.08	ppb	
5) Freon 114	4.82	85	185325m	1.25	ppb	
6) Vinyl Chloride	4.95	62	45978	1.12	ppb	88
7) Butane	5.13	43	47604	1.14	ppb	91
8) 1,3-butadiene	5.08	39	30333m	1.12	ppb	
9) Bromomethane	5.36	94	67550	1.23	ppb	87
10) Chloroethane	5.53	64	19623	1.15	ppb	# 81
11) Ethanol	5.64	45	13180m	0.99	ppb	
12) Acrolein	6.00	56	15386	1.10	ppb	83
13) Vinyl Bromide	5.89	106	65815	1.19	ppb	94
14) Freon 11	6.32	101	243134m	1.27	ppb	
15) Acetone	6.15	58	18392	1.23	ppb	# 73
16) Pentane	6.69	42	27574	0.94	ppb	# 32
17) Isopropyl alcohol	6.38	45	54446	1.09	ppb	# 100
18) 1,1-dichloroethene	6.98	96	31071	1.03	ppb	95
19) Freon 113	7.40	101	75784	1.05	ppb	90
20) t-Butyl alcohol	7.03	59	60918	0.83	ppb	# 90
21) Methylene chloride	7.11	84	30497	1.04	ppb	# 74
22) Allyl chloride	7.24	41	32834	0.88	ppb	92
23) Carbon disulfide	7.46	76	95964	1.02	ppb	82
24) trans-1,2-dichloroethene	8.14	61	49908	0.98	ppb	94
25) methyl tert-butyl ether	8.45	73	88965	0.92	ppb	89
26) 1,1-dichloroethane	8.38	63	62978	1.00	ppb	94
27) Vinyl acetate	8.54	43	73585m	0.91	ppb	
28) Methyl Ethyl Ketone	8.83	72	13784	0.89	ppb	96
29) cis-1,2-dichloroethene	9.35	61	46549	0.97	ppb	97
30) Hexane	9.61	57	42588	0.96	ppb	88
31) Ethyl acetate	9.62	43	91942	0.93	ppb	97
32) Chloroform	9.71	83	91036	1.03	ppb	97
33) Tetrahydrofuran	10.23	42	29641	0.90	ppb	89
34) 1,2-dichloroethane	10.62	62	63539	0.98	ppb	98
36) 1,1,1-trichloroethane	10.95	97	99692	1.14	ppb	95
37) Cyclohexane	11.87	56	44460	1.01	ppb	96
38) Carbon tetrachloride	11.71	117	100429	1.09	ppb	97
39) Benzene	11.52	78	99793	1.09	ppb	90
40) Methyl methacrylate	13.02	41	41807	0.98	ppb	96
41) 1,4-dioxane	12.82	58	18454	0.98	ppb	# 62
42) 2,2,4-trimethylpentane	12.83	57	151031	1.10	ppb	88
43) Heptane	13.14	43	52704	1.11	ppb	96
44) Trichloroethene	12.79	130	49825	1.12	ppb	93
45) 1,2-dichloropropane	12.50	63	38777	1.15	ppb	97

(#)=qualifier out of range (m)=manual integration

A0040702.D A331\_1UG.M

Thu May 04 12:01:45 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO040702.D

Vial: 2

Acq On : 7 Apr 2017 11:20 am

Operator: RJP

Sample : A1UG\_1.0

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 07 12:17:46 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

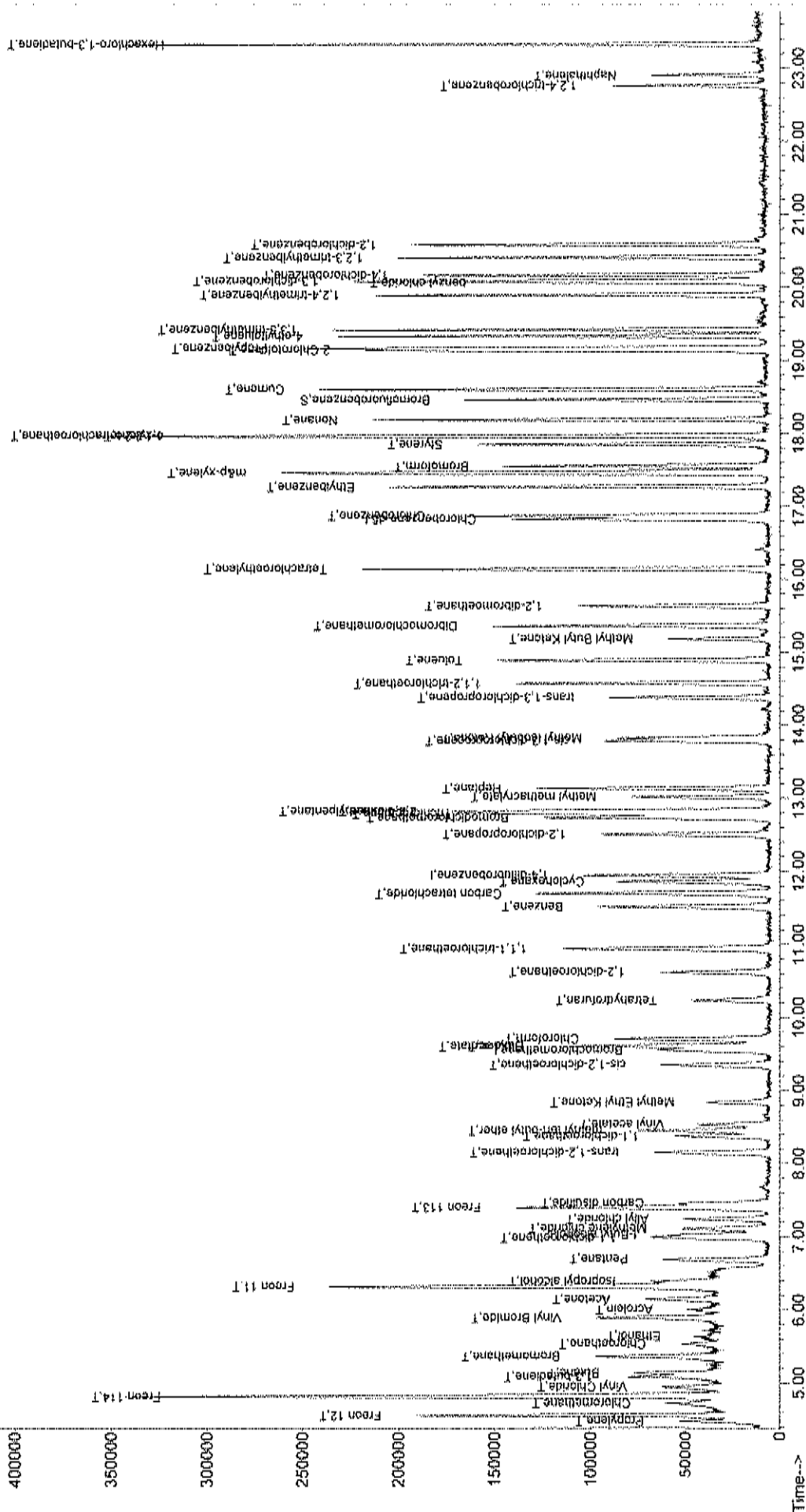
Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.74	83	96927	1.12	ppb	97
47) cis-1,3-dichloropropene	13.79	75	60416	1.09	ppb	96
48) trans-1,3-dichloropropene	14.38	75	57284	1.09	ppb	93
49) 1,1,2-trichloroethane	14.58	97	45788	1.13	ppb	96
51) Toluene	14.90	92	69963	1.05	ppb	# 79
52) Methyl Isobutyl Ketone	13.82	43	64328	0.93	ppb	90
53) Dibromochloromethane	15.36	129	89092	1.09	ppb	90
54) Methyl Butyl Ketone	15.18	43	44289	0.71	ppb	91
55) 1,2-dibromoethane	15.64	107	73621	1.09	ppb	92
56) Tetrachloroethylene	16.14	164	50632	1.08	ppb	98
57) Chlorobenzene	16.87	112	103742	1.12	ppb	98
58) Ethylbenzene	17.27	91	166527	1.07	ppb	100
59) m&p-xylene	17.47	91	279261	2.13	ppb	100
60) Nonane	18.19	43	76907	1.06	ppb	95
61) Styrene	17.85	104	86100	1.05	ppb	92
62) Bromoform	17.56	173	75313	1.11	ppb	100
63) o-xylene	17.97	91	136347	1.08	ppb	92
64) Cumene	18.60	105	172917	1.05	ppb	100
66) 1,1,2,2-tetrachloroethane	17.96	83	101634	1.18	ppb	97
67) Propylbenzene	19.17	120	46509	1.10	ppb	# 1
68) 2-Chlorotoluene	19.14	126	42111	1.08	ppb	# 1
69) 4-ethyltoluene	19.33	105	166042	1.08	ppb	99
70) 1,3,5-trimethylbenzene	19.42	105	139523	1.02	ppb	97
71) 1,2,4-trimethylbenzene	19.90	105	130111	1.02	ppb	95
72) 1,3-dichlorobenzene	20.09	146	94874	1.18	ppb	97
73) benzyl chloride	20.06	91	71400	1.06	ppb	93
74) 1,4-dichlorobenzene	20.17	146	83841	1.13	ppb	98
75) 1,2,3-trimethylbenzene	20.41	105	129500	1.01	ppb	96
76) 1,2-dichlorobenzene	20.59	146	86684	1.12	ppb	98
77) 1,2,4-trichlorobenzene	22.76	180	24245	1.06	ppb	95
78) Naphthalene	22.90	128	49826	0.85	ppb	92
79) Hexachloro-1,3-butadiene	23.32	225	61181	1.07	ppb	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed  
 AO040702.D A331\_1UG.M Thu May 04 12:01:45 2017 MSD1

**សេចក្តីផ្តើម**



**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**RAW DATA**

Data File : C:\HPCHEM\1\DATA\AO033101.D

Vial: 1

Acq On : 31 Mar 2017 1:54 pm

Operator: RJP

Sample : BFB1UG

Inst : MSD #1

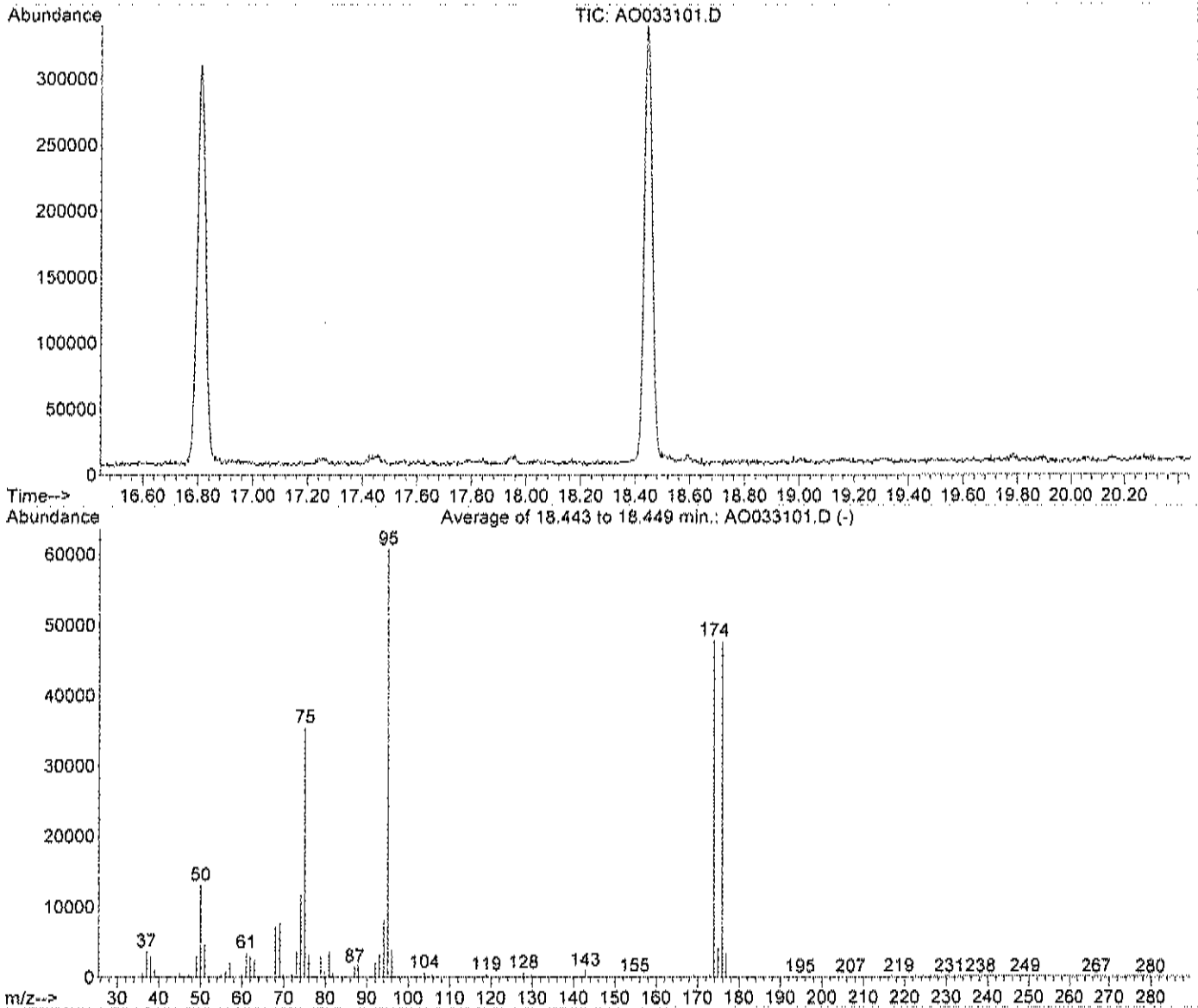
Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration



Spectrum Information: Average of 18.443 to 18.449 min.

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	8	40	21.6	13135	PASS
75	95	30	66	58.4	35541	PASS
95	95	100	100	100.0	60837	PASS
96	95	5	9	6.3	3857	PASS
173	174	0.00	2	0.7	324	PASS
174	95	50	120	78.6	47805	PASS
175	174	4	9	8.5	4060	PASS
176	174	95	101	99.4	47501	PASS
177	176	5	9	7.0	3306	PASS

Data File : C:\HPCHEM\1\DATA\AO040601.D

Vial: 1

Acq On : 6 Apr 2017 8:51 am

Operator: RJP

Sample : BFB1UG

Inst : MSD #1

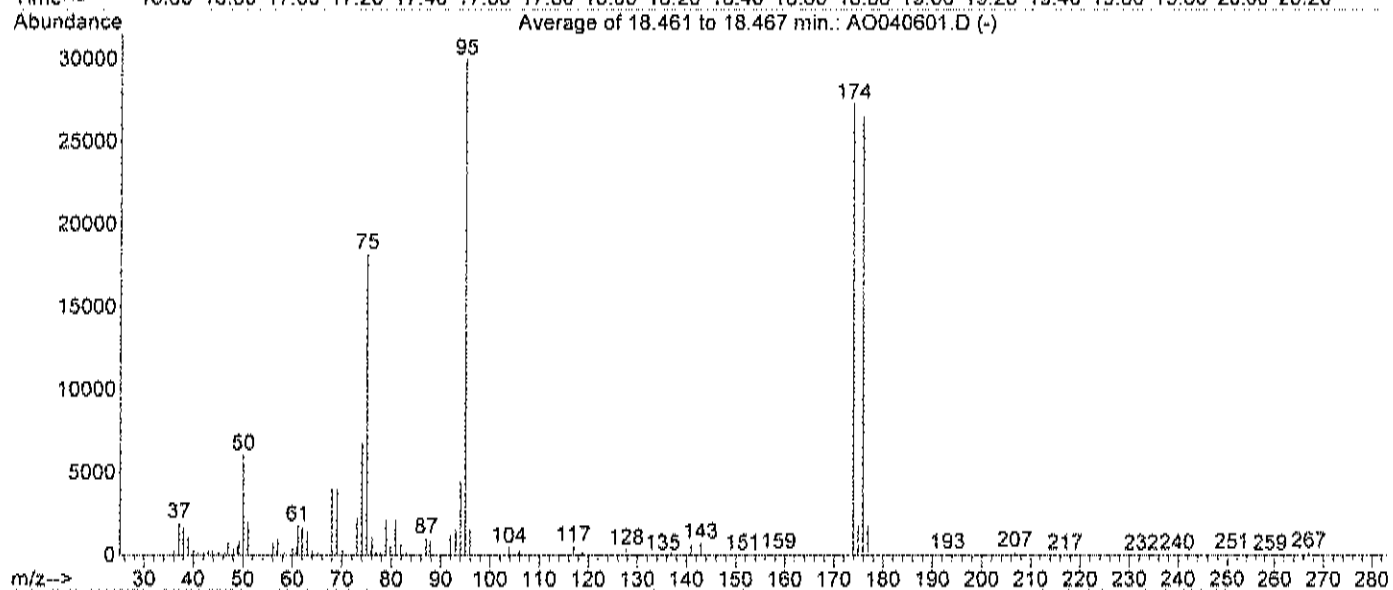
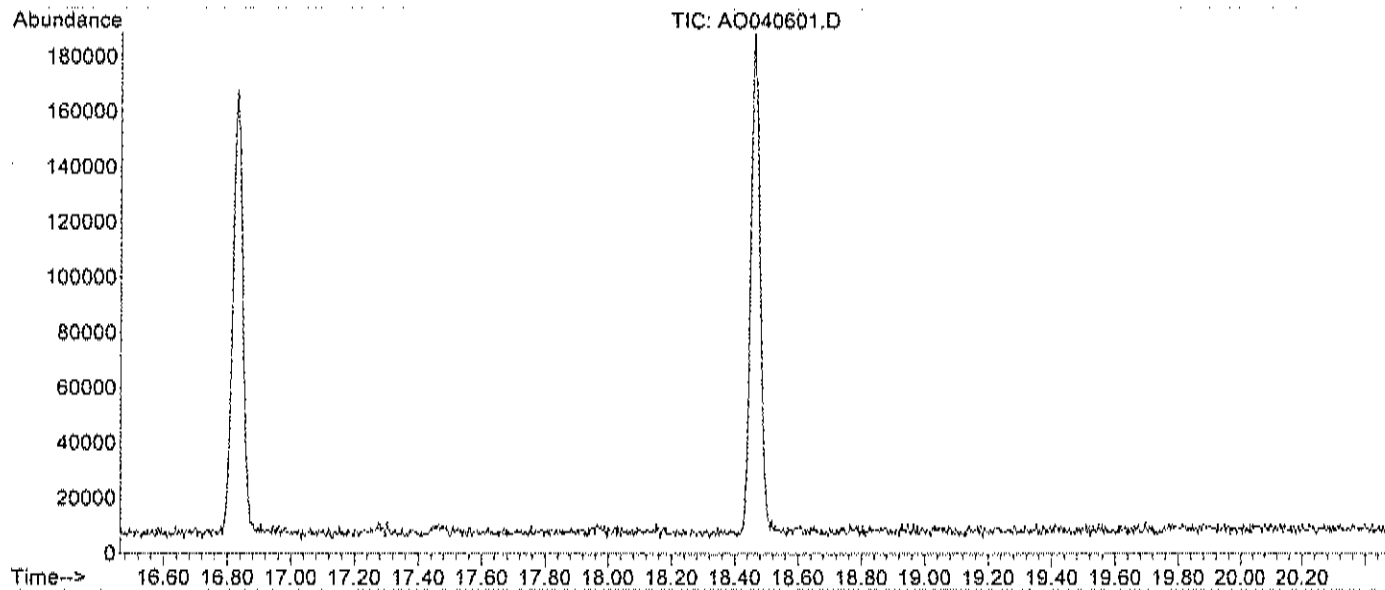
Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration



Spectrum Information: Average of 18.461 to 18.467 min.

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	8	40	20.2	6069	PASS
75	95	30	66	60.7	18221	PASS
95	95	100	100	100.0	30010	PASS
96	95	5	9	5.3	1590	PASS
173	174	0.00	2	0.0	0	PASS
174	95	50	120	91.1	27352	PASS
175	174	4	9	6.7	1838	PASS
176	174	95	101	97.5	26658	PASS
177	176	5	9	6.6	1750	PASS

Data File : C:\HPCHEM\1\DATA\AO040701.D

Vial: 1

Acq On : 7 Apr 2017 10:27 am

Operator: RJP

Sample : BFB1UG

Inst : MSD #1

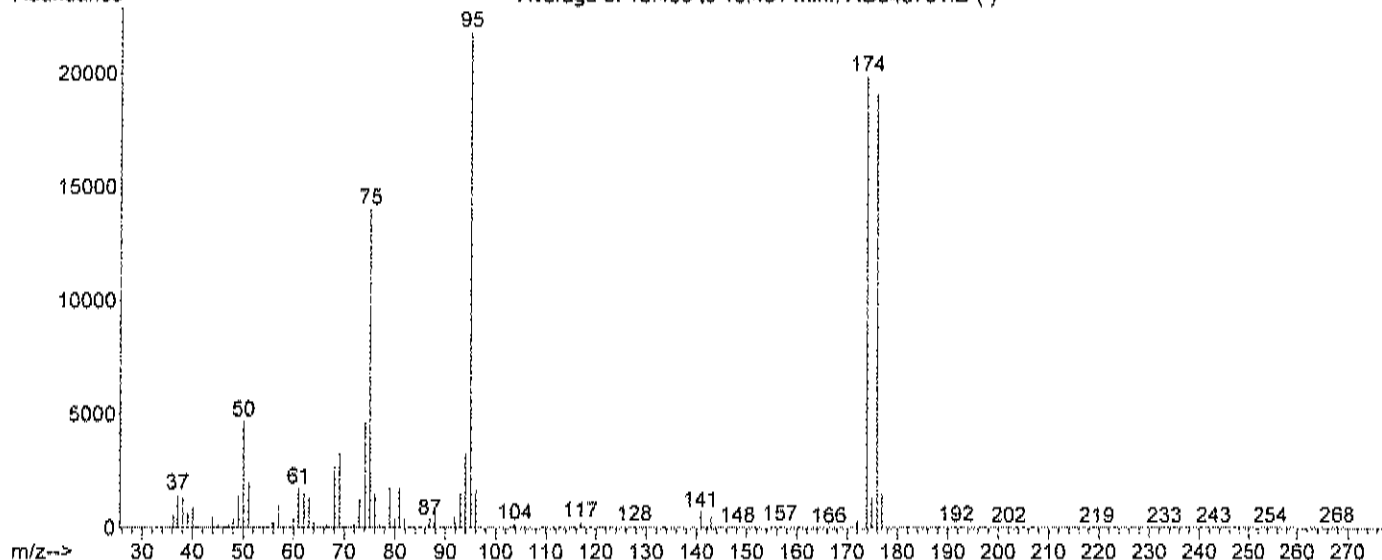
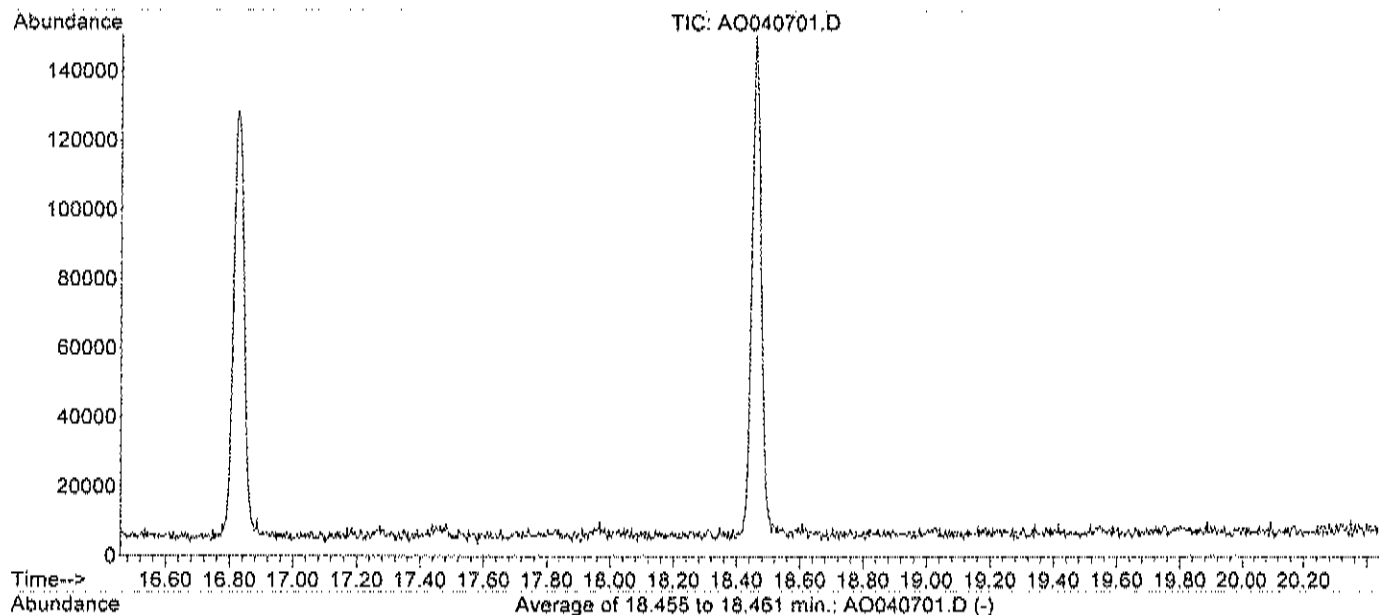
Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration



Spectrum Information: Average of 18.455 to 18.461 min.

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	8	40	21.5	4711	PASS
75	95	30	66	64.2	14056	PASS
95	95	100	100	100.0	21885	PASS
96	95	5	9	7.7	1679	PASS
173	174	0.00	2	0.0	0	PASS
174	95	50	120	91.1	19946	PASS
175	174	4	9	7.2	1427	PASS
176	174	95	101	96.0	19150	PASS
177	176	5	9	8.1	1550	PASS



**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**RAW QC DATA**



Date: 04-May-17

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.

Work Order: C1704014

Project: 691 St Paul Street

TestCode: 0.25CT-TCE-VC

Sample ID	AMB1UG-040617	Sample Type:	MBLK	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12114				
Client ID:	ZZZZZ	Batch ID:	R12114	TestNo: TO-15		Analysis Date: 4/6/2017	SeqNo: 141659				
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,1-Dichloroethene	< 0.15	0.15									
Chloroethane	< 0.15	0.15									
cis-1,2-Dichloroethene	< 0.15	0.15									
trans-1,2-Dichloroethene	< 0.15	0.15									
Trichloroethene	< 0.040	0.040									
Vinyl chloride	< 0.040	0.040									

Sample ID	AMB1UG-040717	Sample Type:	MBLK	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12118				
Client ID:	ZZZZZ	Batch ID:	R12118	TestNo: TO-15		Analysis Date: 4/7/2017	SeqNo: 141730				
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,1-Dichloroethene	< 0.15	0.15									
Chloroethane	< 0.15	0.15									
cis-1,2-Dichloroethene	< 0.15	0.15									
trans-1,2-Dichloroethene	< 0.15	0.15									
Trichloroethene	< 0.040	0.040									
Vinyl chloride	< 0.040	0.040									

Qualifiers:	J	Results reported are not blank corrected	E	Estimated Values above quantitation range	H	Holding times for preparations or analysis exceeded
	S	Analyte detected below quantitation limit	ND	Not Detected at the Limit of Detection	R	RPD outside accepted recovery limits
		Spike Recovery outside accepted recovery limits				

Data File : C:\HPCHEM\1\DATA\AO040604.D

Vial: 4

Acq On : 6 Apr 2017 10:47 am

Operator: RJP

Sample : AMB1UG-040617

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 07 07:18:44 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	26038	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	118374	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	102379	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	63192	0.90	ppb	0.01
Spiked Amount	1.000	Range	70 - 130	Recovery	=	90.00%

Target Compounds

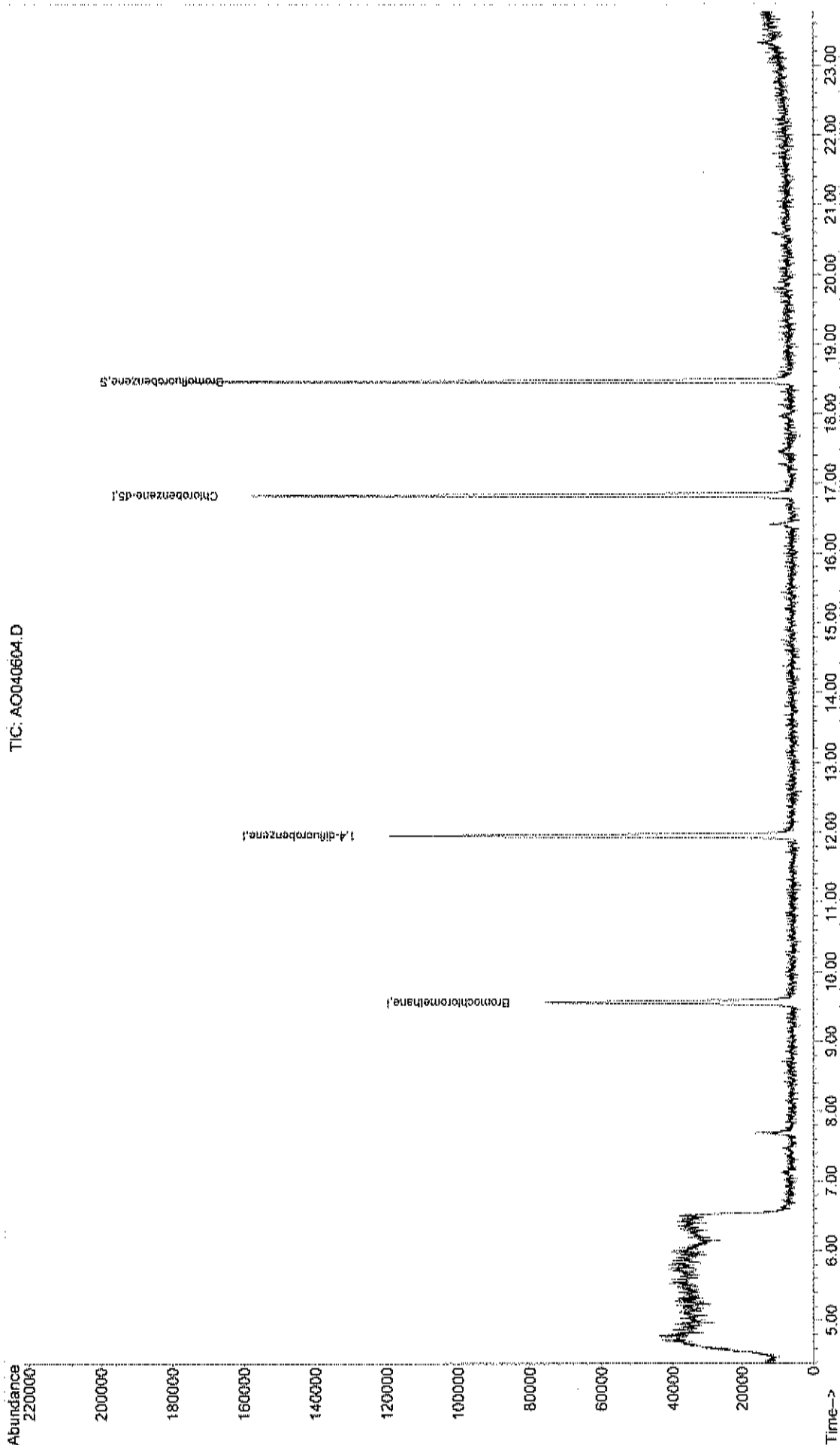
Qvalue

Data File : C:\HPCHEM\1\DATA\AO040604.D  
 Acq On : 6 Apr 2017 10:47 am  
 Sample : AMB1UG-040617  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 10 13:53 2017

Vial: 4  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\AO040704.D  
Acq On : 7 Apr 2017 12:36 pm  
Sample : AMB1UG-040717  
Misc : A331\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Apr 12 08:33:52 2017

Vial: 4  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Apr 03 10:15:59 2017  
Response via : Initial Calibration  
DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.55	128	21100	1.00	ppb	0.01
35) 1,4-difluorobenzene	11.95	114	92197	1.00	ppb	0.01
50) Chlorobenzene-d5	16.82	117	83384	1.00	ppb	0.00

System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	48987	0.86	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	86.00%

Target Compounds

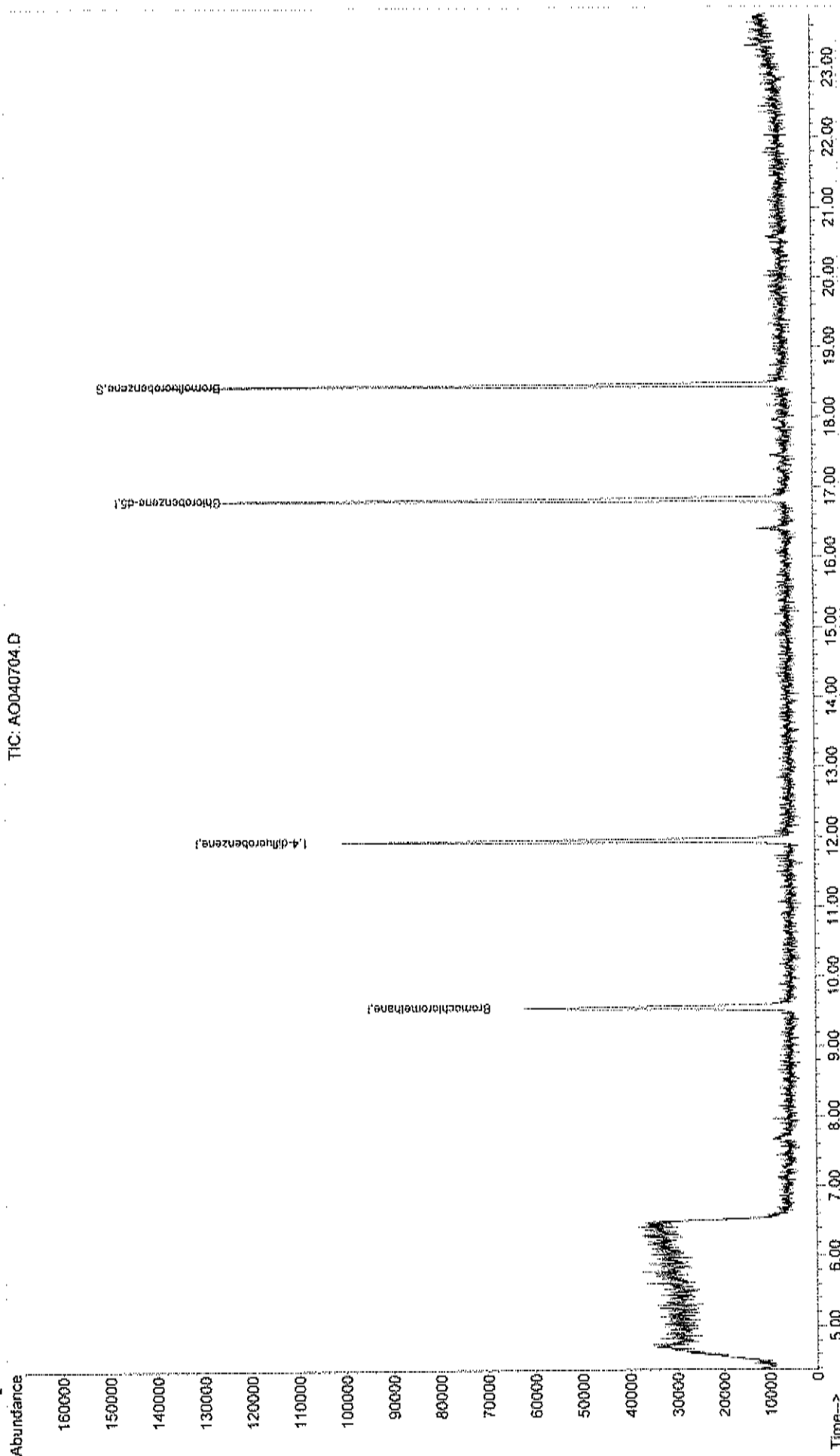
Qvalue

Data File : C:\HPCHEM\1\DATA\AO040704.D  
 Acq On : 7 Apr 2017 12:36 pm  
 Sample : AMB1UG-040717  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 12 8:33 2017

Vial: 4  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration



## CENTEK LABORATORIES, LLC

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.

Work Order: C1704014

Project: 691 St Paul Street

TestCode: 0.25CT-TCE-VC

Sample ID	C1704014-001A MS	SampType: MS	TestCode: 0.25CT-TCE-	Units: ppbv	Prep Date:	RunNo: 12114					
Client ID:	691-B19-1AQ-1	Batch ID: R12114	TestNo: TO-15		Analysis Date: 4/6/2017	SeqNo: 141726					
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	0.9100	0.15	1	0	91.0	70	130				
Chloroethane	1.040	0.15	1	0	104	70	130				
cis-1,2-Dichloroethene	1.230	0.15	1	0.33	90.0	70	130				
trans-1,2-Dichloroethene	0.9500	0.15	1	0	95.0	70	130				
Trichloroethene	1.390	0.040	1	0.4	99.0	70	130				
Vinyl chloride	1.030	0.040	1	0	103	70	130				

Sample ID	C1704014-001A MS	SampType: MSD	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12114					
Client ID:	691-B19-1AQ-1	Batch ID: R12114	TestNo: TO-15		Analysis Date: 4/6/2017	SeqNo: 141727					
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	1.000	0.15	1	0	100	70	130	0.91	9.42	30	
Chloroethane	1.040	0.15	1	0	104	70	130	1.04	0	30	
cis-1,2-Dichloroethene	1.320	0.15	1	0.33	99.0	70	130	1.23	7.06	30	
trans-1,2-Dichloroethene	0.9700	0.15	1	0	97.0	70	130	0.95	2.08	30	
Trichloroethene	1.370	0.040	1	0.4	97.0	70	130	1.39	1.45	30	
Vinyl chloride	1.030	0.040	1	0	103	70	130	1.03	0	30	

Qualifiers: J Results reported are not blank corrected  
 S Analyte detected below quantitation limit  
 ND Spike Recovery outside accepted recovery limits  
 E Estimated Value above quantitation range  
 ND Not Detected at the Limit of Detection  
 H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits



Data File : C:\HPCHEM\1\DATA\AO040622.D

Vial: 22

Acq On : 6 Apr 2017 10:40 pm

Operator: RJP

Sample : C1704014-001A MS

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 07 07:19:02 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	26903	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	116940	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	105999	1.00	ppb	0.01

## System Monitoring Compounds

65) Bromofluorobenzene	18.47	95	69782	0.97	ppb	0.02
Spiked Amount	1.000	Range	70 - 130	Recovery	=	97.00%

## Target Compounds

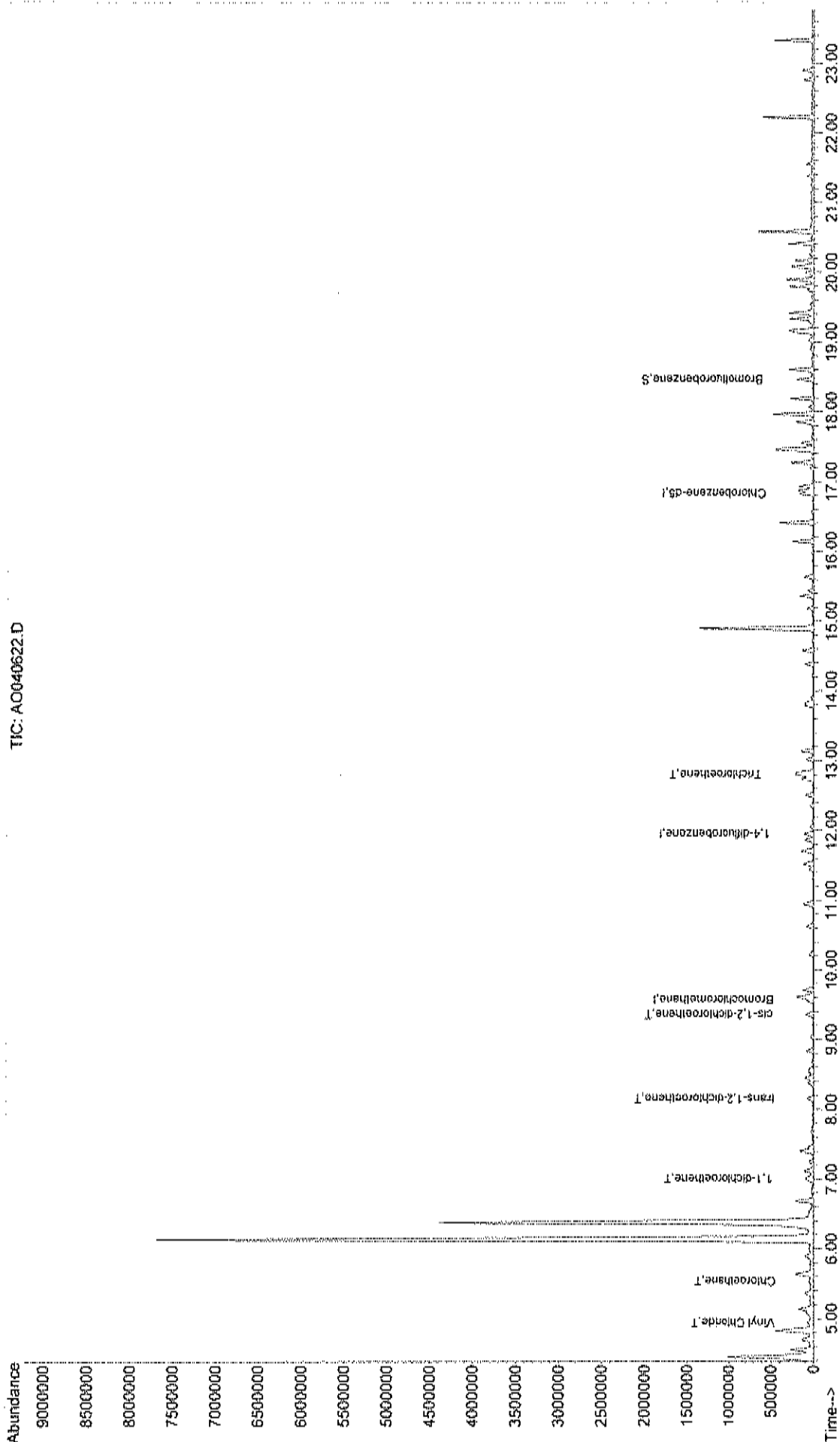
	R.T.	QIon	Response	Conc	Units	Qvalue
6) Vinyl Chloride	4.94	62	48685	1.03	ppb	88
10) Chloroethane	5.54	64	20446	1.04	ppb	# 73
18) 1,1-dichloroethene	6.99	96	31823	0.91	ppb	# 87
24) trans-1,2-dichloroethene	8.16	61	55866	0.95	ppb	93
29) cis-1,2-dichloroethene	9.36	61	68956	1.23	ppb	95
44) Trichloroethene	12.79	130	73147	1.39	ppb	91

Data File : C:\HPCHEM\1\DATA\AO040622.D  
Acq On : 6 Apr 2017 10:40 pm  
Sample : C1704014-001A MS  
Misc : A331\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Apr 11 14:05 2017

Vial: 22  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Thu May 04 11:27:28 2017  
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\AO040623.D

Vial: 23

Acq On : 6 Apr 2017 11:26 pm

Operator: RJP

Sample : C1704014-001A MSD

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 07 07:19:03 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.57	128	26299	1.00	ppb	0.03
35) 1,4-difluorobenzene	11.95	114	120898	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	106126	1.00	ppb	0.01

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	71496	0.99	ppb	0.01
Spiked Amount	1.000	Range	70 - 130	Recovery	=	99.00%

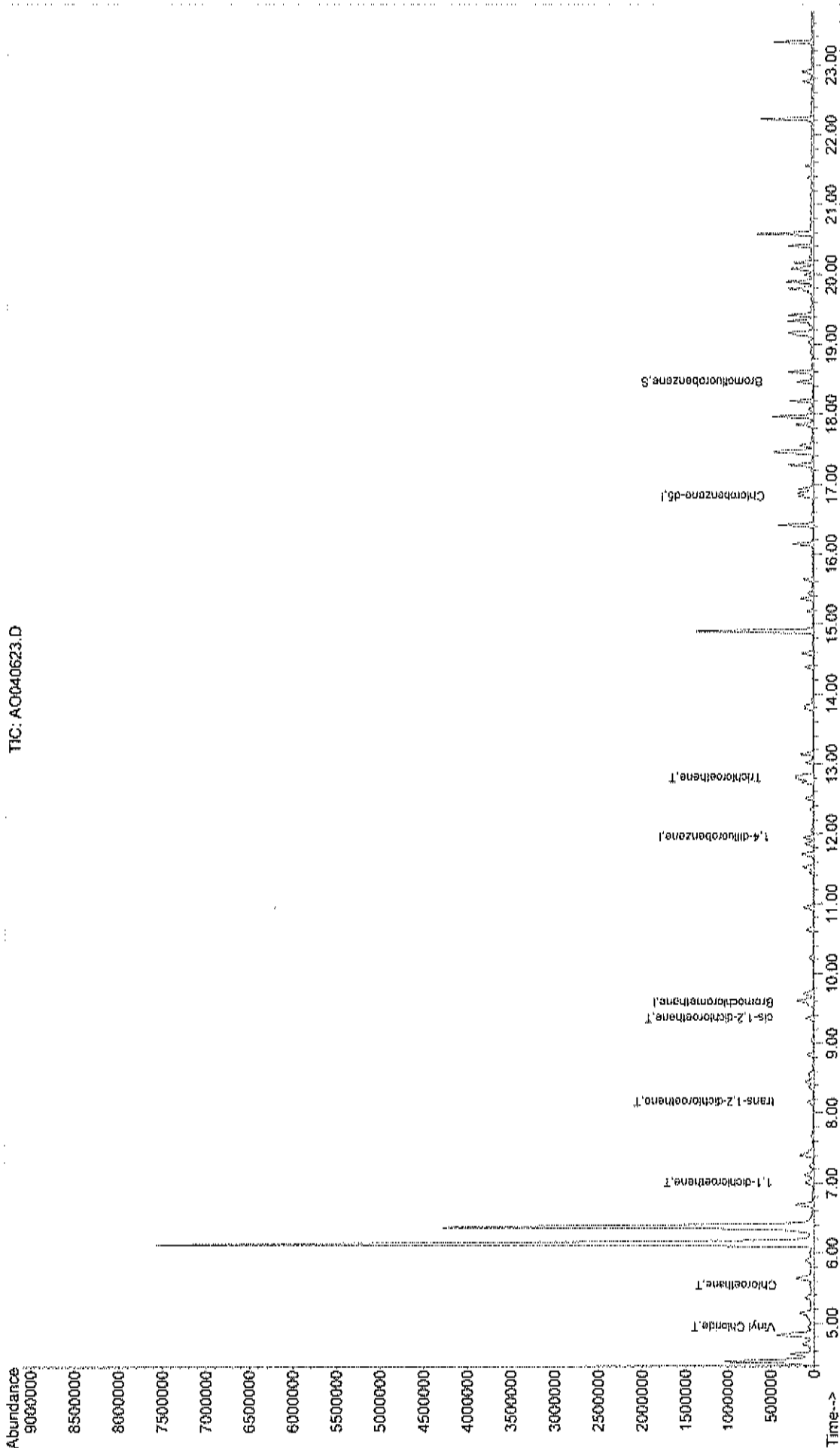
## Target Compounds

						Qvalue
6) Vinyl Chloride	4.94	62	47809	1.03	ppb	87
10) Chloroethane	5.54	64	20007	1.04	ppb	# 79
18) 1,1-dichloroethene	7.00	96	34130	1.00	ppb	94
24) trans-1,2-dichloroethene	8.16	61	55898	0.97	ppb	97
29) cis-1,2-dichloroethene	9.36	61	71906	1.32	ppb	95
44) Trichloroethene	12.79	130	74624	1.37	ppb	92

Data File : C:\HPCHEM\1\DATA\AO040623.D  
 Acq On : 6 Apr 2017 11:26 pm  
 Sample : C1704014-001A MSD  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 11 14:06 2017

Vial: 23  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00  
 Quant Results File: A331\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 11:27:28 2017  
 Response via : Initial Calibration



## CENTEK LABORATORIES, LLC

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.

Work Order: C1704014

Project: 691 St Paul Street

TestCode: 0.25CT-TCE-VC

Sample ID	ALCS1UG-040617	SampType: LCS	TestCode: 0.25CT-TCE-	Units: ppbv	Prep Date:	RunNo: 12114					
Client ID: ZZZZZ		Batch ID: R12114	TestNo: TO-15		Analysis Date: 4/6/2017	SeqNo: 141660					
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	0.9700	0.15	1	0	97.0	70	130				
Chloroethane	1.090	0.15	1	0	109	70	130				
cis-1,2-Dichloroethene	0.9700	0.15	1	0	97.0	70	130				
trans-1,2-Dichloroethene	0.9500	0.15	1	0	95.0	70	130				
Trichloroethene	0.9700	0.040	1	0	97.0	70	130				
Vinyl chloride	0.9700	0.040	1	0	97.0	70	130				

Sample ID	ALCS1UG-040717	SampType: LCS	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12118					
Client ID: ZZZZZ		Batch ID: R12118	TestNo: TO-15		Analysis Date: 4/7/2017	SeqNo: 141731					
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	0.9900	0.15	1	0	99.0	70	130				
Chloroethane	1.180	0.15	1	0	118	70	130				
cis-1,2-Dichloroethene	1.000	0.15	1	0	100	70	130				
trans-1,2-Dichloroethene	0.9800	0.15	1	0	98.0	70	130				
Trichloroethene	1.060	0.040	1	0	106	70	130				
Vinyl chloride	1.150	0.040	1	0	115	70	130				

Sample ID	ALCS1UGD-040617	SampType: LCSD	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12114					
Client ID: ZZZZZ	Batch ID: R12114	TestNo: TO-15	Analysis Date: 4/6/2017	SeqNo: 141661							
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	0.9900	0.15	1	0	99.0	70	130	0.97	2.04	30	
Chloroethane	1.000	0.15	1	0	100	70	130	1.09	8.61	30	

Qualifiers: . Results reported are not blank corrected  
 J Analyte detected below quantitation limit  
 S Spike Recovery outside accepted recovery limits  
 E Estimated Value above quantitation range  
 ND Not Detected at the Limit of Detection  
 H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits

CLIENT: LaBella Associates, P.C.  
 Work Order: C1704014  
 Project: 691 St Paul Street

TestCode: 0.25CT-TCE-VC

Sample ID	ALCS1UGD-040617	SampleType: LCSD	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12114					
Client ID: ZZZZZ		Batch ID: R12114	TestNo: TO-15		Analysis Date: 4/6/2017	SeqNo: 141661					
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
cis-1,2-Dichloroethene	0.9600	0.15	1	0	96.0	70	130	0.97	1.04	30	
trans-1,2-Dichloroethene	0.9800	0.15	1	0	98.0	70	130	0.95	3.11	30	
Trichloroethene	1.000	0.040	1	0	100	70	130	0.97	3.05	30	
Vinyl chloride	1.060	0.040	1	0	106	70	130	0.97	8.87	30	

Sample ID	ALCS1UGD-040717	SampleType: LCSD	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12118					
Client ID: ZZZZZ		Batch ID: R12118	TestNo: TO-15		Analysis Date: 4/8/2017	SeqNo: 141732					
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	1.080	0.15	1	0	108	70	130	0.99	8.70	30	
Chloroethane	1.180	0.15	1	0	118	70	130	1.18	0	30	
cis-1,2-Dichloroethene	1.120	0.15	1	0	112	70	130	1	11.3	30	
trans-1,2-Dichloroethene	1.070	0.15	1	0	107	70	130	0.98	8.78	30	
Trichloroethene	1.100	0.040	1	0	110	70	130	1.06	3.70	30	
Vinyl chloride	1.240	0.040	1	0	124	70	130	1.15	7.53	30	

Qualifiers: J Results reported are not blank corrected E Estimated Value above quantitation range H Holding times for preparation or analysis exceeded  
 S Analyte detected below quantitation limit ND Not Detected at the Limit of Detection R RPD outside accepted recovery limits  
 Spike Recovery outside accepted recovery limits

Data File : C:\HPCHEM\1\DATA\AO040603.D

Vial: 3

Acq On : 6 Apr 2017 10:10 am

Operator: RJP

Sample : ALCS1UG-040617

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 07 07:18:43 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	26450	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	123600	1.00	ppb	0.02
50) Chlorobenzene-d5	16.83	117	109377	1.00	ppb	0.01

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	71003	0.95	ppb	0.01
Spiked Amount	1.000	Range	70 - 130	Recovery	=	95.00%

## Target Compounds

						Qvalue
2) Propylene	4.49	41	40888	1.08	ppb	97
3) Freon 12	4.56	85	251398	1.15	ppb	97
4) Chloromethane	4.72	50	35411m	1.15	ppb	
5) Freon 114	4.83	85	166226	0.98	ppb	85
6) Vinyl Chloride	4.95	62	45262	0.97	ppb	93
7) Butane	5.15	43	46850	0.99	ppb	88
8) 1,3-butadiene	5.10	39	33934m	1.10	ppb	
9) Bromomethane	5.37	94	68097	1.08	ppb	90
10) Chloroethane	5.55	64	21194	1.09	ppb	93
11) Ethanol	5.66	45	13609	0.89	ppb	83
12) Acrolein	6.01	56	13709	0.86	ppb	95
13) Vinyl Bromide	5.90	106	68500	1.09	ppb	94
14) Freon 11	6.33	101	216910	1.00	ppb	98
15) Acetone	6.15	58	18471	1.08	ppb	# 71
16) Pentane	6.70	42	32579	0.98	ppb	# 39
17) Isopropyl alcohol	6.39	45	59472	1.05	ppb	# 100
18) 1,1-dichloroethene	6.99	96	33335	0.97	ppb	93
19) Freon 113	7.41	101	81489	0.99	ppb	91
20) t-Butyl alcohol	7.03	59	81168	0.97	ppb	94
21) Methylene chloride	7.12	84	32341	0.96	ppb	# 80
22) Allyl chloride	7.24	41	39902	0.93	ppb	90
23) Carbon disulfide	7.47	76	106791	0.99	ppb	81
24) trans-1,2-dichloroethene	8.16	61	54775	0.95	ppb	95
25) methyl tert-butyl ether	8.46	73	103728	0.95	ppb	90
26) 1,1-dichloroethane	8.38	63	69940	0.98	ppb	93
27) Vinyl acetate	8.54	43	107835	1.17	ppb	94
28) Methyl Ethyl Ketone	8.84	72	16993	0.97	ppb	99
29) cis-1,2-dichloroethene	9.36	61	53336	0.97	ppb	93
30) Hexane	9.62	57	47571	0.94	ppb	88
31) Ethyl acetate	9.62	43	108734	0.97	ppb	96
32) Chloroform	9.71	83	98677	0.98	ppb	98
33) Tetrahydrofuran	10.23	42	34525	0.92	ppb	86
34) 1,2-dichloroethane	10.62	62	71589	0.96	ppb	98
36) 1,1,1-trichloroethane	10.95	97	108201	0.99	ppb	95
37) Cyclohexane	11.86	56	52233	0.95	ppb	98
38) Carbon tetrachloride	11.71	117	112797	0.98	ppb	99
39) Benzene	11.52	78	111117	0.98	ppb	92
40) Methyl methacrylate	13.02	41	48236	0.91	ppb	94
41) 1,4-dioxane	12.82	58	22028	0.94	ppb	# 57
42) 2,2,4-trimethylpentane	12.84	57	167978	0.98	ppb	90
43) Heptane	13.14	43	56488	0.95	ppb	95
44) Trichloroethene	12.79	130	53930	0.97	ppb	89
45) 1,2-dichloropropane	12.51	63	42172	1.01	ppb	100

(#)= qualifier out of range (m) = manual integration



Data File : C:\HPCHEM\1\DATA\AO040603.D

Vial: 3

Acq On : 6 Apr 2017 10:10 am

Operator: RJP

Sample : ALCS1UG-040617

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 07 07:18:43 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

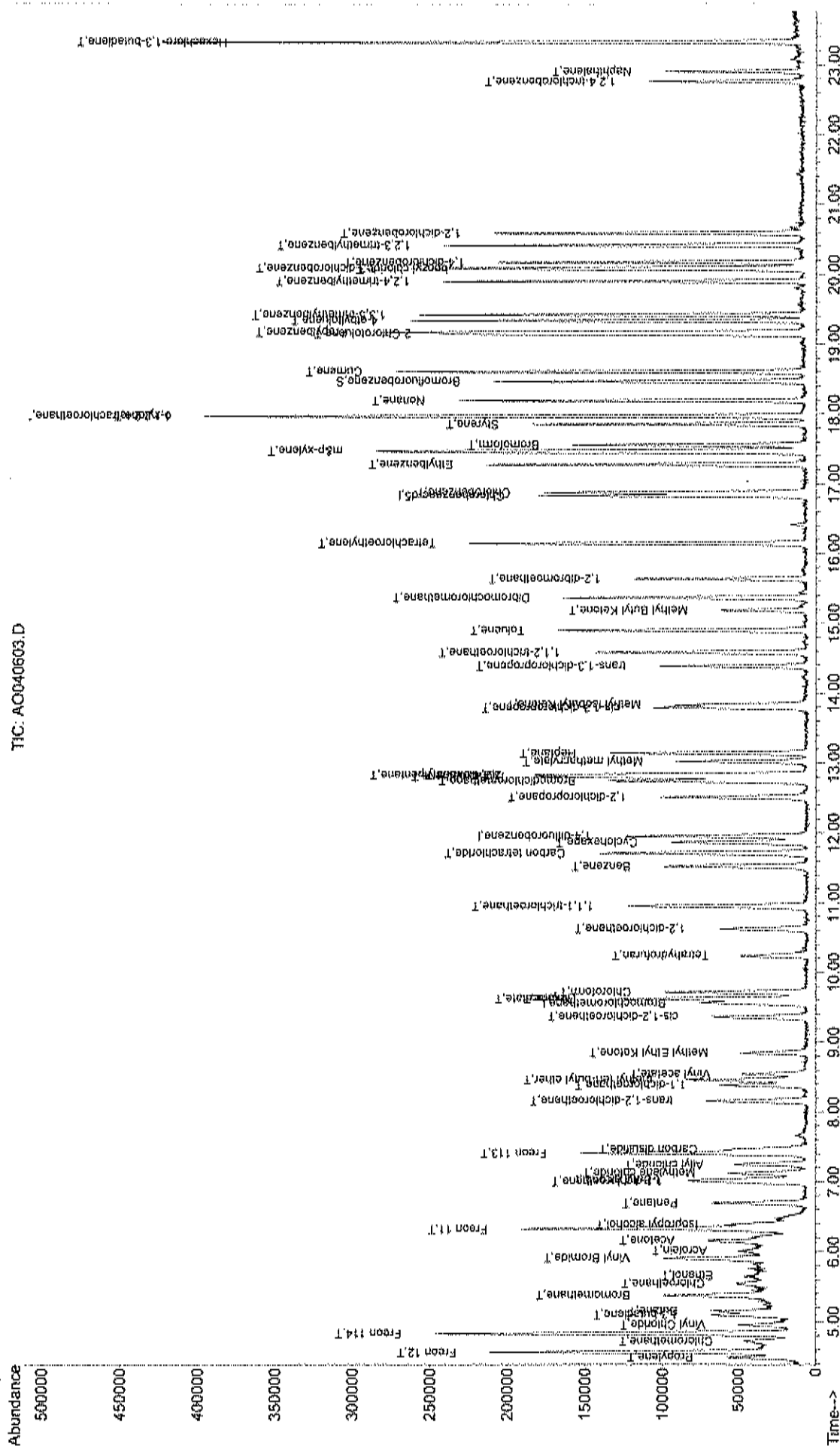
Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.74	83	108252	1.00	ppb	98
47) cis-1,3-dichloropropene	13.79	75	69338	1.00	ppb	97
48) trans-1,3-dichloropropene	14.38	75	64133	0.98	ppb	95
49) 1,1,2-trichloroethane	14.58	97	51263	1.02	ppb	100
51) Toluene	14.90	92	78308	0.96	ppb #	82
52) Methyl Isobutyl Ketone	13.83	43	68416	0.80	ppb	95
53) Dibromochloromethane	15.36	129	100792	1.01	ppb	93
54) Methyl Butyl Ketone	15.19	43	53361m	0.70	ppb	
55) 1,2-dibromoethane	15.64	107	84703	1.02	ppb	98
56) Tetrachloroethylene	16.15	164	55443	0.96	ppb	96
57) Chlorobenzene	16.88	112	113693	1.00	ppb	98
58) Ethylbenzene	17.27	91	183764	0.96	ppb	100
59) m&p-xylene	17.47	91	314573	1.95	ppb	99
60) Nonane	18.19	43	85662	0.96	ppb	94
61) Styrene	17.85	104	95512	0.95	ppb	99
62) Bromoform	17.56	173	82565	0.99	ppb	99
63) o-xylene	17.97	91	148340	0.95	ppb	97
64) Cumene	18.60	105	199558	0.99	ppb	99
66) 1,1,2,2-tetrachloroethane	17.96	83	112277	1.06	ppb	96
67) Propylbenzene	19.17	120	53242	1.02	ppb #	1
68) 2-Chlorotoluene	19.14	126	47557	0.99	ppb #	1
69) 4-ethyltoluene	19.33	105	186873	0.99	ppb	100
70) 1,3,5-trimethylbenzene	19.42	105	161548	0.96	ppb	99
71) 1,2,4-trimethylbenzene	19.90	105	154106	0.98	ppb	95
72) 1,3-dichlorobenzene	20.09	146	106249	1.08	ppb	99
73) benzyl chloride	20.07	91	97783	1.18	ppb	95
74) 1,4-dichlorobenzene	20.17	146	95059	1.04	ppb	98
75) 1,2,3-trimethylbenzene	20.41	105	151247	0.96	ppb	95
76) 1,2-dichlorobenzene	20.59	146	98338	1.04	ppb	97
77) 1,2,4-trichlorobenzene	22.76	180	31814	1.13	ppb	92
78) Naphthalene	22.91	128	73959	1.02	ppb	94
79) Hexachloro-1,3-butadiene	23.32	225	74843	1.06	ppb	95

TIC: A0040603.D



Data File : C:\HPCHEM\1\DATA\AO040703.D

Vial: 3

Acq On : 7 Apr 2017 12:00 pm

Operator: RJP

Sample : ALCS1UG-040717

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 12 08:33:43 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.55	128	22420	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.95	114	98698	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	86489	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.46	95	57254	0.97	ppb	0.01
Spiked Amount	1.000	Range	70 - 130	Recovery	=	97.00%

## Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Propylene	4.48	41	33409	1.05	ppb	76
3) Freon 12	4.56	85	219946	1.19	ppb	99
4) Chloromethane	4.73	50	30609m	1.17	ppb	
5) Freon 114	4.82	85	179639m	1.25	ppb	
6) Vinyl Chloride	4.95	62	45445	1.15	ppb	94
7) Butane	5.15	43	44744	1.11	ppb	# 89
8) 1,3-butadiene	5.09	39	30908m	1.18	ppb	
9) Bromomethane	5.37	94	62521	1.17	ppb	89
10) Chloroethane	5.54	64	19371	1.18	ppb	# 83
11) Ethanol	5.65	45	14756	1.14	ppb	88
12) Acrolein	6.01	56	12973	0.96	ppb	92
13) Vinyl Bromide	5.89	106	63758	1.19	ppb	96
14) Freon 11	6.32	101	238213m	1.29	ppb	
15) Acetone	6.15	58	16292	1.13	ppb	# 64
16) Pentane	6.69	42	27585	0.98	ppb	# 38
17) Isopropyl alcohol	6.40	45	55115	1.15	ppb	# 100
18) 1,1-dichloroethene	6.99	96	28995	0.99	ppb	94
19) Freon 113	7.40	101	71887	1.03	ppb	90
20) t-Butyl alcohol	7.04	59	64341	0.91	ppb	92
21) Methylene chloride	7.12	84	28008	0.99	ppb	# 75
22) Allyl chloride	7.24	41	33731	0.93	ppb	91
23) Carbon disulfide	7.46	76	91473	1.01	ppb	84
24) trans-1,2-dichloroethene	8.15	61	47949	0.98	ppb	93
25) methyl tert-butyl ether	8.45	73	88153	0.95	ppb	87
26) 1,1-dichloroethane	8.37	63	62741	1.03	ppb	96
27) Vinyl acetate	8.54	43	92002	1.17	ppb	92
28) Methyl Ethyl Ketone	8.85	72	15128	1.01	ppb	98
29) cis-1,2-dichloroethene	9.36	61	46370	1.00	ppb	92
30) Hexane	9.60	57	40518	0.95	ppb	87
31) Ethyl acetate	9.62	43	94395	0.99	ppb	97
32) Chloroform	9.71	83	87650	1.03	ppb	96
33) Tetrahydrofuran	10.23	42	30292	0.95	ppb	84
34) 1,2-dichloroethane	10.62	62	59420	0.94	ppb	94
36) 1,1,1-trichloroethane	10.94	97	93614	1.08	ppb	98
37) Cyclohexane	11.86	56	45664	1.04	ppb	95
38) Carbon tetrachloride	11.70	117	97837	1.06	ppb	100
39) Benzene	11.52	78	98582	1.09	ppb	92
40) Methyl methacrylate	13.02	41	43154	1.01	ppb	95
41) 1,4-dioxane	12.81	58	17665	0.94	ppb	# 55
42) 2,2,4-trimethylpentane	12.83	57	144457	1.06	ppb	88
43) Heptane	13.14	43	48601	1.03	ppb	98
44) Trichloroethene	12.79	130	47004	1.06	ppb	91
45) 1,2-dichloropropane	12.50	63	36145	1.08	ppb	100

(#)=qualifier out of range (m)=manual integration

Data File : C:\HPCHEM\1\DATA\AO040703.D

Vial: 3

Acq On : 7 Apr 2017 12:00 pm

Operator: RJP

Sample : ALCS1UG-040717

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 12 08:33:43 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

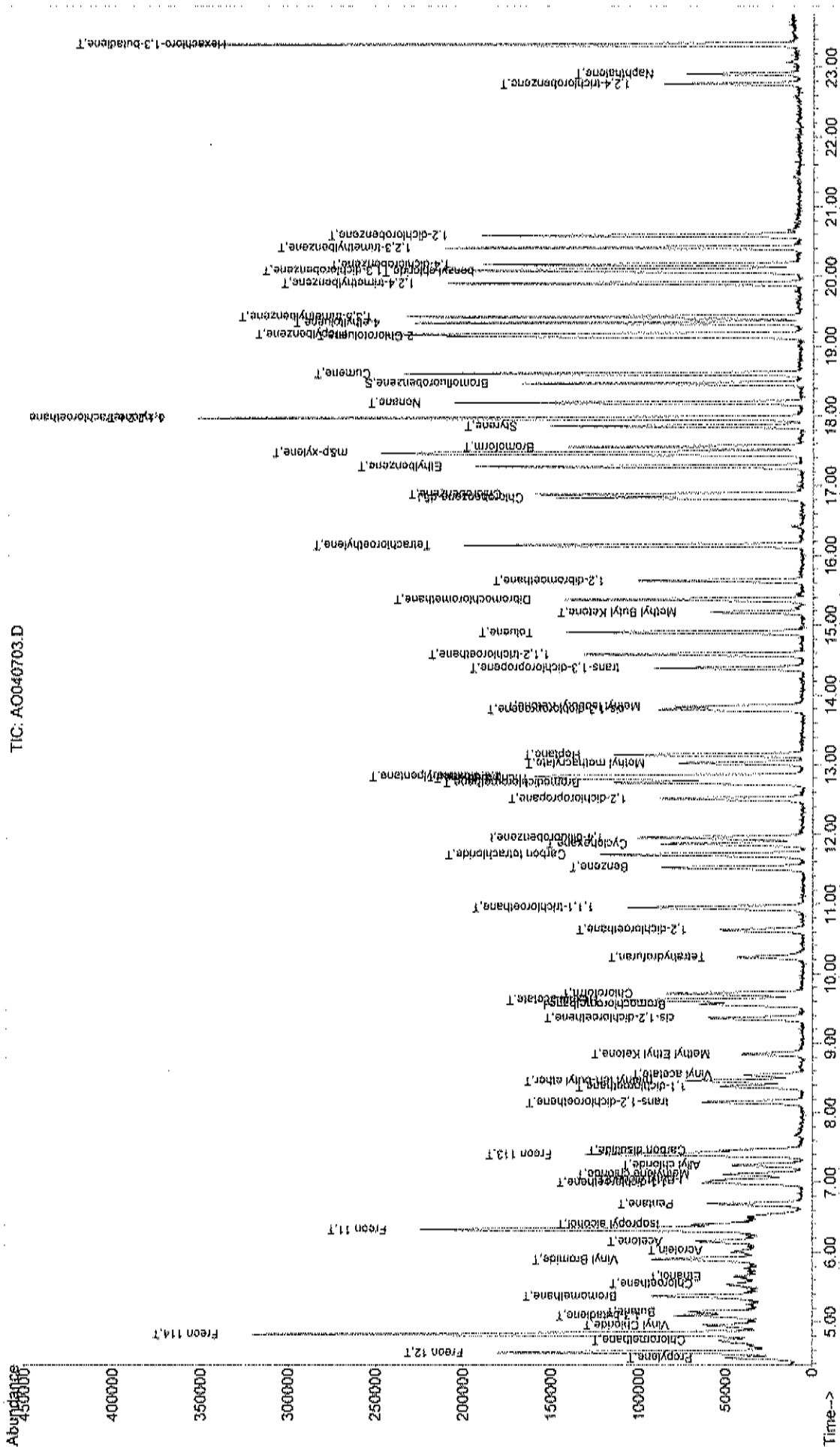
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.74	83	91783	1.06	ppb	96
47) cis-1,3-dichloropropene	13.79	75	58172	1.06	ppb	95
48) trans-1,3-dichloropropene	14.38	75	54481	1.04	ppb	90
49) 1,1,2-trichloroethane	14.58	97	46598	1.16	ppb	94
51) Toluene	14.89	92	63954	0.99	ppb	# 78
52) Methyl Isobutyl Ketone	13.83	43	60946	0.90	ppb	95
53) Dibromochloromethane	15.36	129	84657	1.07	ppb	90
54) Methyl Butyl Ketone	15.18	43	46099	0.76	ppb	93
55) 1,2-dibromoethane	15.63	107	72687	1.10	ppb	96
56) Tetrachloroethylene	16.14	164	49756	1.09	ppb	99
57) Chlorobenzene	16.87	112	96040	1.06	ppb	95
58) Ethylbenzene	17.27	91	157903	1.04	ppb	100
59) m&p-xylene	17.46	91	265326	2.08	ppb	100
60) Nonane	18.18	43	76447	1.08	ppb	93
61) Styrene	17.85	104	81583	1.03	ppb	95
62) Bromoform	17.55	173	71209	1.08	ppb	99
63) o-xylene	17.96	91	127835	1.04	ppb	96
64) Cumene	18.60	105	168618	1.05	ppb	100
66) 1,1,2,2-tetrachloroethane	17.96	83	98561	1.17	ppb	96
67) Propylbenzene	19.17	120	44751	1.09	ppb	# 1
68) 2-Chlorotoluene	19.14	126	40730	1.07	ppb	# 1
69) 4-ethyltoluene	19.33	105	163130	1.09	ppb	97
70) 1,3,5-trimethylbenzene	19.42	105	138712	1.04	ppb	96
71) 1,2,4-trimethylbenzene	19.90	105	126848	1.02	ppb	93
72) 1,3-dichlorobenzene	20.09	146	91327	1.17	ppb	99
73) benzyl chloride	20.06	91	82451	1.26	ppb	94
74) 1,4-dichlorobenzene	20.17	146	83163	1.15	ppb	97
75) 1,2,3-trimethylbenzene	20.41	105	132486	1.06	ppb	95
76) 1,2-dichlorobenzene	20.59	146	88758	1.18	ppb	96
77) 1,2,4-trichlorobenzene	22.76	180	25921	1.17	ppb	95
78) Naphthalene	22.91	128	56238	0.98	ppb	94
79) Hexachloro-1,3-butadiene	23.32	225	68219	1.22	ppb	96

Data File : C:\HPCHEM\1\DATA\AO040703.D  
Acq On : 7 Apr 2017 12:00 pm  
Sample : ALCS1UG-040717  
Misc : A331\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Apr 12 8:34 2017

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Thu May 04 11:27:28 2017  
Response via : Initial Calibration

Vial: 3  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A331\_1UG.RES



Data File : C:\HPCHEM\1\DATA\AO040620.D

Vial: 49

Acq On : 6 Apr 2017 9:15 pm

Operator: RJP

Sample : ALCS1UGD-040617

Inst : MSD #1

Misc : A331\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Apr 07 07:19:00 2017

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Apr 03 10:15:59 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	24809	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	117535	1.00	ppb	0.02
50) Chlorobenzene-d5	16.83	117	100863	1.00	ppb	0.01

## System Monitoring Compounds

65) Bromofluorobenzene	18.47	95	67679	0.98	ppb	0.02
Spiked Amount	1.000	Range	70 - 130	Recovery	=	98.00%

## Target Compounds

						Qvalue
2) Propylene	4.50	41	35590	1.01	ppb	76
3) Freon 12	4.56	85	236111	1.16	ppb	97
4) Chloromethane	4.73	50	33866m	1.17	ppb	
5) Freon 114	4.83	85	190689	1.20	ppb	98
6) Vinyl Chloride	4.95	62	46545	1.06	ppb	92
7) Butane	5.15	43	47237	1.06	ppb	90
8) 1,3-butadiene	5.09	39	31406m	1.09	ppb	
9) Bromomethane	5.37	94	66887	1.14	ppb	88
10) Chloroethane	5.54	64	18161	1.00	ppb	# 73
11) Ethanol	5.65	45	14827	1.04	ppb	84
12) Acrolein	6.01	56	13354	0.89	ppb	98
13) Vinyl Bromide	5.91	106	67026	1.13	ppb	95
14) Freon 11	6.34	101	250288m	1.22	ppb	
15) Acetone	6.16	58	18291	1.14	ppb	# 69
16) Pentane	6.69	42	30218	0.97	ppb	# 44
17) Isopropyl alcohol	6.39	45	57181	1.08	ppb	# 100
18) 1,1-dichloroethene	7.00	96	32066	0.99	ppb	93
19) Freon 113	7.41	101	77004	1.00	ppb	93
20) t-Butyl alcohol	7.04	59	68096	0.87	ppb	96
21) Methylene chloride	7.13	84	31329	1.00	ppb	# 80
22) Allyl chloride	7.25	41	37083	0.93	ppb	90
23) Carbon disulfide	7.47	76	100769	1.00	ppb	83
24) trans-1,2-dichloroethene	8.15	61	53306	0.98	ppb	93
25) methyl tert-butyl ether	8.47	73	94567	0.92	ppb	91
26) 1,1-dichloroethane	8.39	63	68087	1.01	ppb	98
27) Vinyl acetate	8.55	43	100265	1.16	ppb	93
28) Methyl Ethyl Ketone	8.85	72	16009	0.97	ppb	96
29) cis-1,2-dichloroethene	9.36	61	49482	0.96	ppb	95
30) Hexane	9.62	57	43032	0.91	ppb	87
31) Ethyl acetate	9.62	43	99523	0.94	ppb	97
32) Chloroform	9.71	83	94443	1.00	ppb	99
33) Tetrahydrofuran	10.24	42	34468	0.98	ppb	95
34) 1,2-dichloroethane	10.62	62	68052	0.98	ppb	95
36) 1,1,1-trichloroethane	10.95	97	103974	1.00	ppb	96
37) Cyclohexane	11.86	56	49591	0.95	ppb	98
38) Carbon tetrachloride	11.71	117	105869	0.97	ppb	99
39) Benzene	11.52	78	105761	0.98	ppb	91
40) Methyl methacrylate	13.02	41	45439	0.90	ppb	97
41) 1,4-dioxane	12.83	58	18999	0.85	ppb	# 55
42) 2,2,4-trimethylpentane	12.84	57	163528	1.01	ppb	89
43) Heptane	13.14	43	52158	0.92	ppb	95
44) Trichloroethene	12.80	130	52813	1.00	ppb	94
45) 1,2-dichloropropane	12.51	63	40874	1.03	ppb	99

(#)=qualifier out of range (m)=manual integration

AO040620.D A331\_1UG.M

Thu May 04 11:41:32 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO040620.D  
 Acq On : 6 Apr 2017 9:15 pm  
 Sample : ALCS1UGD-040617  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 07 07:19:00 2017

Vial: 49  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Mon Apr 03 10:15:59 2017  
 Response via : Initial Calibration  
 DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.74	83	102280	1.00	ppb	98
47) cis-1,3-dichloropropene	13.79	75	64051	0.98	ppb	98
48) trans-1,3-dichloropropene	14.38	75	60406	0.97	ppb	94
49) 1,1,2-trichloroethane	14.58	97	48605	1.02	ppb	100
51) Toluene	14.90	92	74483	0.99	ppb	85
52) Methyl Isobutyl Ketone	13.84	43	56491m	0.72	ppb	
53) Dibromochloromethane	15.37	129	94470	1.02	ppb	92
54) Methyl Butyl Ketone	15.19	43	39316m	0.56	ppb	
55) 1,2-dibromoethane	15.64	107	78477	1.02	ppb	95
56) Tetrachloroethylene	16.15	164	54001	1.01	ppb	98
57) Chlorobenzene	16.87	112	107457	1.02	ppb	94
58) Ethylbenzene	17.27	91	170983	0.97	ppb	100
59) m&p-xylene	17.47	91	296006	1.99	ppb	99
60) Nonane	18.19	43	81677	0.99	ppb	91
61) Styrene	17.86	104	89618	0.97	ppb	96
62) Bromoform	17.56	173	75354	0.98	ppb	97
63) o-xylene	17.97	91	140353	0.98	ppb	94
64) Cumene	18.60	105	181432	0.97	ppb	99
66) 1,1,2,2-tetrachloroethane	17.96	83	101557	1.04	ppb	94
67) Propylbenzene	19.17	120	46499	0.97	ppb	# 1
68) 2-Chlorotoluene	19.14	126	43318	0.98	ppb	# 1
69) 4-ethyltoluene	19.34	105	172594	0.99	ppb	100
70) 1,3,5-trimethylbenzene	19.42	105	150070	0.97	ppb	100
71) 1,2,4-trimethylbenzene	19.90	105	139431	0.96	ppb	94
72) 1,3-dichlorobenzene	20.09	146	96756	1.06	ppb	96
73) benzyl chloride	20.07	91	78332	1.03	ppb	95
74) 1,4-dichlorobenzene	20.17	146	86547	1.03	ppb	96
75) 1,2,3-trimethylbenzene	20.41	105	138686	0.95	ppb	94
76) 1,2-dichlorobenzene	20.59	146	91284	1.04	ppb	98
77) 1,2,4-trichlorobenzene	22.77	180	26664	1.03	ppb	98
78) Naphthalene	22.91	128	58584	0.88	ppb	94
79) Hexachloro-1,3-butadiene	23.33	225	68422	1.05	ppb	94

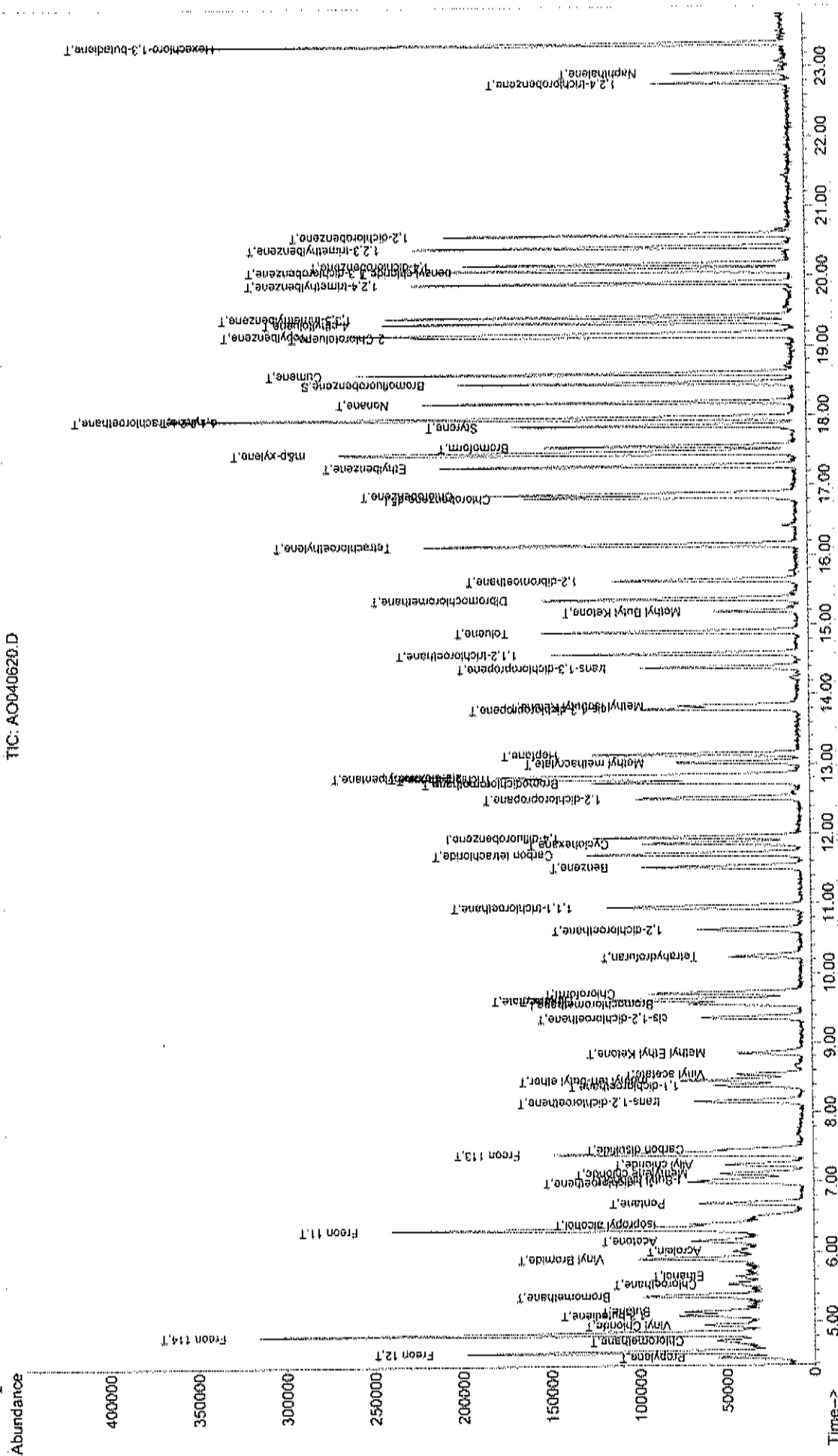


Data File : C:\HPCHEM\1\DATA\A0040620.D  
Acq On : 6 Apr 2017 9:15 pm  
Sample : ALCS1UGD-040617  
Misc : A331\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Apr 10 13:55 2017

Vial: 49  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Thu May 04 11:27:28 2017  
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\AO040725.D  
 Acq On : 8 Apr 2017 2:53 am  
 Sample : ALCS1UGD-040717  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 08 09:00:00 2017

Vial: 25  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Mon Apr 03 10:15:59 2017  
 Response via : Initial Calibration  
 DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.56	128	19313	1.00	ppb	0.02
35) 1,4-difluorobenzene	11.95	114	89890	1.00	ppb	0.01
50) Chlorobenzene-d5	16.83	117	78098	1.00	ppb	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
65) Bromofluorobenzene	18.46	95	51653	0.97	ppb	0.01
Spiked Amount	1.000	Range	70 - 130	Recovery	=	97.00%

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Propylene	4.48	41	31618	1.15	ppb	72
3) Freon 12	4.56	85	204448m	1.29	ppb	
4) Chloromethane	4.72	50	27304m	1.22	ppb	
5) Freon 114	4.82	85	166285m	1.34	ppb	
6) Vinyl Chloride	4.95	62	42200	1.24	ppb	88
7) Butane	5.14	43	45335	1.31	ppb	93
8) 1,3-butadiene	5.08	39	27014m	1.20	ppb	
9) Bromomethane	5.37	94	57912m	1.26	ppb	
10) Chloroethane	5.54	64	16720m	1.18	ppb	
11) Ethanol	5.65	45	14338	1.29	ppb	84
12) Acrolein	6.02	56	12507	1.07	ppb	96
13) Vinyl Bromide	5.90	106	58047m	1.26	ppb	
14) Freon 11	6.32	101	252102	1.58	ppb	99
15) Acetone	6.15	58	12985m	1.04	ppb	
16) Pentane	6.69	42	26726	1.10	ppb	# 43
17) Isopropyl alcohol	6.39	45	51366	1.24	ppb	# 100
18) 1,1-dichloroethene	7.00	96	27027	1.08	ppb	91
19) Freon 113	7.40	101	68890	1.15	ppb	89
20) t-Butyl alcohol	7.03	59	53096	0.87	ppb	# 92
21) Methylene chloride	7.12	84	26922	1.10	ppb	# 78
22) Allyl chloride	7.25	41	32876	1.05	ppb	91
23) Carbon disulfide	7.46	76	86481	1.10	ppb	84
24) trans-1,2-dichloroethene	8.15	61	45121	1.07	ppb	96
25) methyl tert-butyl ether	8.46	73	81727	1.02	ppb	96
26) 1,1-dichloroethane	8.38	63	59764	1.14	ppb	94
27) Vinyl acetate	8.54	43	84681	1.25	ppb	92
28) Methyl Ethyl Ketone	8.84	72	12658	0.99	ppb	99
29) cis-1,2-dichloroethene	9.36	61	45009	1.12	ppb	94
30) Hexane	9.61	57	38800	1.05	ppb	89
31) Ethyl acetate	9.62	43	86633	1.05	ppb	95
32) Chloroform	9.71	83	83163	1.14	ppb	96
33) Tetrahydrofuran	10.23	42	27309	1.00	ppb	87
34) 1,2-dichloroethane	10.62	62	59562	1.10	ppb	96
36) 1,1,1-trichloroethane	10.95	97	86730	1.10	ppb	99
37) Cyclohexane	11.86	56	42448	1.06	ppb	99
38) Carbon tetrachloride	11.71	117	92728	1.11	ppb	100
39) Benzene	11.52	78	90008	1.09	ppb	92
40) Methyl methacrylate	13.02	41	35136	0.91	ppb	90
41) 1,4-dioxane	12.83	58	14989m	0.88	ppb	
42) 2,2,4-trimethylpentane	12.83	57	134803	1.08	ppb	88
43) Heptane	13.14	43	45450	1.05	ppb	95
44) Trichloroethene	12.79	130	44637	1.10	ppb	92
45) 1,2-dichloropropane	12.51	63	34787	1.14	ppb	100

(#) = qualifier out of range (m) = manual integration  
 AO040725.D A331\_1UG.M Thu May 04 11:41:49 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO040725.D  
 Acq On : 8 Apr 2017 2:53 am  
 Sample : ALCS1UGD-040717  
 Misc : A331\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Apr 08 09:00:00 2017

Vial: 25  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A331\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Mon Apr 03 10:15:59 2017  
 Response via : Initial Calibration  
 DataAcq Meth : 1UG\_RUN

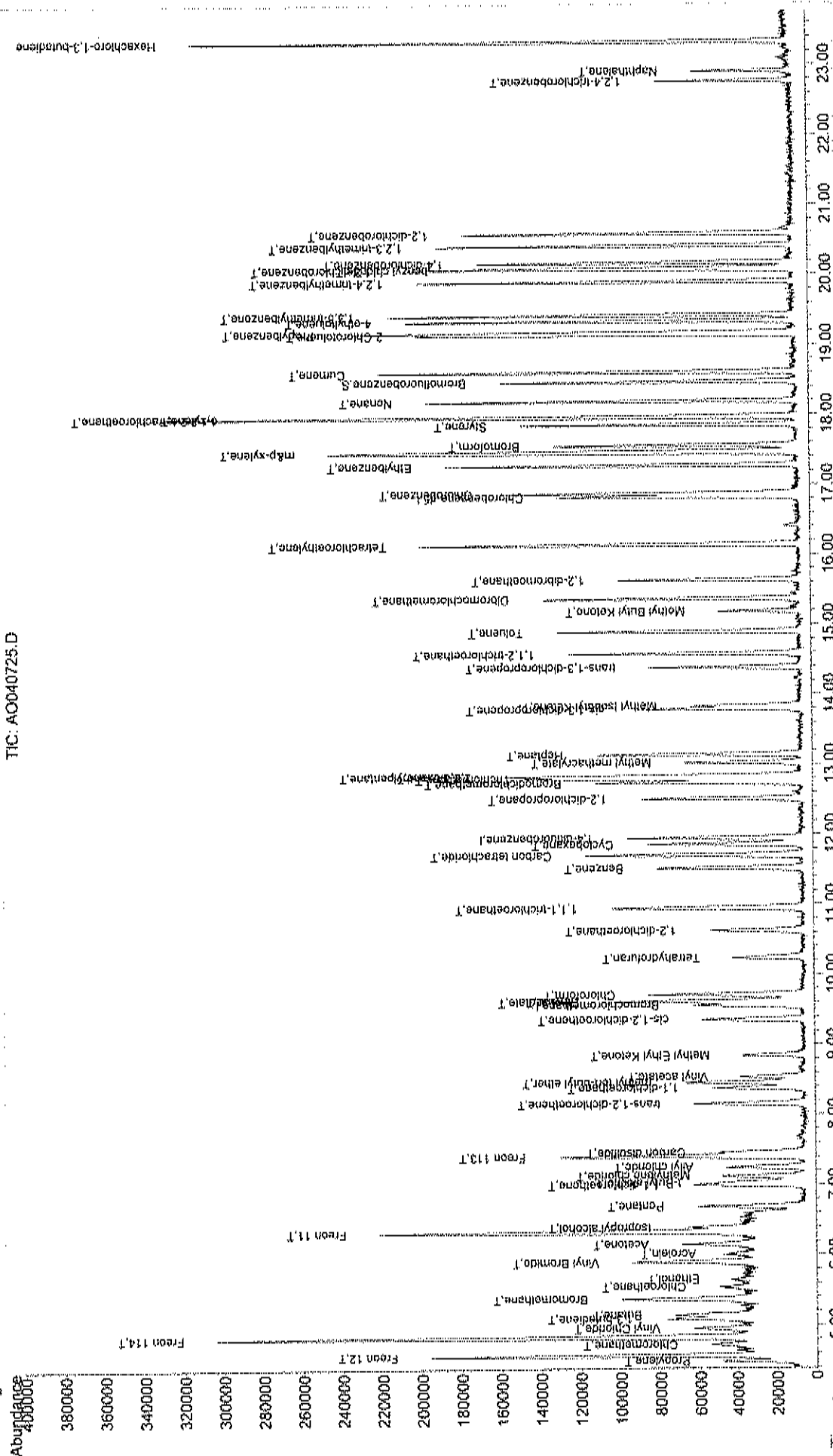
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.73	83	86612	1.10	ppb	95
47) cis-1,3-dichloropropene	13.79	75	56238	1.12	ppb	98
48) trans-1,3-dichloropropene	14.38	75	52555	1.10	ppb	97
49) 1,1,2-trichloroethane	14.58	97	42686	1.17	ppb	98
51) Toluene	14.90	92	62097	1.06	ppb	# 81
52) Methyl Isobutyl Ketone	13.83	43	44573	0.73	ppb	92
53) Dibromochloromethane	15.37	129	80598	1.13	ppb	88
54) Methyl Butyl Ketone	15.19	43	38670m	0.71	ppb	
55) 1,2-dibromoethane	15.64	107	68997	1.16	ppb	95
56) Tetrachloroethylene	16.14	164	45478	1.10	ppb	95
57) Chlorobenzene	16.88	112	93752	1.15	ppb	97
58) Ethylbenzene	17.27	91	149543	1.09	ppb	98
59) m&p-xylene	17.46	91	261402	2.27	ppb	99
60) Nonane	18.19	43	70846	1.11	ppb	93
61) Styrene	17.86	104	79302	1.10	ppb	91
62) Bromoform	17.55	173	64959	1.09	ppb	97
63) o-xylene	17.97	91	123539	1.11	ppb	96
64) Cumene	18.60	105	158840	1.10	ppb	100
66) 1,1,2,2-tetrachloroethane	17.96	83	91941	1.21	ppb	94
67) Propylbenzene	19.17	120	41612	1.12	ppb	# 1
68) 2-Chlorotoluene	19.14	126	39508	1.15	ppb	# 1
69) 4-ethyltoluene	19.34	105	146026	1.08	ppb	99
70) 1,3,5-trimethylbenzene	19.42	105	131377	1.09	ppb	99
71) 1,2,4-trimethylbenzene	19.90	105	118807	1.06	ppb	96
72) 1,3-dichlorobenzene	20.09	146	82665	1.17	ppb	99
73) benzyl chloride	20.07	91	58401	0.99	ppb	90
74) 1,4-dichlorobenzene	20.17	146	75820	1.16	ppb	93
75) 1,2,3-trimethylbenzene	20.41	105	116294	1.03	ppb	95
76) 1,2-dichlorobenzene	20.59	146	81959	1.21	ppb	96
77) 1,2,4-trichlorobenzene	22.76	180	21679	1.08	ppb	92
78) Naphthalene	22.91	128	42909	0.83	ppb	94
79) Hexachloro-1,3-butadiene	23.33	225	54811	1.09	ppb	99

Data File : C:\HPCHEM\1\DATA\A0040725.D  
Acq On : 8 Apr 2017 2:53 am  
Sample : ALC51UGD-040717  
Misc : A331\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Apr 12 8:36 2017

Method : C:\HPCHEM\1\METHODS\A331\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Thu May 04 11:27:28 2017  
Response via : Initial Calibration

Quant Results File: A331\_1UG.RES

TIC: A0040725.D



**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**INJECTION LOG**

Directory: C:\HPCHEM\1\DATA2

## Injection Log

Instrument # 1  
 Internal Standard Stock # A1937  
 Standard Stock # 1938  
 LCS Stock # 1939  
 Method Ref: EPA TO-15 / Injected

ie	Vial	FileName	Multiplier	SampleName	Misc Info	Ref: EPA TO-15 / Injected
1	4	Ao032904.d	1.	ALCS1UG-032917	A312_1UG	29 Mar 2017 11:59
2	5	Ao032905.d	1.	AMB1UG-032917	A312_1UG	29 Mar 2017 12:35
3	6	Ao032906.d	1.	C1703058-002A	A312_1UG	29 Mar 2017 13:15
4	7	Ao032907.d	1.	C1703058-004A	A312_1UG	29 Mar 2017 13:55
5	8	Ao032908.d	1.	C1703058-005A	A312_1UG	29 Mar 2017 14:35
6	9	Ao032909.d	1.	C1703058-002A 20X	A312_1UG	29 Mar 2017 15:19
7	10	Ao032910.d	1.	C1703058-004A 20X	A312_1UG	29 Mar 2017 15:56
8	11	Ao032911.d	1.	C1703058	A312_1UG	29 Mar 2017 16:33
9	12	Ao032912.d	1.	C1703078-001A 4X	A312_1UG	29 Mar 2017 17:11
0	13	Ao032913.d	1.	C1703078-002A 4X	A312_1UG	29 Mar 2017 17:49
1	14	Ao032914.d	1.	C1703078-003A 4X	A312_1UG	29 Mar 2017 18:27
2	15	Ao032915.d	1.	C1703079-001A 4X	A312_1UG	29 Mar 2017 19:04
3	16	Ao032916.d	1.	C1703079-002A 4X	A312_1UG	29 Mar 2017 19:42
4	17	Ao032917.d	1.	C1703079-003A 4X	A312_1UG	29 Mar 2017 20:20
5	18	Ao032918.d	1.	C1703080-001A 4X	A312_1UG	29 Mar 2017 20:57
6	19	Ao032919.d	1.	C1703080-002A 4X	A312_1UG	29 Mar 2017 21:35
7	20	Ao032920.d	1.	C1703080-003A 4X	A312_1UG	29 Mar 2017 22:13
8	21	Ao032921.d	1.	C1703058	A312_1UG	29 Mar 2017 22:50
9	21	Ao032922.d	1.	ALCS1UGD-032917	A312_1UG	29 Mar 2017 23:29
0	30	Ao032923.d	1.	C1703058-005A 5X	A312_1UG	30 Mar 2017 09:49
1	31	Ao032924.d	1.		A312_1UG	30 Mar 2017 15:15
2	32	Ao032925.d	1.		A312_1UG	30 Mar 2017 15:55
3	1	Ao033101.d	1.	BFB1UG	A331_1UG	31 Mar 2017 13:54
4	1	Ao033102.d	1.	A1UG	A331_1UG	31 Mar 2017 14:52
5	2	Ao033103.d	1.	A1UG_2.0	A331_1UG	31 Mar 2017 15:35
6	3	Ao033104.d	1.	A1UG_1.5	A331_1UG	31 Mar 2017 16:16
7	4	Ao033105.d	1.	A1UG_1.25	A331_1UG	31 Mar 2017 16:57
8	5	Ao033106.d	1.	A1UG_1.0	A331_1UG	31 Mar 2017 17:36
9	6	Ao033107.d	1.	A1UG_0.75	A331_1UG	31 Mar 2017 18:15
0	7	Ao033108.d	1.	A1UG_0.50	A331_1UG	31 Mar 2017 18:53
1	8	Ao033109.d	1.	A1UG_0.30	A331_1UG	31 Mar 2017 19:30
2	9	Ao033110.d	1.	A1UG_0.15	A331_1UG	31 Mar 2017 20:07
3	10	Ao033111.d	1.	A1UG_0.10	A331_1UG	31 Mar 2017 20:44
4	11	Ao033112.d	1.	A1UG_0.04	A331_1UG	31 Mar 2017 21:20
5	1	Ao033113.d	1.	IDL1UG	A331_1UG 0.50PPB	31 Mar 2017 22:32
6	2	Ao033114.d	1.	IDL1UG	A331_1UG 0.50PPB	31 Mar 2017 23:10
7	3	Ao033115.d	1.	IDL1UG	A331_1UG 0.50PPB	31 Mar 2017 23:48
8	4	Ao033116.d	1.	IDL1UG	A331_1UG 0.50PPB	1 Apr 2017 00:26
9	5	Ao033117.d	1.	IDL1UG	A331_1UG 0.50PPB	1 Apr 2017 01:04
0	6	Ao033118.d	1.	IDL1UG	A331_1UG 0.50PPB	1 Apr 2017 01:42
1	7	Ao033119.d	1.	IDL1UG	A331_1UG 0.50PPB	1 Apr 2017 02:20
2	8	Ao033120.d	1.	IDL1UG	A331_1UG 0.50PPB	1 Apr 2017 02:58
3	1	Ao033121.d	1.	IDL1UG	A331_1UG 0.10PPB	1 Apr 2017 03:35
4	2	Ao033122.d	1.	IDL1UG	A331_1UG 0.10PPB	1 Apr 2017 04:11
5	3	Ao033123.d	1.	IDL1UG	A331_1UG 0.10PPB	1 Apr 2017 04:48
6	4	Ao033124.d	1.	IDL1UG	A331_1UG 0.10PPB	1 Apr 2017 05:25
7	5	Ao033125.d	1.	IDL1UG	A331_1UG 0.10PPB	1 Apr 2017 06:01
8	6	Ao033126.d	1.	IDL1UG	A331_1UG 0.10PPB	1 Apr 2017 06:38
9	7	Ao033127.d	1.	IDL1UG	A331_1UG 0.10PPB	1 Apr 2017 07:14
0	8	Ao033128.d	1.	IDL1UG	A331_1UG 0.10PPB	1 Apr 2017 07:51
1	1	Ao033129.d	1.	IDL1UG	A331_1UG 0.30PPB	1 Apr 2017 08:28
2	2	Ao033130.d	1.	IDL1UG	A331_1UG 0.30PPB	1 Apr 2017 09:06
3	3	Ao033131.d	1.	IDL1UG	A331_1UG 0.30PPB	1 Apr 2017 09:43
4	4	Ao033132.d	1.	IDL1UG	A331_1UG 0.30PPB	1 Apr 2017 10:20
5	5	Ao033133.d	1.	IDL1UG	A331_1UG 0.30PPB	1 Apr 2017 10:58

# Injection Log

Instrument # 1  
 Internal Standard Stock # A1949  
 Standard Stock # A1950  
 LCS Stock # A1951  
 Method Ref: EPA 70-15 / Jan 1990

Directory: C:\HPCHEM\1\DATA2

ne	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
3	22	Ao040422.d	1.	ALCS1UGD-040417	A331_1UG	5 Apr 2017 00:07
7	23	Ao040423.d	1.	C1704004-005A 10X	A331_1UG	5 Apr 2017 00:46
3	24	Ao040424.d	1.	C1704004-001A 10X	A331_1UG	5 Apr 2017 01:26
9	25	Ao040425.d	1.	C1704004-001A 40X	A331_1UG	5 Apr 2017 02:05
0		Ao040426.d	1.	No MS or GC data present		
1	1	Ao040501.d	1.	BFB1UG	A331_1UG	5 Apr 2017 09:11
2	2	Ao040502.d	1.	A1UG_1.0	A331_1UG	5 Apr 2017 10:02
3	3	Ao040503.d	1.	ALCS1UG-040517	A331_1UG	5 Apr 2017 10:42
4	4	Ao040504.d	1.	AMB1UG-040517	A331_1UG	5 Apr 2017 11:18
5	5	Ao040505.d	1.	C1704004-002A 10X	A331_1UG	5 Apr 2017 11:58
6	6	Ao040506.d	1.	C1704004-002A 40X	A331_1UG	5 Apr 2017 12:34
7	7	Ao040507.d	1.	C1704005-001A	A331_1UG	5 Apr 2017 13:14
8	8	Ao040508.d	1.	C1704005-001A 20X	A331_1UG	5 Apr 2017 14:14
9	9	Ao040509.d	1.	WAC040517A	A331_1UG	5 Apr 2017 14:52
10	10	Ao040510.d	1.	WAC040517B N	A331_1UG	5 Apr 2017 15:45
11	11	Ao040511.d	1.	WAC040517C N	A331_1UG	5 Apr 2017 16:27
12	12	Ao040512.d	1.	WAC040517D	A331_1UG	5 Apr 2017 17:04
13	13	Ao040513.d	1.	WAC040517E	A331_1UG	5 Apr 2017 17:42
14	14	Ao040514.d	1.	WAC040517F	A331_1UG	5 Apr 2017 18:20
15	15	Ao040515.d	1.	WAC040517G	A331_1UG	5 Apr 2017 18:57
1	1	Ao040516.d	1.	C1704006-002A	A331_1UG	5 Apr 2017 19:37
2	2	Ao040517.d	1.	C1704006-003A	A331_1UG	5 Apr 2017 20:17
3	3	Ao040518.d	1.	C1704007-002A	A331_1UG	5 Apr 2017 20:57
4	4	Ao040519.d	1.	C1704007-004A	A331_1UG	5 Apr 2017 21:37
5	5	Ao040520.d	1.	C1704007-005A	A331_1UG	5 Apr 2017 22:17
6	6	Ao040521.d	1.	C1704008-002A	A331_1UG	5 Apr 2017 22:58
7	7	Ao040522.d	1.	C1704008-004A	A331_1UG	5 Apr 2017 23:39
8	8	Ao040523.d	1.	C1704008-005A	A331_1UG	6 Apr 2017 00:19
9	9	Ao040524.d	1.	C1704006-001A	A331_1UG	6 Apr 2017 01:00
10	10	Ao040525.d	1.	C1704007-001A	A331_1UG	6 Apr 2017 01:40
11	11	Ao040526.d	1.	C1704007-003A	A331_1UG	6 Apr 2017 02:20
12	12	Ao040527.d	1.	C1704008-001A	A331_1UG	6 Apr 2017 03:00
13	13	Ao040528.d	1.	C1704008-003A	A331_1UG	6 Apr 2017 03:42
14	14	Ao040529.d	1.	ALCS1UGD-040517	A331_1UG	6 Apr 2017 04:21
15	15	Ao040530.d	1.	C1704006-002A 10X	A331_1UG	6 Apr 2017 04:58
16	16	Ao040531.d	1.	C1704006-003A 10X	A331_1UG	6 Apr 2017 05:35
17	17	Ao040532.d	1.	C1704007-002A 10X	A331_1UG	6 Apr 2017 06:12
18	18	Ao040533.d	1.	C1704007-004A 10X	A331_1UG	6 Apr 2017 06:49
19	19	Ao040534.d	1.	C1704007-005A 10X	A331_1UG	6 Apr 2017 07:25
		Ao040535.d	1.	No MS or GC data present		
1	1	Ao040601.d	1.	BFB1UG	A331_1UG	6 Apr 2017 08:51
2	2	Ao040602.d	1.	A1UG_1.0	A331_1UG	6 Apr 2017 09:31
3	3	Ao040603.d	1.	ALCS1UG-040617	A331_1UG	6 Apr 2017 10:10
4	4	Ao040604.d	1.	AMB1UG-040617	A331_1UG	6 Apr 2017 10:47
0	5	Ao040605.d	1.	C1704008-002A 10X	A331_1UG	6 Apr 2017 11:24
1	6	Ao040606.d	1.	C1704008-004A 10X	A331_1UG	6 Apr 2017 12:01
2	7	Ao040607.d	1.	C1704008-005A 10X	A331_1UG	6 Apr 2017 12:38
3	8	Ao040608.d	1.	C1704006-001A 10X	A331_1UG	6 Apr 2017 13:15
4	9	Ao040609.d	1.	C1704007-001A 10X	A331_1UG	6 Apr 2017 13:59
5	10	Ao040610.d	1.	C1704007-003A 10X	A331_1UG	6 Apr 2017 14:36
6	11	Ao040611.d	1.	C1704008-001A 10X	A331_1UG	6 Apr 2017 15:13
7	12	Ao040612.d	1.	C1704008-003A 10X	A331_1UG	6 Apr 2017 15:50
3	42	Ao040613.d	1.	C1704013-001A	A331_1UG	6 Apr 2017 16:32
9	43	Ao040614.d	1.	C1704013-002A	A331_1UG	6 Apr 2017 17:12
0	44	Ao040615.d	1.	C1704013-003A	A331_1UG	6 Apr 2017 17:54



Directory: C:\HPCHEM\1\DATA2

## Injection Log

Instrument # 1  
 Internal Standard Stock # A1949  
 Standard Stock # A1550  
 LCS Stock # A1551  
 Method Ref: EPA TO-15 / Jan. 1996

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	45	Ao040616.d	1.	C1704013-004A	A331_1UG	6 Apr 2017 18:33
2	46	Ao040617.d	1.	C1704013-005A	A331_1UG	6 Apr 2017 19:14
3	47	Ao040618.d	1.	C1704013-007A	A331_1UG	6 Apr 2017 19:55
4	48	Ao040619.d	1.	C1704013-008A	A331_1UG	6 Apr 2017 20:35
5	49	Ao040620.d	1.	ALCS1UGD-040617	A331_1UG	6 Apr 2017 21:15
6	21	Ao040621.d	1.	C1704014-001A	A331_1UG	6 Apr 2017 21:55
7	22	Ao040622.d	1.	C1704014-001A MS	A331_1UG	6 Apr 2017 22:40
8	23	Ao040623.d	1.	C1704014-001A MSD	A331_1UG	6 Apr 2017 23:26
9	22	Ao040624.d	1.	C1704014-002A	A331_1UG	7 Apr 2017 00:09
0	23	Ao040625.d	1.	C1704014-004A	A331_1UG	7 Apr 2017 00:50
1	24	Ao040626.d	1.	C1704014-006A	A331_1UG	7 Apr 2017 01:31
2	25	Ao040627.d	1.	C1704014-007A	A331_1UG	7 Apr 2017 02:12
3	26	Ao040628.d	1.	C1704014-008A	A331_1UG	7 Apr 2017 02:52
4	27	Ao040629.d	1.	C1704014-010A	A331_1UG	7 Apr 2017 03:33
5	28	Ao040630.d	1.	C1704014-012A	A331_1UG	7 Apr 2017 04:13
6	29	Ao040631.d	1.	C1704014-013A	A331_1UG	7 Apr 2017 04:53
7	30	Ao040632.d	1.	C1704013-001A 10X	A331_1UG	7 Apr 2017 05:30
8	31	Ao040633.d	1.	C1704013-002A 10X	A331_1UG	7 Apr 2017 06:07
9	32	Ao040634.d	1.	C1704013-003A 10X	A331_1UG	7 Apr 2017 06:44
0	33	Ao040635.d	1.	C1704013-004A 10X	A331_1UG	7 Apr 2017 07:21
1	34	Ao040636.d	1.	C1704013-005A 10X	A331_1UG	7 Apr 2017 07:58
2	35	Ao040637.d	1.	C1704013-007A 10X	A331_1UG	7 Apr 2017 08:35
3	36	Ao040638.d	1.	C1704013-008A 10X	A331_1UG	7 Apr 2017 09:12
4		Ao040639.d	1.	No MS or GC data present		
5	1	Ao040701.d	1.	BFB1UG	A331_1UG	7 Apr 2017 10:27
6	2	Ao040702.d	1.	A1UG_1.0	A331_1UG	7 Apr 2017 11:20
7	3	Ao040703.d	1.	ALCS1UG-040717	A331_1UG	7 Apr 2017 12:00
8	4	Ao040704.d	1.	AMB1UG-040717	A331_1UG	7 Apr 2017 12:36
9	5	Ao040705.d	1.	C1704013-001A 90X	A331_1UG	7 Apr 2017 13:13
0	6	Ao040706.d	1.	C1704014-012A 10X	A331_1UG	7 Apr 2017 13:50
1	7	Ao040707.d	1.	C1704014-013A 10X	A331_1UG	7 Apr 2017 14:27
2	8	Ao040708.d	1.	C1704014-003A	A331_1UG	7 Apr 2017 15:23
3	9	Ao040709.d	1.	C1704014-005A	A331_1UG	7 Apr 2017 16:03
4	10	Ao040710.d	1.	C1704014-009A	A331_1UG	7 Apr 2017 17:28
5	11	Ao040711.d	1.	C1704014-011A	A331_1UG	7 Apr 2017 18:09
6	12	Ao040712.d	1.	C1704014-014A	A331_1UG	7 Apr 2017 18:50
7	13	Ao040713.d	1.	C1704014-015A	A331_1UG	7 Apr 2017 19:30
8	14	Ao040714.d	1.	C1704014-003A 10x	A331_1UG	7 Apr 2017 20:07
9	15	Ao040715.d	1.	C1704014	A331_1UG -005A 9x	7 Apr 2017 20:44
0	16	Ao040716.d	1.	C1704014-005A 90x	A331_1UG	7 Apr 2017 21:20
1	17	Ao040717.d	1.	C1704014	A331_1UG -009A 10x	7 Apr 2017 21:57
2	18	Ao040718.d	1.	C1704014	A331_1UG -009A 10x	7 Apr 2017 22:33
3	19	Ao040719.d	1.	C1704014-011A 10x	A331_1UG	7 Apr 2017 23:10
4	20	Ao040720.d	1.	C1704014	A331_1UG -011A 40x	7 Apr 2017 23:47
5	21	Ao040721.d	1.	C1704014-014A 9x	A331_1UG	8 Apr 2017 00:23
6	22	Ao040722.d	1.	C1704014-014A 90x	A331_1UG	8 Apr 2017 01:00
7	23	Ao040723.d	1.	C1704014-015A 10x	A331_1UG	8 Apr 2017 01:37
8	24	Ao040724.d	1.	C1704014	A331_1UG -015A 40x	8 Apr 2017 02:13
9	25	Ao040725.d	1.	ALCS1UGD-040717	A331_1UG	8 Apr 2017 02:53
0		Ao040726.d	1.	No MS or GC data present		
1	1	Ao041001.d	1.	BFB1UG	A331_1UG	10 Apr 2017 11:01
2	2	Ao041002.d	1.	A1UG_1.0	A331_1UG	10 Apr 2017 11:44
3	3	Ao041003.d	1.	ALCS1UG-041017	A331_1UG	10 Apr 2017 12:39
4	4	Ao041004.d	1.	AMB1UG-041017	A331_1UG	10 Apr 2017 13:19
5	5	Ao041005.d	1.	WAC041017A	A331_1UG	10 Apr 2017 14:09

**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**STANDARDS LOG**

Std #	Date Prep	Date Exp	Description	Stock #	Stock Conc	Initial Vol (psig)	Final Vol (psia)	Final Conc (ppb)	Prep by	Chkd b
A-1788	12/22/16	12/29/16	TO15 STLX	<del>A1088</del> A1089	500 ppb	3.0		50	LL	
A-1789			↓ SULF	A0270	1 ppm	1.5		50		
A-1790			↓ H2S	A0269	10 ppm	1.5		500		
A-1791			TO15 NG IS	A1782	50 ppb	0.9		1		
A-1792			↓ STD	A1783	↓	↓		↓		
A-1793			↓ LCS	A1784	↓	↓		↓		
A-1794	12/29/16	1/5/17	TO15 IS	A1289	1 ppm	1.5		50	WD	
A-1795			STD	A1203	↓	↓		↓		
A-1796			LCS	A1204	↓	↓		↓		
A-1797			4PCH	9519	↓	↓		↓		
A-1798			4PCH5	A1797	50 ppb	3.0		5		
A-1799			FORM	A0974	11.5 ppm	0.20		50		
A-1800			SILOX	<del>A1088</del> A1089	500 ppb	3.0		50		
A-1801			SULF	A0270	1 ppm	1.5		50		
A-1802			H2S	A0269	10 ppm	1.5		500		
A-1803			TO15 NG IS	A1794	50 ppb	0.9		1		
A-1804			↓ STD	A1795	↓	↓		↓		
A-1805			↓ LCS	A1796	↓	↓		↓		
A-1806	1/5/17	1/5/18	TO15 IS	FF-47206	LINDE			1 ppm	WD	
A-1807	1/5/17	1/5/18	STOCK TO15 STD	FF-45347	LINDE			1 ppm	WD	
A-1808	1/6/17	1/6/18	TO15 <del>SS</del> LCS	<del>A1796</del> A1203	1 ppm	1.5		50 ppb	M	

FORM 153

A1203 570 IS NOW LCS

Page #

7

Std #	Date Prep	Date Exp	Description	Stock #	Stock Conc	Initial Vol (psig)	Final Vol (psia)	Final Conc (ppb)	Prep by	Chkd by
A-1934	3/23/17	4/4/17	TO15	<del>A1088</del> A1089	500ppb	3.0	30	50	WJD	
A-1935			↓	A0270	1ppm	1.5	30	50		
A-1936			↓	A0269	10ppm	1.5	30	500		
A-1937			TO15 149	A1928	50ppb	0.9	45	4		
A-1938			↓	A1929	↓	↓	↓	↓		
A-1939			↓	A1930	↓	↓	↓	↓		
A-1940	4/5/17	4/15/17	TO15	A1806	1ppm	1.5	30.0ppm	50	WJD	
A-1941			↓	A1807	↓	↓	↓	↓		
A-1942			↓	A1808	↓	↓	↓	↓		
A-1943			4PC4	9515	↓	↓	↓	↓		
A-1944			4PC45	A1543	50	3.0	↓	↓		
A-1945			FOAM	A0174	11.5ppm	0.2	45	50		
A-1946			SILOX	A1088/1089	500ppb	3.0	36	50		
A-1947			↓	A0270	1ppm	1.5	↓	↓		
A-1948			↓	A0269	10ppm	↓	↓	500		
A-1949			TO15 109	A1940	50ppb	0.9	45	↓		
A-1950			↓	A1541	↓	↓	↓	↓		
A-1951			↓	A1542	↓	↓	↓	↓		
A-1952										
A-1953										
A-1954										

**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**CANISTER CLEANING LOG**

Instrument: Entech 3100

Canister Number	Canister Size	QC Can Number	# of Cycles	Int & Date	QC Batch Number	Detection Limits	Leak Test 24hr Int & Date
163	1L	94	20	3/8/17	WAC031317A	1 µg +0.25	+ 30 + 3/20/17 30
550							+ 30
248							+ 30
240							+ 30
94		✓					+ 30
162		327			B		+ 30
362							+ 30
170							+ 30
95							+ 30
327		✓					+ 30
1318		1193			C		+ 30
569							+ 30
552							+ 30
317							+ 30
1193		✓					+ 30
193		1190			D		+ 30
556							+ 30
229							+ 30
131							+ 30
1190		✓					+ 30
1191		1184			E		+ 30
93							+ 30
599		✓					+ 30
562							+ 30
1184		✓					+ 30

**entek Laboratories, LLC**

Instrument: Entech 3100

[illegible]



Instrument: Entech 3100

Canister Number	Canister Size	QC Can Number	# of Cycles	Int & Date Cleaned	QC Batch Number	Detection Limits	Leak Test 24hr Int & Date
359	1L	89	20		WAC033017A	1ug/m <sup>3</sup> + 0.25	+ 30 RSP 03/20/17
222							+ 30
546							+ 30
363							+ 30
89							+ 30
96		136			WAC032017B		+ 30
1177							+ 30
245							+ 30
1195							+ 30
136							+ 30
83		137			WAC032017C		+ 30
171							+ 30
242							+ 30
233							+ 30
137							+ 30
106		285			WAC032017D		+ 30
467							+ 30
1188							+ 30
130							+ 30
285							+ 30
367		1186			WAC032017E		+ 30
328							+ 30
139							+ 30
286							+ 30
1186							+ 30



Data File : C:\HPCHEM\1\DATA2\AO031310.D

Vial: 26

Acq On : 13 Mar 2017 1:56 pm

Operator: RJP

Sample : WAC031317A

Inst : MSD #1

Misc : A312\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 14 10:02:15 2017

Quant Results File: A312\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Sun Mar 12 22:26:37 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	60201	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.92	114	294080	1.00	ppb	-0.01
50) Chlorobenzene-d5	16.81	117	241427	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.45	95	175772	0.97	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	97.00%

Target Compounds

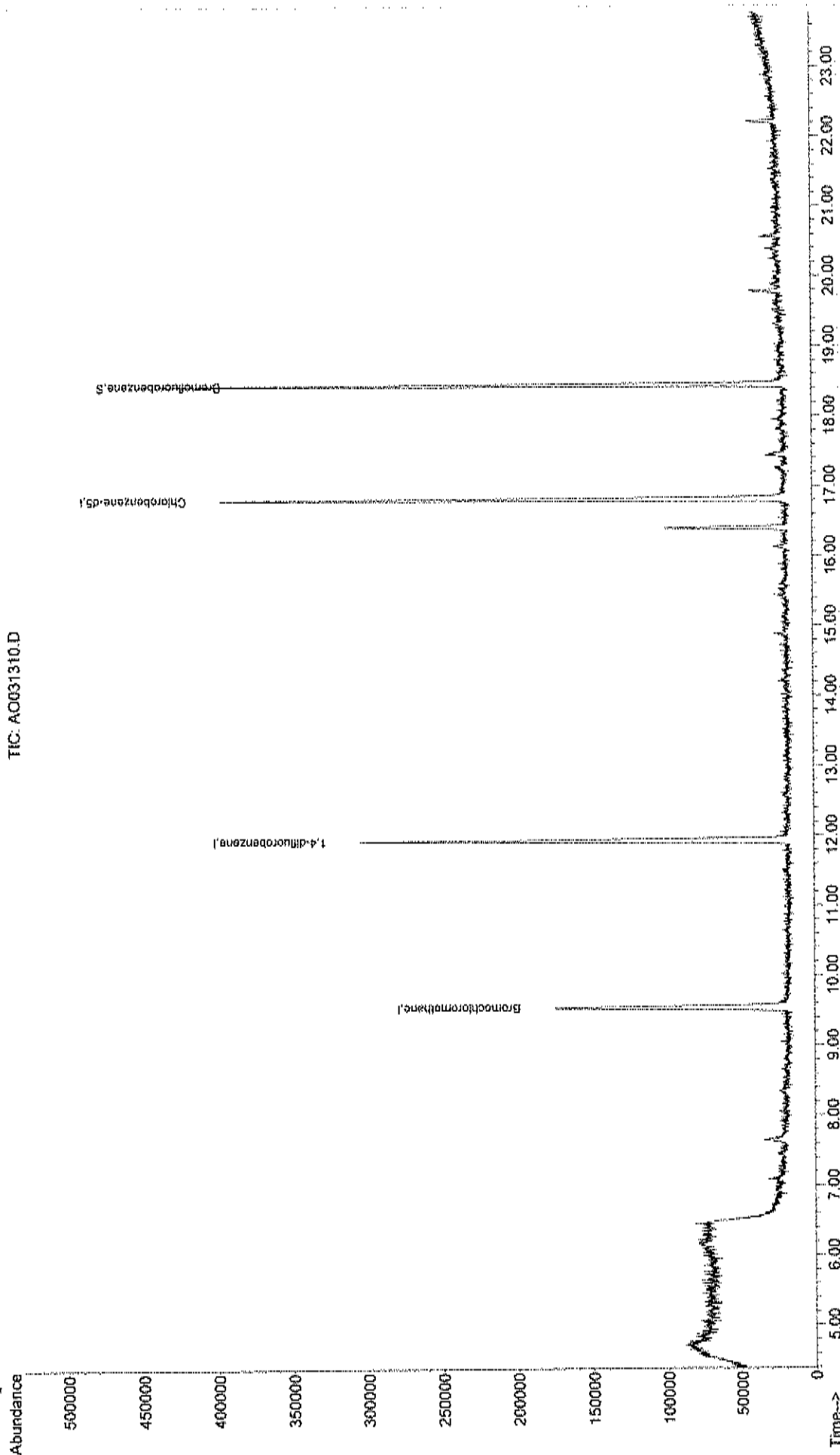
Qvalue

Data File : C:\HPCHEM\1\DATA2\AO031310.D  
 Acq On : 13 Mar 2017 1:56 pm  
 Sample : WAC031317A  
 Misc : A312\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Mar 14 10:18 2017

Vial: 26  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A312\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A426\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 14:17:03 2017  
 Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA2\AO031311.D  
Acq On : 13 Mar 2017 2:34 pm  
Sample : WAC031317B  
Misc : A312\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 14 10:02:16 2017

Vial: 27  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A312\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Sun Mar 12 22:26:37 2017  
Response via : Initial Calibration  
DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	62194	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.93	114	285996	1.00	ppb	0.00
50) Chlorobenzene-d5	16.81	117	238418	1.00	ppb	0.00

System Monitoring Compounds

65) Bromofluorobenzene	18.45	95	171452	0.96	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	96.00%

Target Compounds

Qvalue

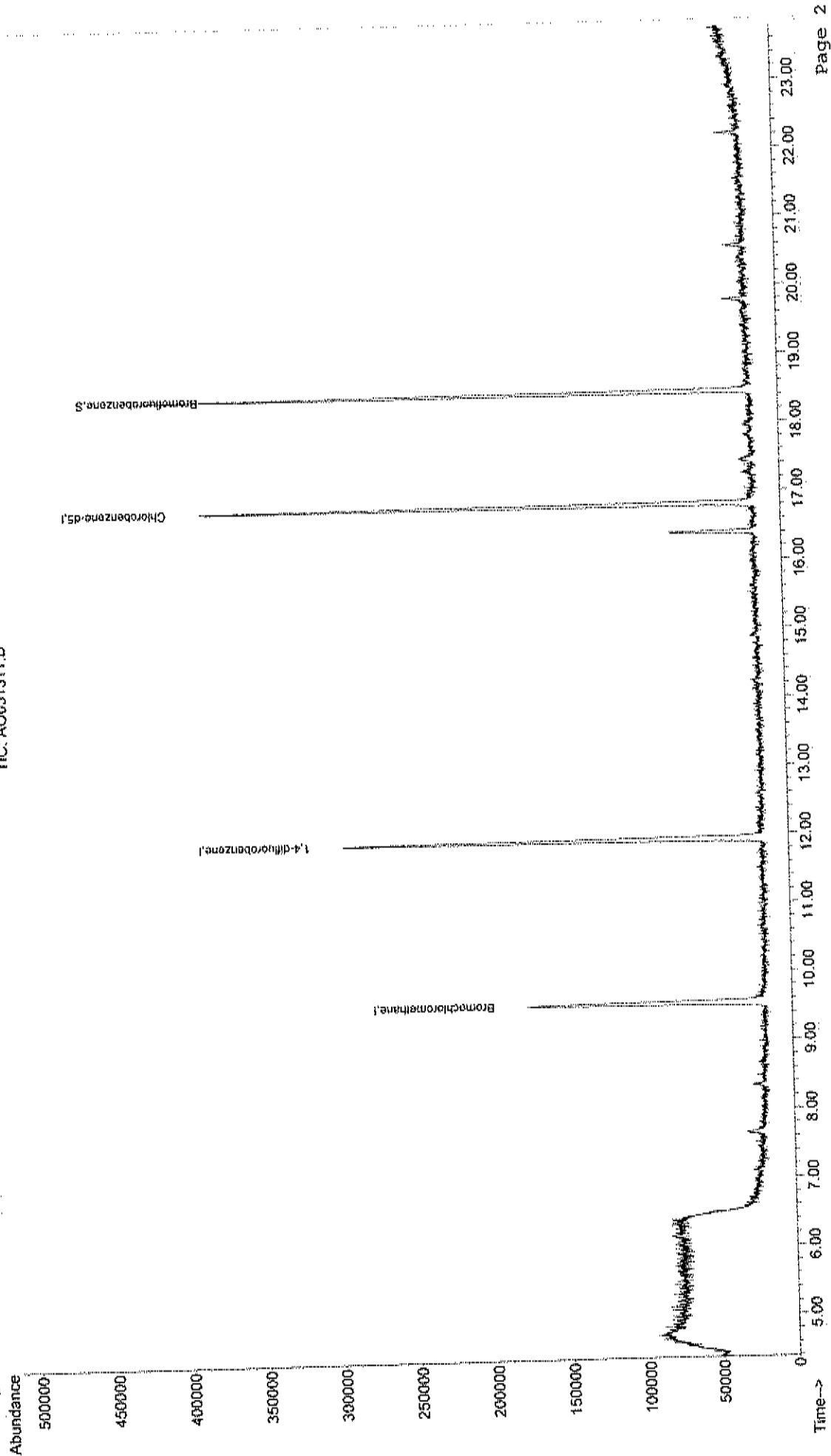
Data File : C:\HPCHEM\1\DATA2\AO031311.D  
Acq On : 13 Mar 2017 2:34 pm  
Sample : WAC031317B  
Misc : A312\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 14 10:18 2017

Vial: 27  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A312\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A426\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Thu May 04 14:17:03 2017  
Response via : Initial Calibration

TIC: AO031311.D



MSD1

Mon May 08 10:44:04 2017

AO031311.D A426\_1UG.M

Data File : C:\HPCHEM\1\DATA2\AO031312.D

Vial: 28

Acq On : 13 Mar 2017 3:12 pm

Operator: RJP

Sample : WAC031317C

Inst : MSD #1

Misc : A312\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 14 10:02:17 2017

Quant Results File: A312\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Sun Mar 12 22:26:37 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	58213	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.93	114	283013	1.00	ppb	0.00
50) Chlorobenzene-d5	16.81	117	232802	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.45	95	174331	1.00	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	100.00%

Target Compounds

Qvalue

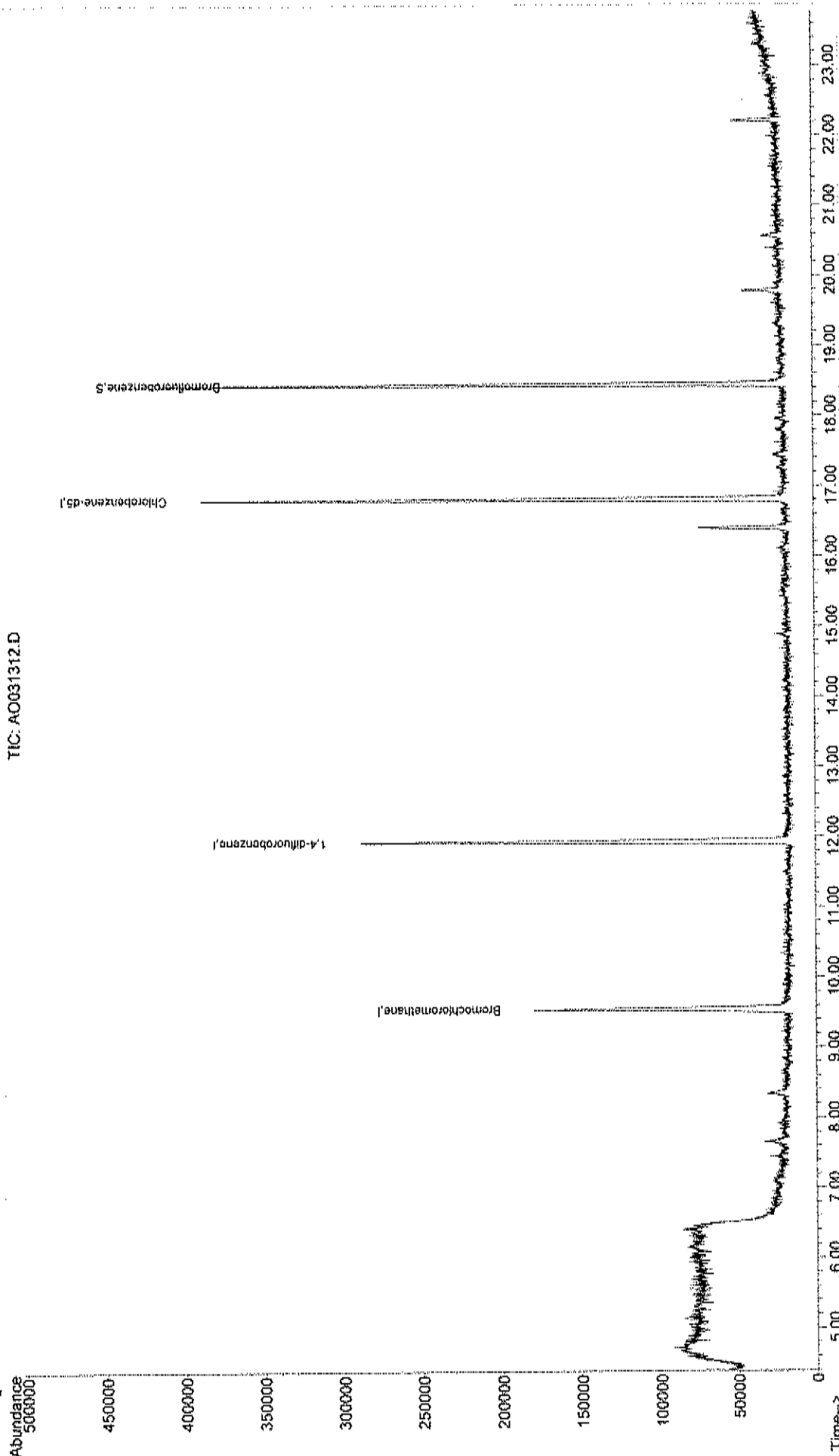


Data File : C:\HPCHEM\1\DATA2\AO031312.D  
 Acq On : 13 Mar 2017 3:12 pm  
 Sample : WAC031317C  
 Misc : A312\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Mar 14 10:02 2017

Vial: 28  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A312\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A426\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 14:17:03 2017  
 Response via : Initial Calibration



TIC: AO031312.D

Data File : C:\HPCHEM\1\DATA2\AO031313.D

Vial: 29

Acq On : 13 Mar 2017 3:49 pm

Operator: RJP

Sample : WAC031317D

Inst : MSD #1

Misc : A312\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 14 10:02:18 2017

Quant Results File: A312\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Sun Mar 12 22:26:37 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	56611	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.93	114	277027	1.00	ppb	0.00
50) Chlorobenzene-d5	16.81	117	225435	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.44	95	169439	1.01	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	101.00%

Target Compounds

Qvalue

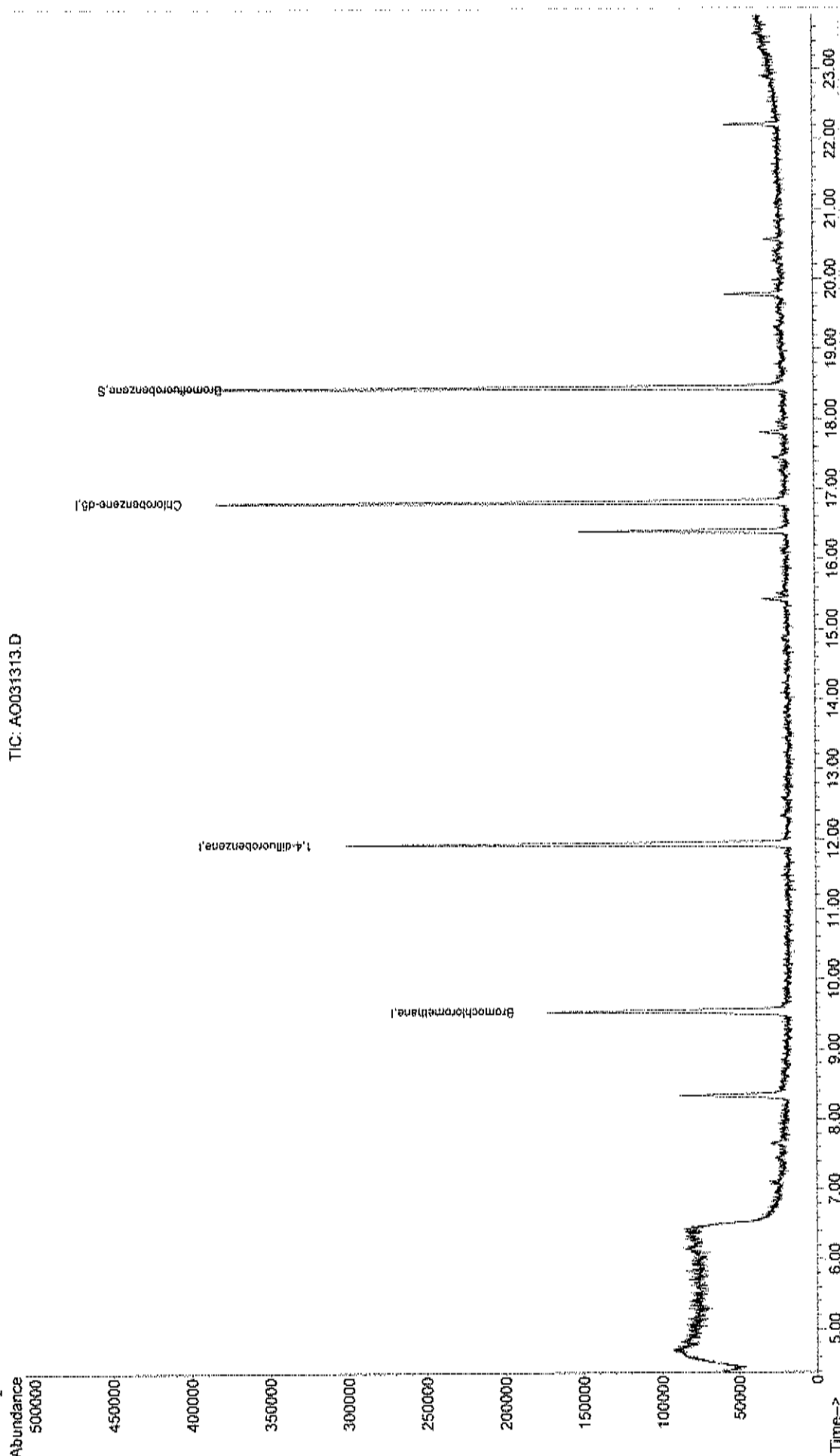
Data File : C:\HPCHEM\1\DATA2\AO031313.D  
Acq On : 13 Mar 2017 3:49 pm  
Sample : WAC031317D  
Misc : A312\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 14 10:19 2017

Vial: 29  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A312\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A426\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Thu May 04 14:17:03 2017  
Response via : Initial Calibration

TIC: AO031313.D



Data File : C:\HPCHEM\1\DATA2\AO031314.D

Vial: 30

Acq On : 13 Mar 2017 4:27 pm

Operator: RJP

Sample : WAC031317E

Inst : MSD #1

Misc : A312\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 14 10:02:19 2017

Quant Results File: A312\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Sun Mar 12 22:26:37 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	56965	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.92	114	274359	1.00	ppb	-0.01
50) Chlorobenzene-d5	16.81	117	226756	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.45	95	169835	1.00	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	100.00%

Target Compounds

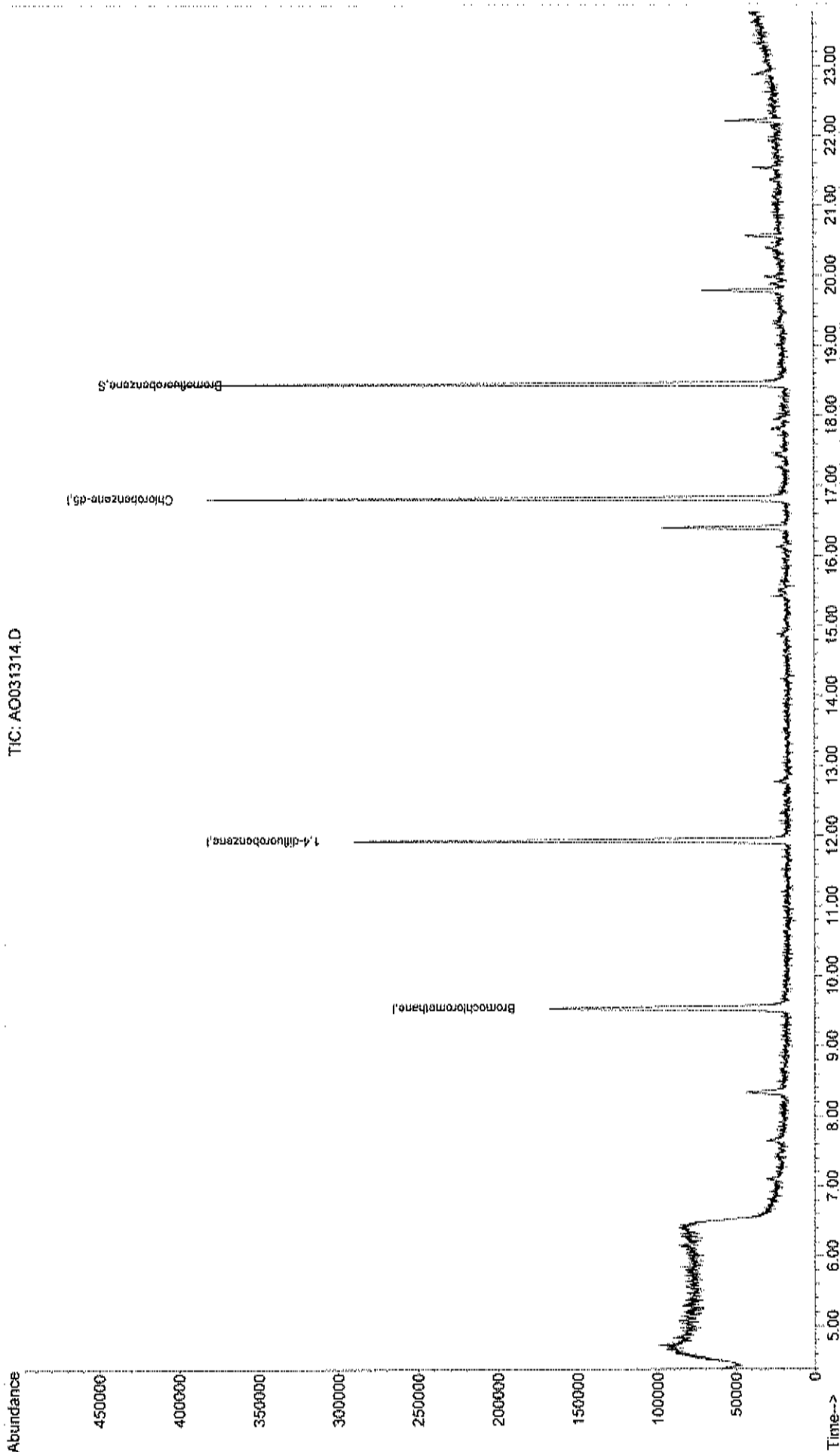
Qvalue

Data File : C:\HPCHEM\1\DATA2\AO031314.D  
 Acq On : 13 Mar 2017 4:27 pm  
 Sample : WAC031317E  
 Misc : A312\_IUG  
 MS Integration Params: RTEINT.P  
 Quant Time: Mar 14 10:19 2017

Vial: 30  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A312\_IUG.RES

Method : C:\HPCHEM\1\METHODS\A426\_IUG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 14:17:03 2017  
 Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA2\AO031315.D

Vial: 31

Acq On : 13 Mar 2017 5:05 pm

Operator: RJP

Sample : WAC031317F

Inst : MSD #1

Misc : A312\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 14 10:02:20 2017

Quant Results File: A312\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Sun Mar 12 22:26:37 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.53	128	55487	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.93	114	273365	1.00	ppb	0.00
50) Chlorobenzene-d5	16.81	117	226312	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.45	95	163837	0.97	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	97.00%

Target Compounds

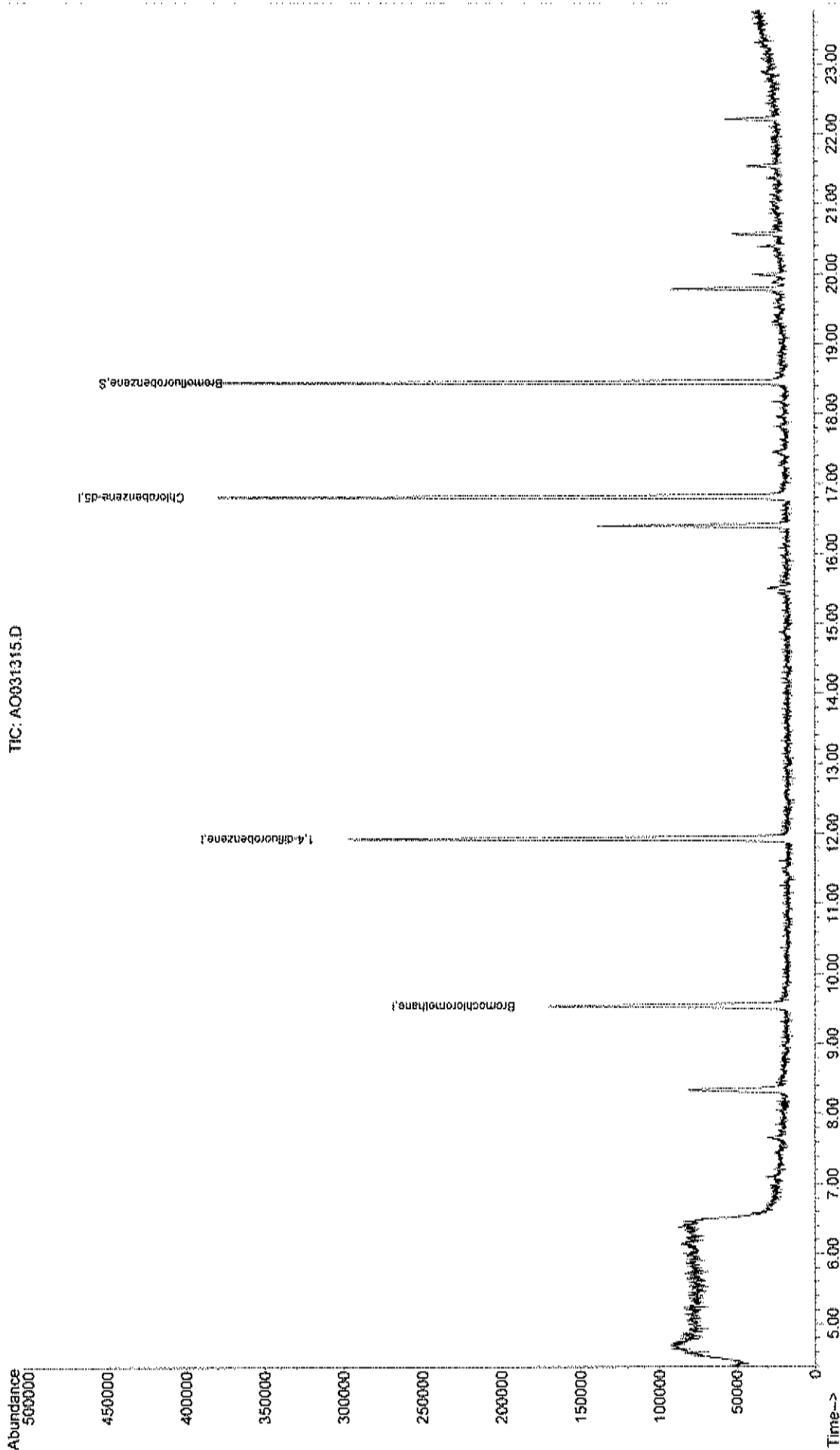
Qvalue

Data File : C:\HPCHEM\1\DATA2\AO031315.D  
 Acq On : 13 Mar 2017 5:05 pm  
 Sample : WAC031317F  
 Misc : A312\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Mar 14 10:02 2017

Vial: 31  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A312\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A426\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 14:17:03 2017  
 Response via : Initial Calibration





Data File : C:\HPCHEM\1\DATA2\AO032006.D  
Acq On : 20 Mar 2017 2:43 pm  
Sample : WAC032017A  
Misc : A312\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 21 09:03:02 2017

Vial: 1  
Operator: RJP  
Inst : MSD #1  
Multiplier: 1.00

Quant Results File: A312\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Wed Mar 15 10:58:20 2017  
Response via : Initial Calibration  
DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.53	128	15562	1.00	ppb	-0.02
35) 1,4-difluorobenzene	11.93	114	73218	1.00	ppb	-0.01
50) Chlorobenzene-d5	16.80	117	61485	1.00	ppb	-0.01

System Monitoring Compounds

65) Bromofluorobenzene	18.44	95	42443	0.92	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	92.00%

Target Compounds	Qvalue
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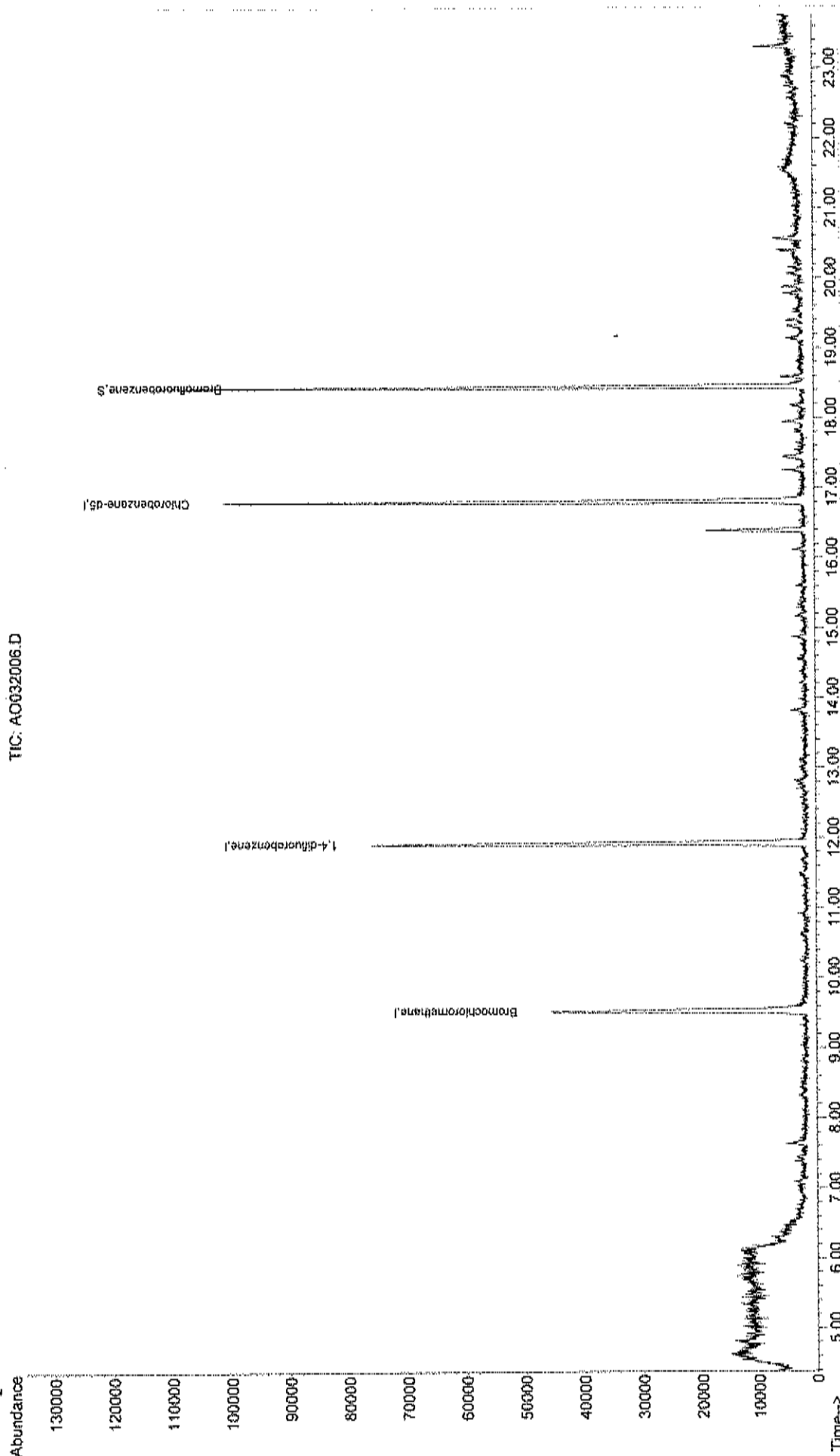
Data File : C:\HPCHEM\1\DATA2\AO032006.D  
Acq On : 20 Mar 2017 2:43 pm  
Sample : WAC032017A  
Misc : A312\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 21 9:03 2017

Vial: 1  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A312\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A426\_1UG.M (RTE Integrator)  
Title : FO-15 VOA Standards for 5 point calibration  
Last Update : Thu May 04 14:17:03 2017  
Response via : Initial Calibration

TIC: AO032006.D



Data File : C:\HPCHEM\1\DATA2\AO032007.D

Vial: 2

Acq On : 20 Mar 2017 3:21 pm

Operator: RJP

Sample : WAC032017B

Inst : MSD #1

Misc : A312\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 21 09:03:03 2017

Quant Results File: A312\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Wed Mar 15 10:58:20 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	14949	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.93	114	69738	1.00	ppb	0.00
50) Chlorobenzene-d5	16.81	117	57843	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.44	95	39683	0.92	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	92.00%

Target Compounds

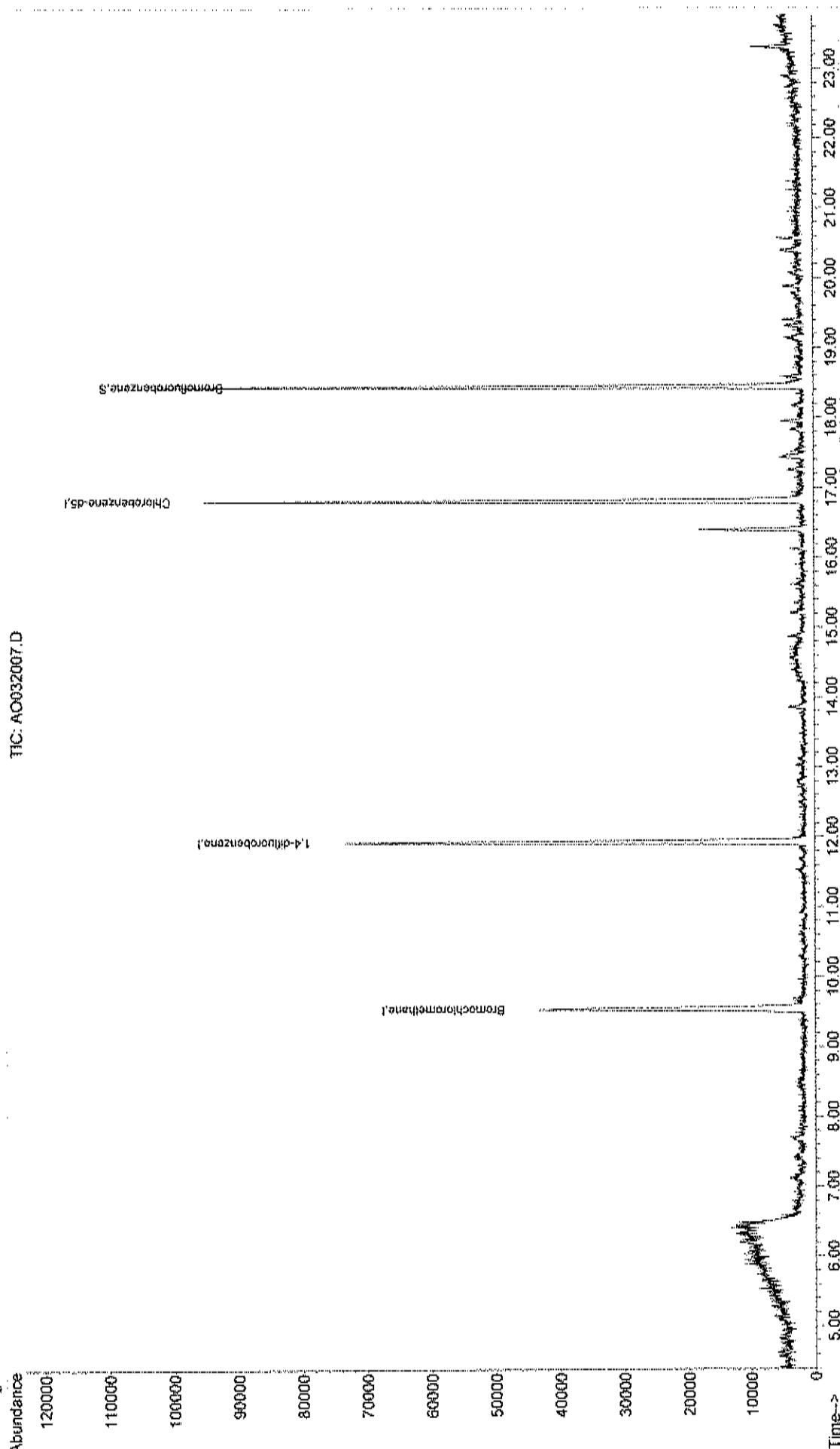
Qvalue

Data File : C:\HPCHEM\1\DATA2\AO032007.D  
 Acq On : 20 Mar 2017 3:21 pm  
 Sample : WAC032017B  
 Misc : A312\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Mar 21 9:03 2017

Vial: 2  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00  
 Quant Results File: A312\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A426\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 14:17:03 2017  
 Response via : Initial Calibration

HC: AO032007.D



Data File : C:\HPCHEM\1\DATA2\AO032008.D

Vial: 3

Acq On : 20 Mar 2017 4:02 pm

Operator: RJP

Sample : WAC032017C

Inst : MSD #1

Misc : A312\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 21 09:03:04 2017

Quant Results File: A312\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Wed Mar 15 10:58:20 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.52	128	14291	1.00	ppb	-0.02
35) 1,4-difluorobenzene	11.92	114	66450	1.00	ppb	-0.02
50) Chlorobenzene-d5	16.80	117	55920	1.00	ppb	-0.02

## System Monitoring Compounds

65) Bromofluorobenzene	18.44	95	38009	0.91	ppb	-0.01
Spiked Amount	1.000	Range	70 - 130	Recovery	=	91.00%

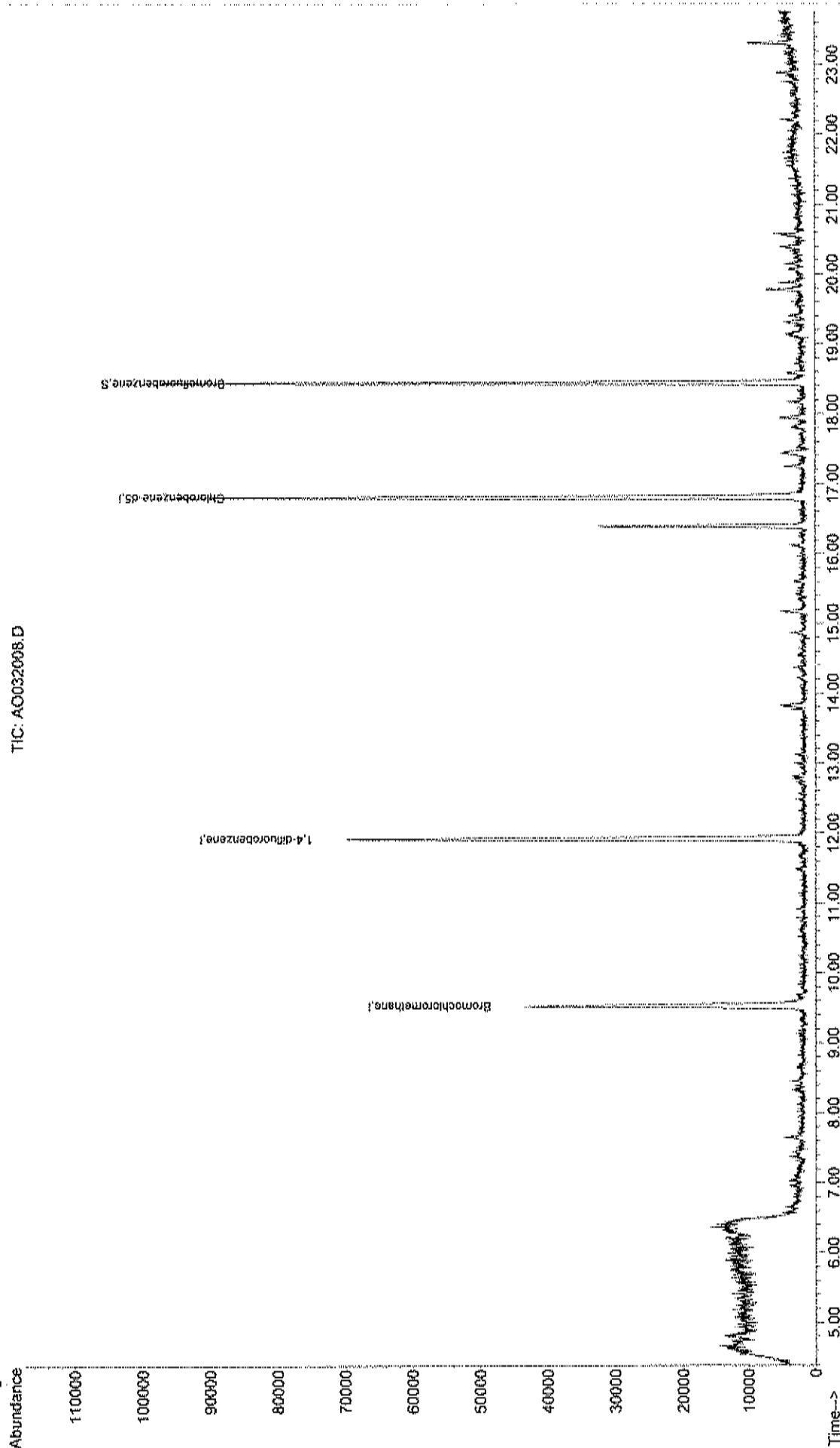
Target Compounds

Qvalue

Data File : C:\HPCHEM\1\DATA2\AO032008.D  
 Acq On : 20 Mar 2017 4:02 pm  
 Sample : WAC032017C  
 Misc : A312\_1UG  
 MS Integration Params: RPRINT.P  
 Quant Time: Mar 21 12:11 2017

Vial: 3  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00  
 Quant Results File: A312\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A426\_1UG.M (RTE Integration)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 14:17:03 2017  
 Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA2\AO032009.D

Vial: 4

Acq On : 20 Mar 2017 4:40 pm

Operator: RJP

Sample : WAC032017D

Inst : MSD #1

Misc : A312\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 21 09:03:05 2017

Quant Results File: A312\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Wed Mar 15 10:58:20 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	14630	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.92	114	66585	1.00	ppb	-0.02
50) Chlorobenzene-d5	16.81	117	54294	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.44	95	37248	0.92	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	92.00%

Target Compounds

Qvalue

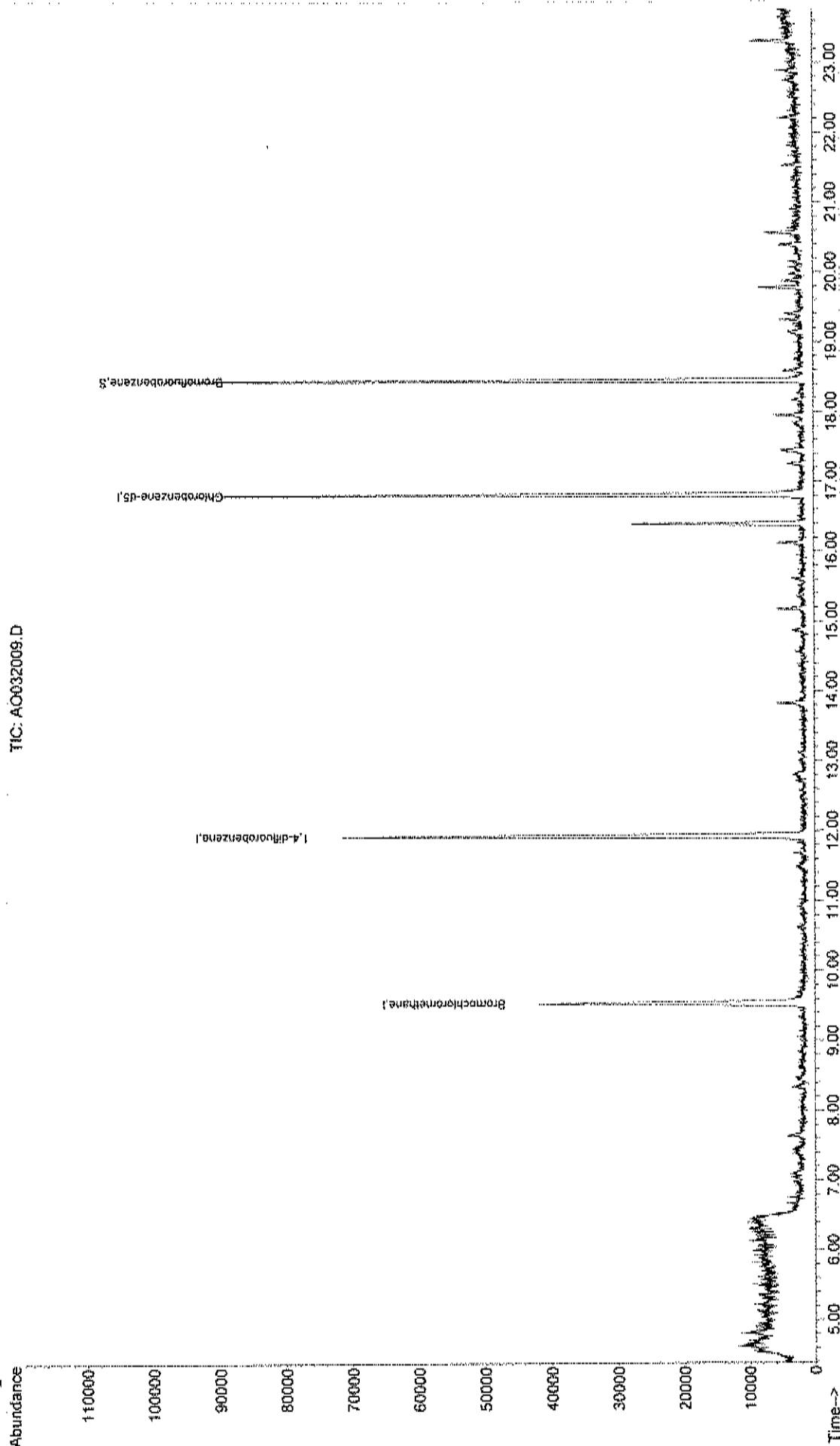


Data File : C:\HPCHEM\1\DATA2\AO032009.D  
Acq On : 20 Mar 2017 4:40 pm  
Sample : WAC032017D  
Misc : A312\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 21 9:03 2017

Vial: 4  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A312\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A426\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Thu May 04 14:17:03 2017  
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA2\AO032010.D

Vial: 5

Acq On : 20 Mar 2017 5:17 pm

Operator: RJP

Sample : WAC032017E

Inst : MSD #1

Misc : A312\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 21 09:03:06 2017

Quant Results File: A312\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Wed Mar 15 10:58:20 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	13448	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.92	114	65306	1.00	ppb	-0.02
50) Chlorobenzene-d5	16.81	117	53539	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.44	95	36660	0.92	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	92.00%

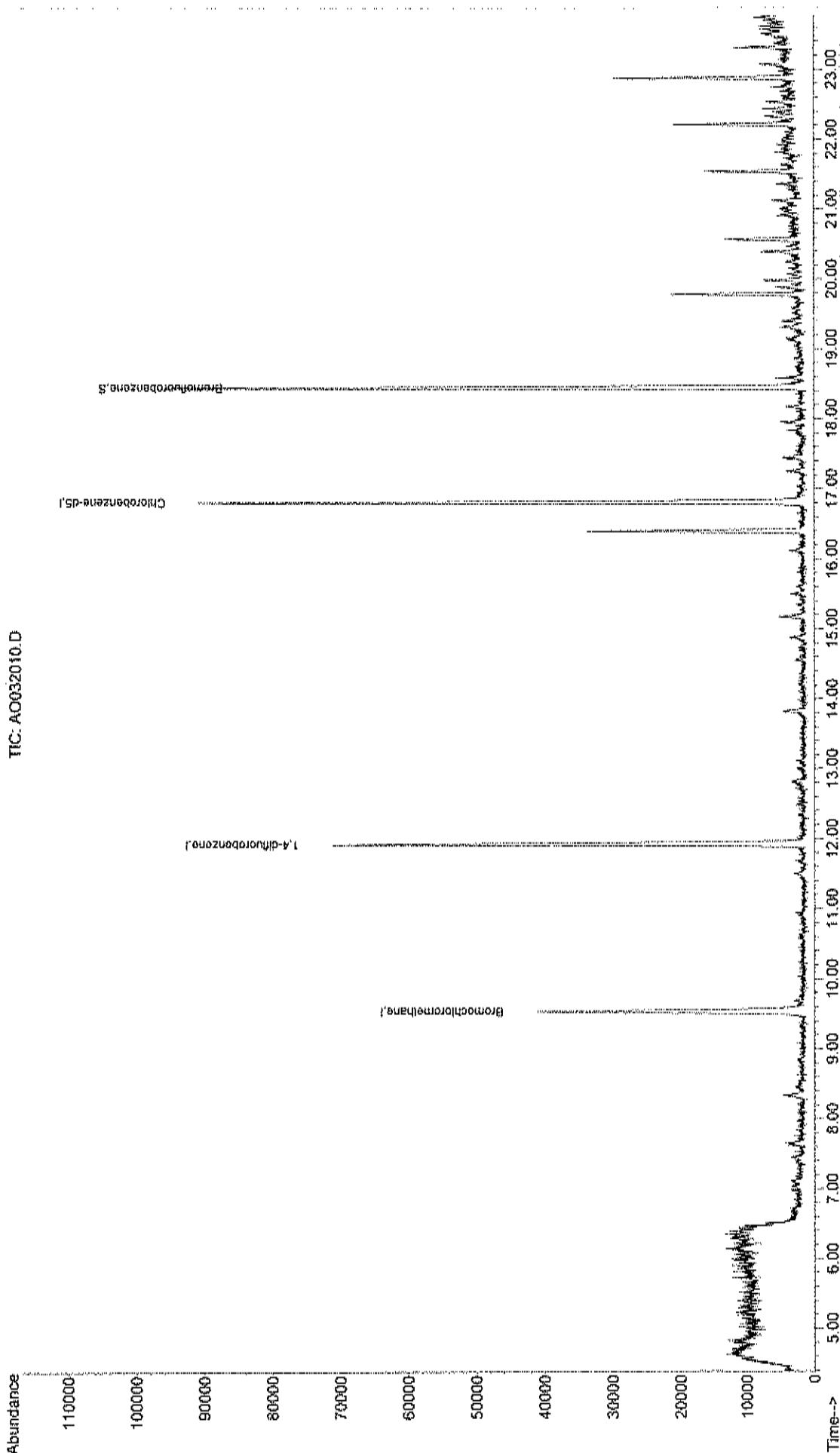
Target Compounds

Qvalue

Data File : C:\HPCHEM\1\DATA2\AO032010.D  
 Acq On : 20 Mar 2017 5:17 pm  
 Sample : WAC032017E  
 Misc : A312\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Mar 21 12:11 2017

Vial: 5  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00  
 Quant Results File: A312\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A426\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 14:17:03 2017  
 Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA2\AO032011.D

Vial: 6

Acq On : 20 Mar 2017 5:55 pm

Operator: RJP

Sample : WAC032017F

Inst : MSD #1

Misc : A312\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 21 09:03:07 2017

Quant Results File: A312\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Wed Mar 15 10:58:20 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.53	128	14055	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.92	114	62760	1.00	ppb	-0.01
50) Chlorobenzene-d5	16.81	117	51932	1.00	ppb	0.00

## System Monitoring Compounds

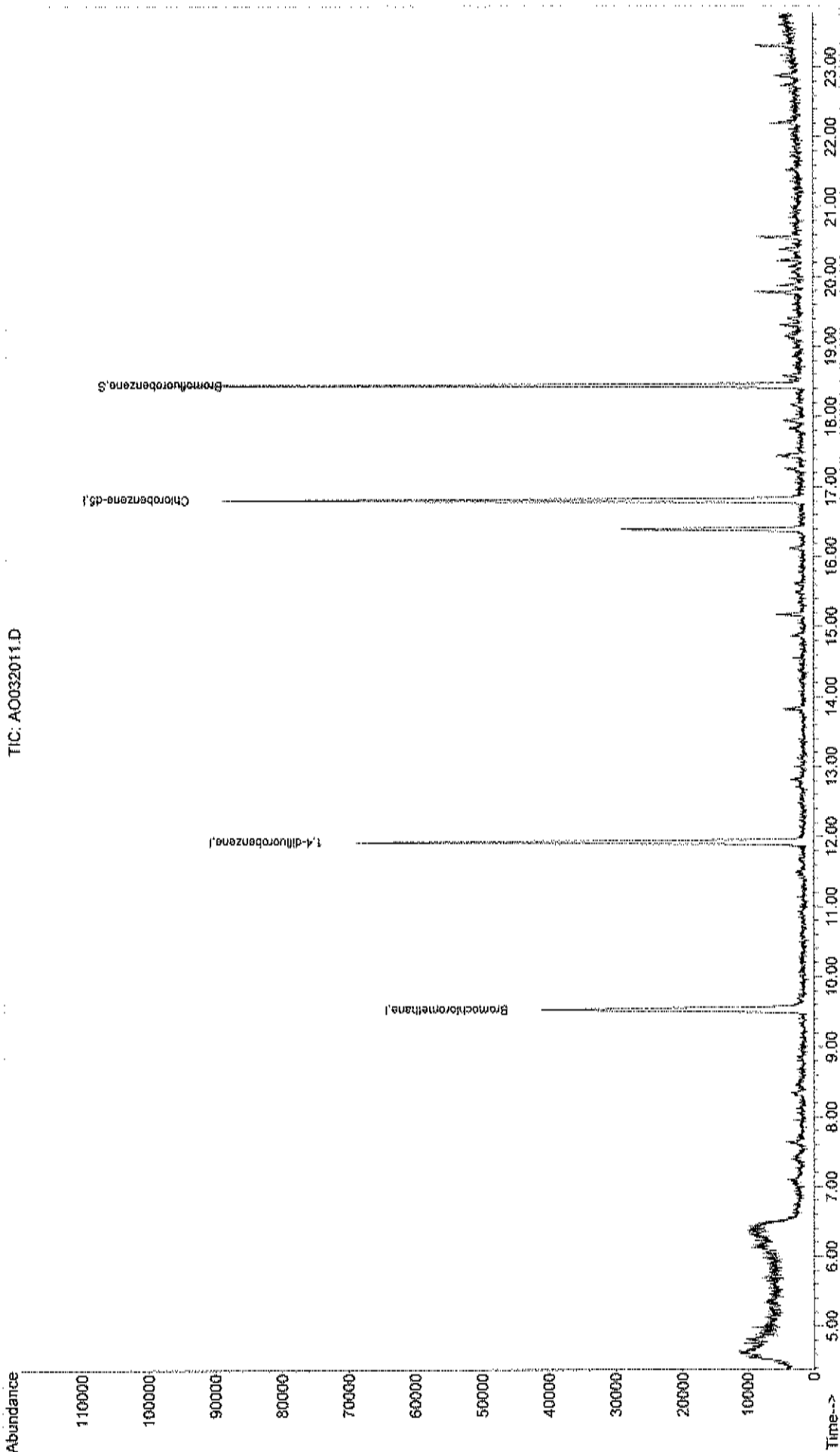
65) Bromofluorobenzene	18.44	95	35901	0.92	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	92.00%

Target Compounds

Qvalue

Data File : C:\HPCHEM\1\DATA2\AO032011.D  
 Acq On : 20 Mar 2017 5:55 pm  
 Sample : WAC032017F  
 Misc : A312\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Mar 21 12:11 2017  
 Quant Results File: A312\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A426\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 14:17:03 2017  
 Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA2\AO032012.D

Vial: 7

Acq On : 20 Mar 2017 6:32 pm

Operator: RJP

Sample : WAC032017G

Inst : MSD #1

Misc : A312\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 21 09:03:08 2017

Quant Results File: A312\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Wed Mar 15 10:58:20 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.54	128	13214	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.93	114	61386	1.00	ppb	0.00
50) Chlorobenzene-d5	16.81	117	51318	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.45	95	34657	0.90	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	90.00%

Target Compounds

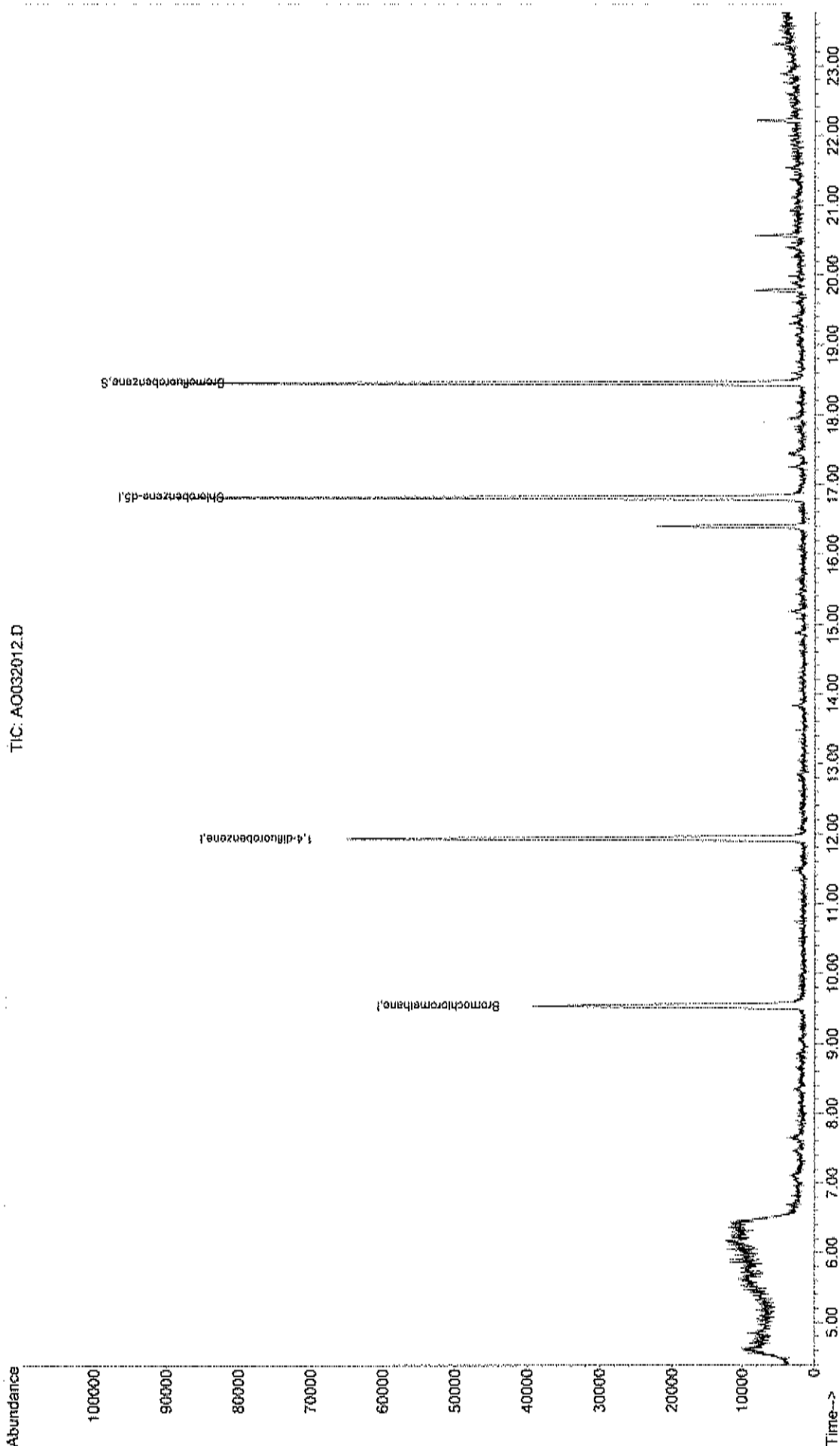
Qvalue

Data File : C:\HPCHEM\1\DATA2\AO032012.D  
 Acq On : 20 Mar 2017 6:32 pm  
 Sample : WAC032017G  
 Misc : A312\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Mar 21 9:03 2017  
 Quant Results File: A312\_1UG.RES

Vial: 7  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\A426\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 14:17:03 2017  
 Response via : Initial Calibration

TIC: AO032012.D



Data File : C:\HPCHEM\1\DATA2\AO032013.D  
 Acq On : 20 Mar 2017 7:10 pm  
 Sample : WAC032017H  
 Misc : A312\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Mar 21 09:03:09 2017

Vial: 8  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A312\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Wed Mar 15 10:58:20 2017  
 Response via : Initial Calibration  
 DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.53	128	13439	1.00	ppb	-0.01
35) 1,4-difluorobenzene	11.92	114	62774	1.00	ppb	-0.01
50) Chlorobenzene-d5	16.80	117	50944	1.00	ppb	-0.01

#### System Monitoring Compounds

65) Bromofluorobenzene	18.45	95	35665	0.94	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	94.00%

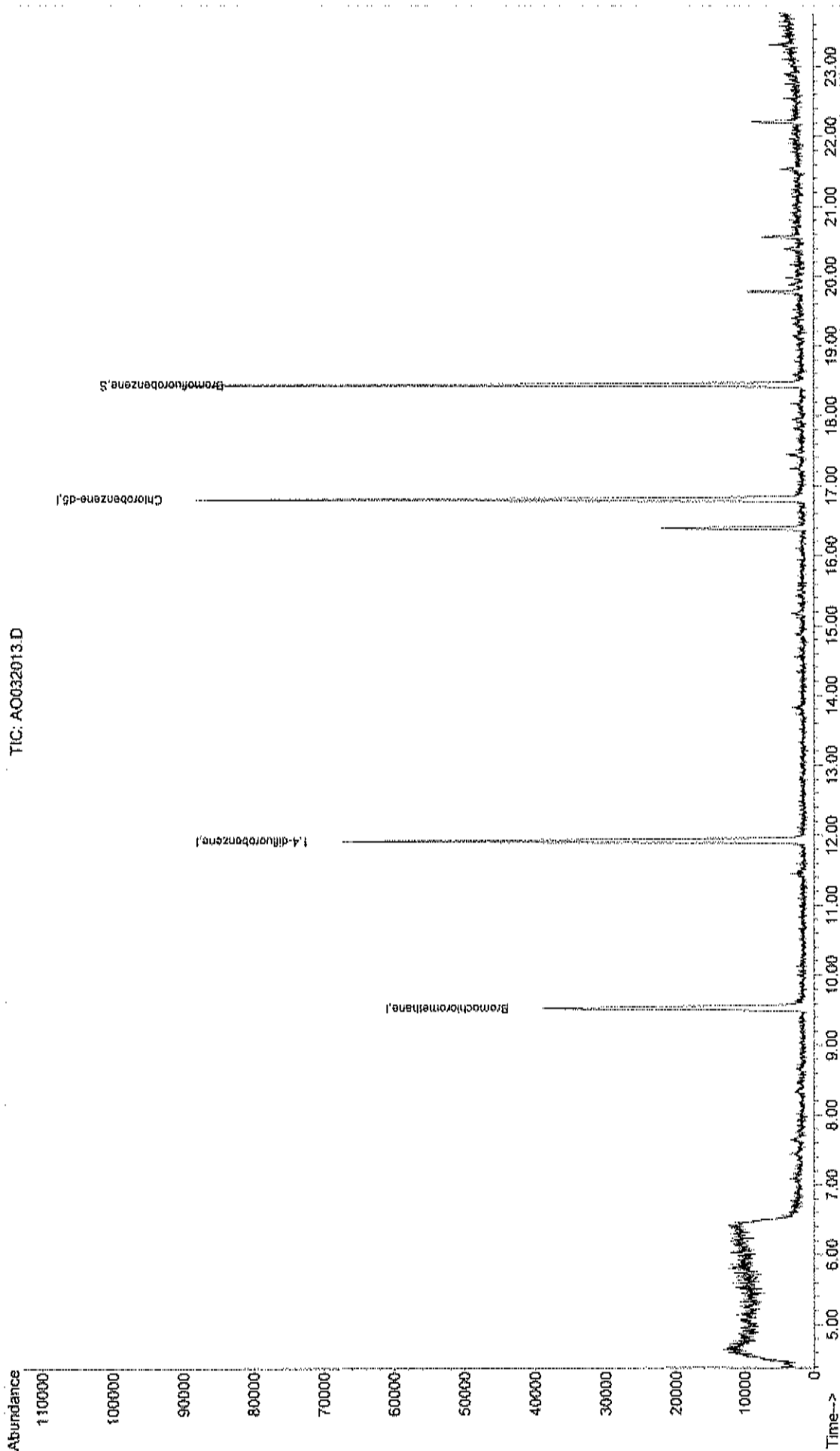
Target Compounds

Qvalue



Data File : C:\HPCHEM\1\DATA2\AO032013.D  
 Acq On : 20 Mar 2017 7:10 pm  
 Sample : WAC032017H  
 Misc : A312 1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Mar 21 9:03 2017  
 Quant Results File: A312 1UG.RES

Method : C:\HPCHEM\1\METHODS\A426\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Thu May 04 14:17:03 2017  
 Response via : Initial Calibration



## TO-15 Package Review Checklist

Client: LaBellaProject: 691-705 St. PaulSDG: C17R3015

		YES	NO	NA
Analytical Results	Present and Complete	/		
TIC's present	Present and Complete	/		
	Holding Times Met	/		

Comments: \_\_\_\_\_

\_\_\_\_\_

Chain-of-Custody	Present and Complete	/		
Surrogate Recovery	Present and Complete	/		
	Recoveries within limits	/		
	Sample(s) reanalyzed	/		
Internal Standards Recovery	Present and Complete	/		
	Recoveries within limits	/		
	Sample(s) reanalyzed	/		

Comments: \*SEE CASE NARRATIVE

\_\_\_\_\_

\_\_\_\_\_

Lab Control Sample (LCS)	Present and Complete	/		
	Recoveries within limits	/		
Lab Control Sample Dupe (LCSD)	Present and Complete	/		
	Recoveries within limits	/		
MS/MSD	Present and Complete	/		
	Recoveries within limits	/		

Comments: \_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

Sample Raw Data	Present and Complete	/		
	Spectra present for all samples	/		

Comments: \_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

## TO-15 Package Review Checklist

Client: LaBella Project: 691-705 St. Paul SDG: C1703015

		YES	NO	NA
Standards Data				
Initial Calibration Summary	Present and Complete	/		
	Calibration(s) met criteria	/		
Continuing Calibration Summary	Present and Complete	/		
	Calibration(s) met criteria	/		
Standards Raw Data	Present and Complete	/		

Comments: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

## Raw Quality Control Data

Tune Criteria Report	Present and Complete	/		
Method Blank Data	MB Results <PQL	/		
	Associated results flagged "B"	/		
LCS sample data	Present and Complete	/		
LCSD sample data	Present and Complete	/		
MS/MSD sample data	Present and Complete	/		

Comments: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

## Logbooks

Injection Log	Present and Complete	/		
Standards Log	Present and Complete	/		
Can Cleaning Log	Present and Complete	/		
	Raw Data Present	/		
Calculation sheet	Present and Complete	/		
IDL's	Present and Complete	/		
Bottle Order Form	Present and Complete	/		
Sample Tracking Form	Present and Complete	/		

Additional Comments: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

Section Supervisor: Will Doherty Date: 3/29/17

QC Supervisor: Will Doherty Date: 3/29/17



## CEN TEK LABORATORIES, LLC

143 Midler Park Drive \* Syracuse, NY 13206

Phone (315) 431-9730 \* Emergency 24/7 (315) 416-2752

NYSDOH ELAP

Certificate No. 11830

### **Analytical Report**

Michael Pelychaty  
LaBella Associates, P.C.  
300 State Street, Suite 201  
Rochester, NY 14614

Friday, March 10, 2017  
Order No.: C1703015

TEL: (585) 454-6110

FAX (585) 454-3066

RE: 691 and 705 St Paul St

Dear Michael Pelychaty:

Centek Laboratories, LLC received 6 sample(s) on 3/7/2017 for the analyses presented in the following report.

I certify that this data package is in compliance with the terms and conditions of the Contract, both technically and for completeness. Release of the data contained in this hardcopy data package and/or in the computer readable data submitted has been authorized by the Laboratory Manager or his designee, as verified by the following signature.

All method blanks, laboratory spikes, and/or matrix spikes met quality assurance objective except as indicated in the case narrative. All samples were received and analyzed within the EPA recommended holding times. Test results are not Method Blank (MB) corrected for contamination.

Centek Laboratories is distinctively qualified to meet your needs for precise and timely volatile organic compound analysis. We perform all analyses according to EPA, NIOSH or OSHA-approved analytical methods. Centek Laboratories is dedicated to providing quality analyses and exceptional customer service. Samples were analyzed using the methods outlined in the following references:

Compendium of Methods for the Determination of Toxic Organic Compounds, Compendium Method TO-15, January 1999.

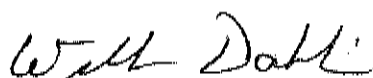
Centek Laboratories SOP TS-80

Analytical results relate to samples as received at laboratory. We do our best to make our reporting format clear and understandable and hope you are thoroughly satisfied with our services.

Please contact your client service representative at (315) 431-9730 or myself, if you would like any additional information regarding this report.

This report cannot be reproduced except in its entirety, without prior written authorization.

Sincerely,



William Dobbin  
Lead Technical Director

~~Disclaimer:~~ The test results and procedures utilized, and laboratory interpretations of the data obtained by Centek as contained in this report are believed by Centek to be accurate and reliable for sample(s) tested. In accepting this report, the customer agrees that the full extent of any and all liability for actual and consequential damages of Centek for the services performed shall be equal to the fee charged to the customer for the services as liquidated damages. ELAP does not offer certification for the following parameters by this method at present time, they are: 4-ethyltoluene, ethyl acetate, propylene, tetrahydrofuran, 4-PCH, sulfur derived and silicon series compounds.

#### Centek Laboratories, LLC Terms and Conditions

##### Sample Submission

All samples sent to Centek Laboratories should be accompanied by our Request for Analysis Form or Chain of Custody Form. A Chain of Custody will be provided with each order shipped for all sampling events, or if needed, one is available at our website [www.CentekLabs.com](http://www.CentekLabs.com). Samples received after 3:00pm are considered to be a part of the next day's business.

##### Sample Media

Samples can be collected in an canister or a Tedlar bag. Depending on your analytical needs, Centek Laboratories may receive a bulk, liquid, soil or other matrix sample for headspace analysis.

##### Blanks

Every sample is run with a surrogate or tracer compound at a pre-established concentration. The surrogate compound run with each sample is used as a standard to measure the performance of each run of the instrument. If required, a Minican can be provided containing nitrogen to be run as a trip blank with your samples.

##### Sampling Equipment

Centek Laboratories will be happy to provide the canisters to carry-out your sampling event at no charge. The necessary accessories, such as regulators, tubing or personal sampling belts, are also provided to meet your sampling needs. The customer is responsible for all shipping charges to the client's destination and return shipping to the laboratory. Client assumes all responsibility for lost, stolen and any damages of equipment.

##### Turn Around time (TAT)

Centek Laboratories will provide results to its clients in one business-week by 6:00pm EST after receipt of samples. For example, if samples are received on a Monday they are due on the following Monday by 6:00pm EST. Results are faxed or emailed to the requested location indicated on the Chain of Custody. Non-routine analysis may require more than the one business-week turnaround time. Please confirm non-routine sample turnaround times.

#### Reporting

Results are emailed or faxed at no additional charge. A hard copy of the result report is mailed within 24 hours of the faxing or emailing of your results. Cat "B" like packages are within 3-4 weeks from time of analysis. Standard Electronic Disk Deliverables (EDD) is also available at no additional charge.

#### Payment Terms

Payment for all purchases shall be due within 30 days from date of invoice. The client agrees to pay a finance charge of 1.5% per month on the overdue balance and cost of collection, including attorney fees, if collection proceedings are necessary. You must have a completed credit application on file to extend credit. Purchase orders or checks information must be submitted for us to release results

#### Rush Turnaround Samples

Expedited turn around times is available. Please confirm rush turnaround times with Client Services before submitting samples.

Applicable Surcharges for Rush Turnaround Samples:

Same day TAT = 200%

Next business day TAT by Noon = 150%

Next business day TAT by 6:00pm = 100%

Second business day TAT by 6:00pm = 75%

Third business day TAT by 6:00pm = 50%

Fourth business day TAT by 6:00pm = 35%

Fifth business day = Standard

#### Statement of Confidentiality

Centek Laboratories, LLC is aware of the importance of the confidentiality of results to many of our clients. Your name and data will be held in the strictest of confidence. We will not accept business that may constitute a conflict of interest. We commonly sign Confidential Nondisclosure Agreements with clients prior to beginning work. All research, results and reports will be kept strictly confidential. Secrecy Agreements and Disclosure Statements will be signed for the client if so specified. Results will be provided only to the addressee specified on the Chain of Custody Form submitted with the samples unless law requires release. Written permission is required from the addressee to release results to any other party.

#### Limitation on Liability

Centek Laboratories, LLC warrants the test results to be accurate to the methodology and sample type for each sample submitted to Centek Laboratories, LLC. In no event shall Centek Laboratories, LLC be liable for direct, indirect, special, punitive, incidental, exemplary or consequential damages, or any damages whatsoever, even if Centek Laboratories, LLC has been previously advised of the possibility of such damages whether in an action under contract, negligence, or any other theory, arising out of or in connection with the use, inability to use or performance of the information, services, products and materials available from the laboratory or this site. These limitations shall apply notwithstanding any failure of essential purpose of any limited remedy. Because some jurisdictions do not allow limitations on how long an implied warranty lasts, or the exclusion or limitation of liability for consequential or incidental damages, the above limitations may not apply to you. This is a comprehensive limitation of

liability that applies to all damages of any kind, including (without limitation) compensatory, direct, indirect or consequential damages, loss of data, income or profit and or loss of or damage to property and claims of third parties.

## **ASP CAT B DELIVERABLE PACKAGE Table of Contents**

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  - b. IS Summary Report**
  - c. MB Summary Report**
  - d. LCS Summary Report**
  - e. MSD Summary Report**
  - f. IDL's**
  - g. Calculation**
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  - b. Quantitation Report with Spectra**
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**CEN TEK LABORATORIES, LLC**

**Date:** 29-Mar-17

**CLIENT:** LaBella Associates, P.C.

**Project:** 691 and 705 St Paul St

**Lab Order:** C1703015

## **CASE NARRATIVE**

---

Samples were analyzed using the methods outlined in the following references:

Centek Laboratories, LLC SOP TS-80

Compendium of Methods for the Determination of Toxic Organic Compounds, Compendium Method TO-15, January 1999

All method blanks, laboratory spikes, and/or matrix spikes met quality assurance objective except as indicated in the corrective action report(s). All samples were received and analyzed within the EPA recommended holding times. Test results are not Method Blank (MB) corrected for contamination.

### **NYSDEC ASP samples:**

Canisters should be evacuated to a reading of less than or equal to 50 millitorr prior to shipment to sampling personnel. The vacuum in the canister will be field checked prior to sampling, and must read 28" of Hg ( $\pm 2$ ", vacuum, absolute) before a sample can be collected. After the sample has been collected, the pressure of the canister will be read and recorded again, and must be 5" of Hg ( $\pm 1$ ", vacuum, absolute) for the sample to be valid. Once received at the laboratory, the canister vacuum should be confirmed to be 5" of Hg,  $\pm 1$ ". Please record and report the pressure/vacuum of received canisters on the sample receipt paperwork. A pressure/vacuum reading should also be taken just prior to the withdrawal of sample from the canister, and recorded on the sample preparation log sheet. All regulators are calibrated to meet these requirements before they leave the laboratory. However, due to environmental conditions and use of the equipment Centek can not guarantee that this criteria can always be achieved.

See Corrective Action: [3488] The 3rd IS & Surrogate did not meet criteria.

## Centek Laboratories, LLC

## Corrective Action Report

Date Initiated: 08-Mar-17

Corrective Action Report ID: 3488

Initiated By: Russell Pellegrino

Department: MSVOA

## Corrective Action Description

**CAR Summary:** The 3rd IS & Surrogate did not meet criteria.**Description of Nonconformance Root/Cause(s):** The 3rd IS & Surrogate were high and did not meet criteria for samples C1703015-001A - 005A & 003A MS/MSD. Based on the chromatographic evidence, it appears that the contamination is from the presence of very high petroleum pattern. All Compounds of concern fall under the first 2 Internal standards which did meet criteria**Description of Corrective Action w/Proposed C.A.:** Samples C1703015-003A - 005A were analyzed further as dilutions with criteria being met. Due to matrix being in a canister it is difficult to see any signs of problems. All sets of data submitted.**Performed By:** Russell Pellegrino**Completion Date:** 09-Mar-17

## Client Notification

**Client Notification Required:** No**Notified By:****Comment:**

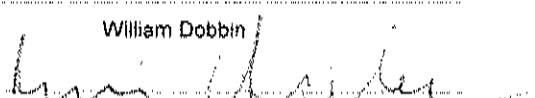
## Quality Assurance Review

**Nonconformance Type:** Deficiency**Further Action required by QA:** Since the compounds of concern fall under the first 2 internal standards and they did meet criteria and all other QC meets requirements report the 1x results. Submit all sets of data.

## Approval and Closure

**Technical Director /  
Deputy Tech. Dir.:****Close Date:** 10-Mar-17

William Dobbin

**QA Officer Approval:**

Nick Scala

**QA Date:** 10-Mar-17**Last Updated BY** russ**Updated:** 29-Mar-2017 9:25 AM**Reported:** 29-Mar-2017 9:25 AM





CENTEK LABORATORIES, LLC

Date: 29-Mar-17

CLIENT: LaBella Associates, P.C.

Project: 691 and 705 St Paul St

Lab Order: C1703015

## Work Order Sample Summary

Lab Sample ID	Client Sample ID	Tag Number	Collection Date	Date Received
C1703015-001A	SV-1	542.256	3/3/2017	3/7/2017
C1703015-002A	SV-2	237.402	3/3/2017	3/7/2017
C1703015-003A	SV-3	1206.249	3/3/2017	3/7/2017
C1703015-004A	SV-4	290.299	3/3/2017	3/7/2017
C1703015-005A	Duplicate	1184.299	3/3/2017	3/7/2017
C1703015-006A	Ambient Air	556.267	3/3/2017	3/7/2017



## CENTEK LABORATORIES, LLC

## Sample Receipt Checklist

Client Name LABELLA - ROCHESTER

Date and Time Receive

3/7/2017

Work Order Number C1703015

Received by NM

Checklist completed by

Signature

Date

Reviewed by

Initials

Date

Matrix:

Carrier name: FedEx Ground

Shipping container/cooler in good condition?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	Not Present <input type="checkbox"/>
Custody seals intact on shipping container/cooler?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	Not Present <input checked="" type="checkbox"/>
Custody seals intact on sample bottles?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	Not Present <input checked="" type="checkbox"/>
Chain of custody present?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Chain of custody signed when relinquished and received?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Chain of custody agrees with sample labels?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Samples in proper container/bottle?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Sample containers intact?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Sufficient sample volume for indicated test?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
All samples received within holding time?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Container/Temp Blank temperature in compliance?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Water - VOA vials have zero headspace?	No VOA vials submitted <input checked="" type="checkbox"/>	Yes <input type="checkbox"/>	No <input type="checkbox"/>
Water - pH acceptable upon receipt?	Yes <input type="checkbox"/>	No <input checked="" type="checkbox"/>	

Adjusted?

Checked by

Any No and/or NA (not applicable) response must be detailed in the comments section be

Client contacted \_\_\_\_\_ Date contacted: \_\_\_\_\_ Person contacted \_\_\_\_\_

Contacted by: \_\_\_\_\_ Regarding: \_\_\_\_\_

Comments: \_\_\_\_\_

Corrective Action \_\_\_\_\_

## Centek Laboratories, LLC

29-Mar-17

Lab Order: C1703015  
 Client: LaBella Associates, P.C.  
 Project: 691 and 705 St Paul St

## DATES REPORT

Sample ID	Client Sample ID	Collection Date	Matrix	Test Name	TCLP Date	Prep Date	Analysis Date
C1703015-001A	SV-1	3/3/2017	Air	1ug/M3 by Method TO15			3/8/2017
C1703015-002A	SV-2			1ug/M3 by Method TO15			3/8/2017
C1703015-003A	SV-3			1ug/M3 by Method TO15			3/8/2017
				1ug/M3 by Method TO15			3/8/2017
C1703015-004A	SV-4			1ug/M3 by Method TO15			3/9/2017
				1ug/M3 by Method TO15			3/8/2017
C1703015-005A	Duplicate			1ug/M3 by Method TO15			3/9/2017
				1ug/M3 by Method TO15			3/8/2017
C1703015-006A	Ambient Air			1ug/m3 w/ 0.25ug/M3 CT-TCE-VC			3/8/2017

**CENTEK LABORATORIES, LLC***Air Quality Testing...It's a Gas*

143 Midler Park Drive \* Syracuse, NY 13206

TEL: 315-431-9730 \* FAX: 315-431-9731

**CANISTER ORDER****6352**

29-Mar-17

**SHIPPED TO:**

Company: LaBella Associates, P.C.

Contact: Michael Pelychaty

Address: 300 State Street, Suite 201

Rochester, NY 14614

Phone: (585) 454-6110

Quote ID: 0

Project:

PO:

Submitted By:

MadeBy: NM

Ship Date: 2/28/2017

VIA: FedEx Ground

Due Date: 3/1/2017

Bottle Code	Bottle Type	TEST(s)	QTY
MC1400CC	1.4L Mini-Can	1ug/M3 by Method TO15	1
MC1000CC	1L Mini-Can	1ug/M3 by Method TO15	6

Can / Reg ID	Description
237	1L Mini-Can - 1168 VI
249	Time-Set Reg - 687 VI
256	Time-Set Reg - 694 VI
267	Time-Set Reg - 705 VI
290	1L Mini-Can - 1266 VI
339	Time-Set Reg - 736 VI
340	Time-Set Reg - 737 VI
402	Time-Set Reg - 781 VI
459	1L Mini-Can - 1362 VI
556	1L Mini-Can - 124 VI
1184	1L Mini-Can - 1248 VI
1206	1.4L Mini-Can - 1376 VI

Comments: (6) 1L @ 8hrs + (1) 1.4L @ 8hrs + dupe W/ He shroud &amp; Meter, tubing WAC 110316 A-B, 021717A-C

**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**ANALYTICAL RESULTS**



## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-001A

Client Sample ID: SV-1  
 Tag Number: 542.256  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-1			"Hg		3/7/2017
Lab Vacuum Out	-30			"Hg		3/7/2017
<b>1UG/M3 BY METHOD TO15</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 1:28:00 PM
Chloroethane	0.13	0.15	J	ppbV	1	3/8/2017 1:28:00 PM
cis-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 1:28:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 1:28:00 PM
Trichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 1:28:00 PM
Vinyl chloride	< 0.15	0.15		ppbV	1	3/8/2017 1:28:00 PM
Surr; Bromofluorobenzene	239	70-130	S	%REC	1	3/8/2017 1:28:00 PM

**NOTES:**

Surrogate did not meet criteria due to severe matrix interference.

<b>Qualifiers:</b>	<b>**</b>	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte, Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

**Centek Laboratories, LLC**

Date: 27-Mar-17

**CLIENT:** LaBella Associates, P.C.  
**Lab Order:** C1703015  
**Project:** 691 and 705 St Paul St  
**Lab ID:** C1703015-001A

**Client Sample ID:** SV-1  
**Tag Number:** 542.256  
**Collection Date:** 3/3/2017  
**Matrix:** AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>1UG/M3 BY METHOD TO15</b>		<b>TO-15</b>		<b>Analyst: RJP</b>		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 1:28:00 PM
Chloroethane	0.34	0.40	J	ug/m3	1	3/8/2017 1:28:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 1:28:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 1:28:00 PM
Trichloroethene	< 0.81	0.81		ug/m3	1	3/8/2017 1:28:00 PM
Vinyl chloride	< 0.38	0.38		ug/m3	1	3/8/2017 1:28:00 PM

**NOTES:**

Surrogate did not meet criteria due to severe matrix interference.

<b>Qualifiers:</b>	<b>**</b>	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-002A

Client Sample ID: SV-2  
 Tag Number: 237.402  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-1			"Hg		3/7/2017
Lab Vacuum Out	-30			"Hg		3/7/2017
<b>1UG/M3 BY METHOD TO15</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 2:09:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	3/8/2017 2:09:00 PM
cis-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 2:09:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 2:09:00 PM
Trichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 2:09:00 PM
Vinyl chloride	< 0.15	0.15		ppbV	1	3/8/2017 2:09:00 PM
Surr: Bromofluorobenzene	235	70-130	S	%REC	1	3/8/2017 2:09:00 PM
<b>NOTES:</b>						
Surrogate did not meet criteria due to severe matrix interference.						

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits  
 . Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

Page 2 of 6

## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-002A

Client Sample ID: SV-2  
 Tag Number: 237.402  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY METHOD TO15		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 2:09:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	3/8/2017 2:09:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 2:09:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 2:09:00 PM
Trichloroethene	< 0.81	0.81		ug/m3	1	3/8/2017 2:09:00 PM
Vinyl chloride	< 0.38	0.38		ug/m3	1	3/8/2017 2:09:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-003A

Client Sample ID: SV-3  
 Tag Number: 1206.249  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	1			"Hg		3/7/2017
Lab Vacuum Out	-30			"Hg		3/7/2017
<b>1UG/M3 BY METHOD TO15</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 2:48:00 PM
Chloroethane	0.11	0.15	J	ppbV	1	3/8/2017 2:48:00 PM
cis-1,2-Dichloroethene	0.18	0.15		ppbV	1	3/8/2017 2:48:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 2:48:00 PM
Trichloroethene	7.4	1.5		ppbV	10	3/8/2017 11:34:00 PM
Vinyl chloride	< 0.15	0.15		ppbV	1	3/8/2017 2:48:00 PM
Surr: Bromofluorobenzene	97.0	70-130		%REC	10	3/8/2017 11:34:00 PM
Surr: Bromofluorobenzene	166	70-130	S	%REC	1	3/8/2017 2:48:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

Qualifiers:	**	Quantitation Limit	,	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte, Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-003A

Client Sample ID: SV-3  
 Tag Number: 1206.249  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY METHOD TO15		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 2:48:00 PM
Chloroethane	0.29	0.40	J	ug/m3	1	3/8/2017 2:48:00 PM
cis-1,2-Dichloroethene	0.71	0.59		ug/m3	1	3/8/2017 2:48:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 2:48:00 PM
Trichloroethene	40	8.1		ug/m3	10	3/8/2017 11:34:00 PM
Vinyl chloride	< 0.38	0.38		ug/m3	1	3/8/2017 2:48:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

Qualifiers:	** Quantitation Limit	.	Results reported are not blank corrected
	B Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S Spike Recovery outside accepted recovery limits		

## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-004A

Client Sample ID: SV-4  
 Tag Number: 290.299  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
				FLD		Analyst:
Lab Vacuum In	-3			"Hg		3/7/2017
Lab Vacuum Out	-30			"Hg		3/7/2017
<b>1UG/M3 BY METHOD TO15</b>						
				TO-15		Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 4:59:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	3/8/2017 4:59:00 PM
cis-1,2-Dichloroethene	1.0	0.15		ppbV	1	3/8/2017 4:59:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 4:59:00 PM
Trichloroethene	8.9	1.5		ppbV	10	3/9/2017 12:47:00 AM
Vinyl chloride	< 0.15	0.15		ppbV	1	3/8/2017 4:59:00 PM
Surr: Bromofluorobenzene	100	70-130		%REC	10	3/9/2017 12:47:00 AM
Surr: Bromofluorobenzene	190	70-130	S	%REC	1	3/8/2017 4:59:00 PM
<b>NOTES:</b>						
Surrogate did not meet criteria due to severe matrix interference.						

Qualifiers:	**	Quantitation Limit	,	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-004A

Client Sample ID: SV-4  
 Tag Number: 290.299  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY METHOD TO15		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 4:59:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	3/8/2017 4:59:00 PM
cis-1,2-Dichloroethene	4.1	0.59		ug/m3	1	3/8/2017 4:59:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 4:59:00 PM
Trichloroethene	48	8.1		ug/m3	10	3/9/2017 12:47:00 AM
Vinyl chloride	< 0.38	0.38		ug/m3	1	3/8/2017 4:59:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		



## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-005A

Client Sample ID: Duplicate  
 Tag Number: 1184.299  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-3			"Hg		3/7/2017
Lab Vacuum Out	-30			"Hg		3/7/2017
<b>1UG/M3 BY METHOD TO15</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 5:39:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	3/8/2017 5:39:00 PM
cis-1,2-Dichloroethene	0.99	0.15		ppbV	1	3/8/2017 5:39:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 5:39:00 PM
Trichloroethene	8.8	1.5		ppbV	10	3/9/2017 2:01:00 AM
Vinyl chloride	< 0.15	0.15		ppbV	1	3/8/2017 5:39:00 PM
Surr: Bromofluorobenzene	98.0	70-130		%REC	10	3/9/2017 2:01:00 AM
Surr: Bromofluorobenzene	187	70-130	S	%REC	1	3/8/2017 5:39:00 PM

**NOTES:**

Surrogate did not meet criteria due to severe matrix interference.

Qualifiers:	**	Quantitation Limit	,	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

**Centek Laboratories, LLC**

Date: 27-Mar-17

**CLIENT:** LaBella Associates, P.C.  
**Lab Order:** C1703015  
**Project:** 691 and 705 St Paul St  
**Lab ID:** C1703015-005A

**Client Sample ID:** Duplicate  
**Tag Number:** 1184.299  
**Collection Date:** 3/3/2017  
**Matrix:** AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>1UG/M3 BY METHOD TO15</b>		<b>TO-15</b>		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 5:39:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	3/8/2017 5:39:00 PM
cis-1,2-Dichloroethene	3.9	0.59		ug/m3	1	3/8/2017 5:39:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 5:39:00 PM
Trichloroethene	47	8.1		ug/m3	10	3/9/2017 2:01:00 AM
Vinyl chloride	< 0.38	0.38		ug/m3	1	3/8/2017 5:39:00 PM

**NOTES:**

Surrogate did not meet criteria due to severe matrix interference.

<b>Qualifiers:</b>	<b>** Quantitation Limit</b>	Results reported are not blank corrected
B	Analyte detected in the associated Method Blank	E Estimated Value above quantitation range
H	Holding times for preparation or analysis exceeded	J Analyte detected below quantitation limit
JN	Non-routine analyte. Quantitation estimated.	ND Not Detected at the Limit of Detection
S	Spike Recovery outside accepted recovery limits	

## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-006A

Client Sample ID: Ambient Air  
 Tag Number: 556.267  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses Result \*\*Limit Qual Units DF Date Analyzed

FIELD PARAMETERS		FLD		Analyst:	
Lab Vacuum In	1	"Hg		3/7/2017	
Lab Vacuum Out	-30	"Hg		3/7/2017	
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP	
1,1-Dichloroethene	< 0.15	0.15	ppbV	1	3/8/2017 12:48:00 PM
Chloroethane	< 0.15	0.15	ppbV	1	3/8/2017 12:48:00 PM
cis-1,2-Dichloroethene	< 0.15	0.15	ppbV	1	3/8/2017 12:48:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15	ppbV	1	3/8/2017 12:48:00 PM
Trichloroethene	< 0.040	0.040	ppbV	1	3/8/2017 12:48:00 PM
Vinyl chloride	< 0.040	0.040	ppbV	1	3/8/2017 12:48:00 PM
Surr: Bromofluorobenzene	87.0	70-130	%REC	1	3/8/2017 12:48:00 PM

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

**Centek Laboratories, LLC**

Date: 27-Mar-17

**CLIENT:** LaBella Associates, P.C.  
**Lab Order:** C1703015  
**Project:** 691 and 705 St Paul St  
**Lab ID:** C1703015-006A

**Client Sample ID:** Ambient Air  
**Tag Number:** 556.267  
**Collection Date:** 3/3/2017  
**Matrix:** AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>		<b>TO-15</b>		<b>Analyst: RJP</b>		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 12:48:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	3/8/2017 12:48:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 12:48:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 12:48:00 PM
Trichloroethene	< 0.21	0.21		ug/m3	1	3/8/2017 12:48:00 PM
Vinyl chloride	< 0.10	0.10		ug/m3	1	3/8/2017 12:48:00 PM

<b>Qualifiers:</b>	<b>** Quantitation Limit</b>	, Results reported are not blank corrected
B	Analyte detected in the associated Method Blank	E Estimated Value above quantitation range
H	Holding times for preparation or analysis exceeded	J Analyte detected below quantitation limit
JN	Non-routine analyte, Quantitation estimated.	ND Not Detected at the Limit of Detection
S	Spike Recovery outside accepted recovery limits	

**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**QUALITY CONTROL SUMMARY**

Date: 27-Mar-17



CENTEK LABORATORIES, LLC

# QC SUMMARY REPORT SURROGATE RECOVERIES

CLIENT: LaBella Associates, P.C.

Work Order: C1703015

Project: 691 and 705 St Paul St

Test No: TO-15

Matrix: A

Sample ID	BR4FBZ								
ALCSIUG-030817	105								
ALCSIUGD-030817	104								
AMBIUG-030817	85.0								
C1703015-001A	239 *								
C1703015-002A	235 *								
C1703015-003A	97.0								
C1703015-003A MS	170 *								
C1703015-003A MSD	173 *								
C1703015-004A	100								
C1703015-005A	98.0								
C1703015-006A	87.0								

Acronym	Surrogate	QC Limits
BR4FBZ	= Bromofluorobenzene	70-130

\* Surrogate recovery outside acceptance limits

1

## GC/MS QA-QC Check Report

Tune File : C:\HPCHEM\1\DATA\AO030803.D

Tune Time : 8 Mar 2017 10:33 am

Daily Calibration File : C:\HPCHEM\1\DATA\AO030803.D

		(BFB)	(IS1)	(IS2)	(IS3)
			51607	232876	196939
File	Sample	DL	Surrogate Recovery %	Internal Standard Responses	
AO030804.D	ALCS1UG-030817	105		54344	232438 196865
AO030805.D	AMB1UG-030817	85		51456	199472 169731
AO030806.D	C1703015-006A	87		47960	186316 173836
AO030807.D	C1703015-001A	239*		59338	270574 278336
AO030808.D	C1703015-002A	235*		72867	317032 316562*
AO030809.D	C1703015-003A	166*		80896	366728 336237*
AO030810.D	C1703015-003A MS	170*		80202	356830 344072*
AO030811.D	C1703015-003A MSD	173*		79391	370639 337498*
AO030812.D	C1703015-004A	190*		76299	343779 323940*
AO030813.D	C1703015-005A	187*		77699	348681 330439*
AO030822.D	C1703015-003A 10X	97		53827	227536 201273
AO030824.D	C1703015-004A 10X	100		46536	210199 197259
AO030826.D	C1703015-005A 10X	98		46525	198463 190799
AO030828.D	ALCS1UGD-030817	104		45106	199550 175594

t - fails 24hr time check \* - fails criteria

Created: Mon Mar 27 11:04:40 2017 MSD #1/

**CENTEK LABORATORIES, LLC**

## ANALYTICAL QC SUMMARY REPORT

**CLIENT:** LaBella Associates, P.C.

Work Order: C1703015

**Project:** 691 and 705 St Paul St

TestCode: 0.25CT-TCE-VC

Sample ID	ALCS1UG-030817	SampleType:	LCS	TestCode:	0.25CT-ICE-	Units:	ppbV	Prep Date:	RunNo:	12014		
Client ID:	ZZZZ	Batch ID:	R12014	TestNo:	TO-15			Analysis Date:	SeqNo:	140535		
Analyte		Result		PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	RPDLimit	Qual

1,1-Dichloroethene	1.060	0.15	1	0	106	70	130
Chloroethane	0.9000	0.15	1	0	90.0	70	130
cis-1,2-Dichloroethene	1.100	0.15	1	0	110	70	130
trans-1,2-Dichloroethene	1.070	0.15	1	0	107	70	130
Trichloroethene	1.060	0.040	1	0	106	70	130
Vinyl chloride	0.8700	0.040	1	0	87.0	70	130

Sample ID	ALCS1UGD-030817	SampleType:	LCSO	TestCode:	0.25CT-ICE-	Units:	ppbv	Prep Date:	RunNo:	12014
Client ID:	ZZZZZ	Batch ID:	R12014	TestNo:	TO-15	Analysis Date:	3/9/2017	SeqNo:	140536	
Analyte		Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	RPDLimit Qual

1,1-Dichloroethene	1	0	110	70	130	1.06	3.70	30
Chloroethane	1	0	100	70	130	0.9	10.5	30
cis-1,2-Dichloroethene	1	0	113	70	130	1.1	2.69	30
trans-1,2-Dichloroethene	1	0	110	70	130	1.07	2.76	30
Trichloroethene	1	0	110	70	130	1.06	3.70	30
Vinyl chloride	1	0	91.0	70	130	0.87	4.49	30

**Qualifiers:** Results reported are not blank corrected

† Analyte detected below quantitation limit

S Spike Recovery outside accepted recovery limits

Estimated Value above qualification range

AND Not Detected at the Limit of Detection

... papaya's skin but is discarded for some time before...

RPD outside accepted recovery limits



Date: 27-Mar-17

## CENTEK LABORATORIES, LLC

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.

Work Order: C1703015

Project: 691 and 705 St Paul St

TestCode: 0.25CT-TCE-VC

Sample ID	AMB1UG-030817	Sample Type:	MBLK	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12014					
Client ID:	ZZZZZ	Batch ID:	R12014	TestNo: TO-15		Analysis Date: 3/8/2017	SeqNo: 140534					
Analyte		Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene		< 0.15	0.15									
Chloroethane		< 0.15	0.15									
cis-1,2-Dichloroethene		< 0.15	0.15									
trans-1,2-Dichloroethene		< 0.15	0.15									
Trichloroethene		< 0.040	0.040									
Vinyl chloride		< 0.040	0.040									

Qualifiers:		Results reported are not blank corrected	E	Estimated Value above quantitation range	H	Holding times for preparation or analysis exceeded
J		Analyte detected below quantitation limit	ND	Not Detected at the Limit of Detection	R	RPD outside accepted recovery limits
S		Spike Recovery outside accepted recovery limits				

Page 1 of 1

Date: 27-Mar-17

## CENTEK LABORATORIES, LLC

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.  
 Work Order: C1703015  
 Project: 691 and 705 St Paul St

TestCode: 1ugM3\_TO15

Sample ID	C1703015-003A MS	SampleType: MS	Batch ID: R12014	TestCode: 1ugM3_TO15	Units: ppbV	Prep Date:	RunNo: 12014				
Client ID:	SV-3			TestNo: TO-15		Analysis Date: 3/8/2017	SeqNo: 140546				
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	1.090	0.15	1	0	109	70	130				
Chloroethane	0.8100	0.15	1	0.11	70.0	70	130				
cis-1,2-Dichloroethene	1.400	0.15	1	0.18	122	70	130				
trans-1,2-Dichloroethene	1.140	0.15	1	0	114	70	130				
Trichloroethene	9.890	0.15	1	8.34	155	70	130				S
Vinyl chloride	0.7200	0.15	1	0	72.0	70	130				

Sample ID	C1703015-003A MS	SampType: MSD	Batch ID: R12014	TestCode: 1ugM3_TO15	Units: ppbV	Prep Date:	RunNo: 12014				
Client ID:	SV-3			TestNo: TO-15		Analysis Date: 3/8/2017	SeqNo: 140547				
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	1.160	0.15	1	0	116	70	130	1.09	6.22	30	
Chloroethane	0.8600	0.15	1	0.11	75.0	70	130	0.81	5.99	30	
cis-1,2-Dichloroethene	1.440	0.15	1	0.18	126	70	130	1.4	2.82	30	
trans-1,2-Dichloroethene	1.180	0.15	1	0	118	70	130	1.14	3.45	30	
Trichloroethene	9.550	0.15	1	8.34	121	70	130	9.89	3.50	30	
Vinyl chloride	0.7500	0.15	1	0	75.0	70	130	0.72	4.08	30	

Qualifiers: J Results reported are not blank corrected  
 S Analyte detected below quantitation limit  
 R Spike Recovery outside accepted recovery limits  
 E Estimated Value above quantitation range  
 ND Not Detected at the Limit of Detection  
 H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits

Page 1 of 1

Centek Laboratories  
IDL Study1ug/M3 Detection Limit  
January 2016Method TO-15A  
Units=ppb

Name	Amount	IDL#1	IDL#2	IDL#3	IDL#4	IDL#5	IDL#6	IDL#7	Average	StdDev	%Rec	IDL
Propylene	0.15	0.16	0.15	0.16	0.14	0.16	0.14	0.16	0.153	0.010	98.1	0.030
Freon 12	0.15	0.18	0.17	0.17	0.17	0.18	0.17	0.17	0.173	0.005	86.8	0.015
Chloromethane	0.15	0.19	0.18	0.16	0.18	0.18	0.2	0.17	0.180	0.013	83.3	0.041
Freon 114	0.15	0.18	0.17	0.16	0.17	0.18	0.17	0.18	0.174	0.005	86.1	0.017
Vinyl Chloride	0.15	0.17	0.16	0.16	0.15	0.16	0.15	0.15	0.157	0.008	95.5	0.024
Butane	0.15	0.18	0.16	0.17	0.18	0.18	0.19	0.19	0.179	0.011	84.0	0.034
1,3-butadiene	0.15	0.21	0.2	0.2	0.22	0.17	0.18	0.23	0.201	0.021	74.5	0.066
Bromomethane	0.15	0.18	0.2	0.21	0.18	0.22	0.16	0.21	0.194	0.021	77.2	0.068
Chloroethane	0.15	0.19	0.19	0.16	0.19	0.19	0.18	0.19	0.184	0.011	81.4	0.036
Ethanol	0.15	0.16	0.16	0.18	0.17	0.19	0.18	0.19	0.176	0.013	85.4	0.040
Acrolein	0.15	0.22	0.17	0.19	0.16	0.18	0.21	0.17	0.186	0.022	80.8	0.070
Vinyl Bromide	0.15	0.17	0.15	0.16	0.16	0.17	0.17	0.17	0.164	0.008	91.3	0.025
Freon 11	0.15	0.18	0.17	0.17	0.18	0.19	0.17	0.18	0.177	0.008	84.7	0.024
Acetone	0.15	0.2	0.17	0.18	0.15	0.15	0.18	0.14	0.167	0.021	89.7	0.067
Pentane	0.15	0.18	0.17	0.18	0.16	0.17	0.2	0.15	0.174	0.014	86.1	0.044
Isopropyl alcohol	0.15	0.22	0.2	0.19	0.2	0.19	0.21	0.19	0.200	0.012	75.0	0.036
1,1-dichloroethane	0.15	0.2	0.17	0.19	0.19	0.19	0.18	0.18	0.186	0.010	80.8	0.031
Freon 113	0.15	0.17	0.16	0.18	0.18	0.18	0.17	0.17	0.173	0.008	86.8	0.024
t-Butyl alcohol	0.15	0.21	0.2	0.2	0.21	0.2	0.2	0.18	0.200	0.010	75.0	0.031
Methylene chloride	0.15	0.2	0.18	0.19	0.18	0.2	0.19	0.17	0.187	0.011	80.2	0.035
Allyl chloride	0.15	0.18	0.17	0.16	0.18	0.18	0.2	0.18	0.179	0.012	84.0	0.038
Carbon disulfide	0.15	0.2	0.17	0.19	0.19	0.2	0.18	0.19	0.189	0.011	79.5	0.034
trans-1,2-dichloroethane	0.15	0.15	0.14	0.14	0.14	0.16	0.14	0.15	0.146	0.008	102.9	0.025
methyl tert-butyl ether	0.15	0.14	0.14	0.14	0.13	0.15	0.14	0.13	0.139	0.007	108.2	0.022
1,1-dichloroethane	0.15	0.17	0.15	0.16	0.15	0.17	0.16	0.16	0.160	0.008	93.8	0.026
Vinyl acetate	0.15	0.14	0.13	0.14	0.13	0.13	0.13	0.12	0.131	0.007	114.1	0.022
Methyl Ethyl Ketone	0.15	0.17	0.17	0.16	0.16	0.15	0.13	0.12	0.151	0.020	99.1	0.061
cis-1,2-dichloroethane	0.15	0.15	0.14	0.16	0.15	0.16	0.15	0.14	0.150	0.008	100.0	0.026
Hexane	0.15	0.12	0.14	0.13	0.13	0.13	0.12	0.12	0.127	0.008	118.0	0.024
Ethyl acetate	0.15	0.16	0.17	0.14	0.15	0.14	0.16	0.13	0.150	0.014	100.0	0.044
Chloroform	0.15	0.16	0.16	0.16	0.16	0.17	0.16	0.17	0.163	0.005	92.1	0.015
Tetrahydrofuran	0.15	0.15	0.13	0.15	0.15	0.15	0.15	0.14	0.146	0.008	102.9	0.025
1,2-dichloroethane	0.15	0.16	0.15	0.16	0.16	0.17	0.16	0.17	0.161	0.007	92.9	0.022
1,1,1-trichloroethane	0.15	0.17	0.16	0.17	0.17	0.16	0.17	0.17	0.167	0.005	89.7	0.015
Cyclohexane	0.15	0.14	0.14	0.14	0.15	0.15	0.14	0.14	0.143	0.005	105.0	0.015
Carbon tetrachloride	0.15	0.13	0.15	0.15	0.15	0.15	0.15	0.16	0.149	0.009	101.0	0.028
Benzene	0.15	0.15	0.16	0.16	0.15	0.16	0.16	0.16	0.157	0.005	95.5	0.015
Methyl methacrylate	0.15	0.15	0.15	0.14	0.14	0.14	0.15	0.11	0.140	0.014	107.1	0.044
1,4-dioxane	0.15	0.18	0.18	0.19	0.18	0.15	0.17	0.12	0.167	0.024	89.7	0.076

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1/8/2016

Centek Laboratories IDL Study		1ug/M3 Detection Limit January 2016							Method TO-15A Units=ppb			
Name	Amount	IDL#1	IDL#2	IDL#3	IDL#4	IDL#5	IDL#6	IDL#7	Average	StdDev	%Rec	IDL
2,2,4-trimethylpentane	0.15	0.15	0.15	0.15	0.16	0.14	0.16	0.15	0.151	0.007	99.1	0.022
Heptane	0.15	0.12	0.13	0.13	0.12	0.13	0.13	0.13	0.127	0.005	118.0	0.015
Trichloroethene	0.15	0.14	0.15	0.14	0.15	0.15	0.14	0.15	0.146	0.005	102.9	0.017
1,2-dichloropropane	0.15	0.16	0.17	0.17	0.16	0.17	0.16	0.16	0.184	0.005	91.3	0.017
Bromodichloromethane	0.15	0.16	0.16	0.16	0.15	0.16	0.17	0.16	0.160	0.006	93.8	0.018
cis-1,3-dichloropropene	0.15	0.13	0.13	0.14	0.14	0.13	0.13	0.13	0.133	0.005	112.9	0.015
trans-1,3-dichloropropene	0.15	0.16	0.13	0.13	0.14	0.14	0.14	0.16	0.143	0.013	105.0	0.039
1,1,2-trichloroethane	0.15	0.16	0.15	0.16	0.15	0.16	0.18	0.17	0.161	0.011	92.9	0.034
Toluene	0.15	0.14	0.14	0.14	0.13	0.16	0.14	0.15	0.143	0.010	105.0	0.030
Methyl Isobutyl Ketone	0.15	0.18	0.18	0.18	0.18	0.16	0.18	0.15	0.173	0.013	86.8	0.039
Dibromochloromethane	0.15	0.16	0.15	0.17	0.16	0.16	0.17	0.18	0.169	0.009	89.0	0.028
Methyl Butyl Ketone	0.15	0.17	0.16	0.18	0.17	0.16	0.17	0.14	0.164	0.013	91.3	0.040
1,2-dibromoethane	0.15	0.16	0.17	0.16	0.16	0.16	0.16	0.17	0.163	0.005	92.1	0.015
Tetrachloroethylene	0.15	0.16	0.17	0.16	0.16	0.16	0.17	0.17	0.164	0.005	91.3	0.017
Chlorobenzene	0.15	0.16	0.16	0.16	0.17	0.15	0.17	0.17	0.163	0.008	92.1	0.024
1,1,1,2-tetrachloroethane	0.15	0.17	0.17	0.17	0.18	0.16	0.18	0.17	0.171	0.007	87.5	0.022
Ethylbenzene	0.15	0.13	0.14	0.14	0.14	0.12	0.14	0.13	0.134	0.008	111.7	0.025
m&p-xylene	0.3	0.25	0.25	0.25	0.23	0.25	0.25	0.25	0.247	0.008	121.4	0.024
Nonane	0.15	0.11	0.11	0.11	0.11	0.1	0.1	0.11	0.107	0.005	140.0	0.015
Styrene	0.15	0.12	0.13	0.13	0.11	0.12	0.13	0.12	0.123	0.008	122.1	0.024
Bromoform	0.15	0.15	0.15	0.16	0.15	0.15	0.17	0.16	0.156	0.008	96.3	0.025
o-xylene	0.15	0.11	0.12	0.12	0.14	0.14	0.12	0.11	0.123	0.013	122.1	0.039
Cumene	0.15	0.12	0.13	0.13	0.12	0.13	0.13	0.13	0.127	0.005	118.0	0.015
Bromofluorobenzene	1	0.88	0.9	0.9	0.87	0.89	0.89	0.9	0.890	0.012	112.4	0.036
1,1,2,2-tetrachloroethane	0.15	0.16	0.16	0.17	0.16	0.17	0.17	0.16	0.164	0.005	91.3	0.017
Propylbenzene	0.15	0.13	0.12	0.13	0.13	0.11	0.13	0.11	0.123	0.010	122.1	0.030
2-Chlorotoluene	0.15	0.13	0.13	0.13	0.14	0.13	0.12	0.13	0.130	0.006	115.4	0.018
4-ethyltoluene	0.15	0.11	0.12	0.12	0.12	0.13	0.13	0.11	0.120	0.008	125.0	0.026
1,3,5-trimethylbenzene	0.15	0.12	0.13	0.14	0.12	0.13	0.13	0.13	0.129	0.007	116.7	0.022
1,2,4-trimethylbenzene	0.15	0.12	0.13	0.12	0.12	0.13	0.12	0.12	0.123	0.005	122.1	0.015
1,3-dichlorobenzene	0.15	0.14	0.14	0.14	0.13	0.14	0.13	0.14	0.137	0.005	109.4	0.015
benzyl chloride	0.15	0.13	0.16	0.13	0.15	0.13	0.15	0.16	0.144	0.014	104.0	0.044
1,4-dichlorobenzene	0.15	0.13	0.11	0.12	0.12	0.12	0.12	0.13	0.121	0.007	123.5	0.022
1,2,3-trimethylbenzene	0.15	0.12	0.11	0.12	0.12	0.12	0.11	0.11	0.116	0.005	129.6	0.017
1,2-dichlorobenzene	0.15	0.13	0.14	0.14	0.14	0.14	0.14	0.13	0.137	0.005	109.4	0.015
1,2,4-trichlorobenzene	0.15	0.1	0.11	0.1	0.11	0.11	0.12	0.1	0.107	0.008	140.0	0.024
Naphthalene	0.15	0.13	0.13	0.14	0.11	0.12	0.14	0.12	0.127	0.011	118.0	0.035
Hexachloro-1,3-butadiene	0.15	0.16	0.17	0.17	0.17	0.16	0.16	0.16	0.164	0.005	91.3	0.017

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1/8/2016

Centek Laboratories  
IDL Study0.25ug/M3 Detection Limit  
January 2016Method TO-15A  
Units=ppb

Name	Amount	IDL#1	IDL#2	IDL#3	IDL#4	IDL#5	IDL#6	IDL#7	Average	StdDev	%Rec	IDL
Vinyl Chloride	0.1	0.11	0.11	0.09	0.09	0.1	0.09	0.1	0.099	0.009	101.4	0.028
Carbon tetrachloride	0.1	0.1	0.11	0.08	0.09	0.09	0.09	0.09	0.093	0.010	107.7	0.030
Trichloroethene	0.1	0.1	0.1	0.07	0.08	0.08	0.08	0.08	0.084	0.011	118.6	0.036
Tetrachloroethylene	0.1	0.11	0.12	0.09	0.09	0.1	0.09	0.09	0.099	0.012	101.4	0.038
Naphthalene	0.1	0.09	0.08	0.07	0.06	0.06	0.07	0.06	0.070	0.012	142.9	0.036

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1/15/2016

## GC/MS-Whole Air Calculations

## Relative Response Factor (RRF)

$$RRF = \frac{A_x * C_{is}}{A_{is} * C_x}$$

where:  $A_x$  = area of the characteristic ion for the compound being measured  
 $A_{is}$  = area of the characteristic ion for the specific internal standard of the compound being measured  
 $C_x$  = concentration of the compound being measured (ppbv)  
 $C_{is}$  = concentration of the internal standard (ppbv)

## Percent Relative Standard Deviation (%RSD)

$$\% RSD = \frac{\text{Standard deviation of RRF values} * 100}{\text{mean RRF}}$$

## Percent Difference (%D)

$$\% D = \frac{(RRF_c - \text{mean RRF}_i) * 100}{\text{mean RRF}_i}$$

where:  $RRF_c$  = relative response factor from the continuing calibration  
 $\text{mean RRF}_i$  = mean relative response factor from the initial calibration

## Sample Calculations

$$\text{ppbv} = \frac{A_x * I_s * D_f}{A_{is} * RRF}$$

where:  $A_x$  = area of the characteristic ion for the compound being measured  
 $A_{is}$  = area of the characteristic ion for the specific internal standard of the compound being measured  
 $I_s$  = Concentration of the internal standard injected (ppbv)  
 $RRF$  = relative response factor for the compound being measured  
 $D_f$  = Dilution factor

**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**SAMPLE DATA**

## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-001A

Client Sample ID: SV-1  
 Tag Number: 542.256  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
----------	--------	---------	------	-------	----	---------------

## FIELD PARAMETERS

## FLD

Analyst:

Lab Vacuum In

-1

"Hg

3/7/2017

Lab Vacuum Out

-30

"Hg

3/7/2017

## 1UG/M3 BY METHOD TO15

## TO-15

Analyst: RJP

1,1-Dichloroethene

&lt; 0.15

0.15

ppbV

1

3/8/2017 1:28:00 PM

Chloroethane

0.13

0.15

J

ppbV

1

3/8/2017 1:28:00 PM

cis-1,2-Dichloroethene

&lt; 0.15

0.15

ppbV

1

3/8/2017 1:28:00 PM

trans-1,2-Dichloroethane

&lt; 0.15

0.15

ppbV

1

3/8/2017 1:28:00 PM

Trichloroethene

&lt; 0.15

0.15

ppbV

1

3/8/2017 1:28:00 PM

Vinyl chloride

&lt; 0.15

0.15

ppbV

1

3/8/2017 1:28:00 PM

Surr: Bromofluorobenzene

239

70-130

S

%REC

1

3/8/2017 1:28:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte, Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits

. Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

Page 1 of 6



## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-001A

Client Sample ID: SV-1  
 Tag Number: 542.256  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>1UG/M3 BY METHOD TO15</b>		<b>TO-15</b>		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 1:28:00 PM
Chloroethane	0.34	0.40	J	ug/m3	1	3/8/2017 1:28:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 1:28:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 1:28:00 PM
Trichloroethene	< 0.81	0.81		ug/m3	1	3/8/2017 1:28:00 PM
Vinyl chloride	< 0.38	0.38		ug/m3	1	3/8/2017 1:28:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte, Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

Data File : C:\HPCHEM\1\DATA\AO030807.D

Vial: 22

Acq On : 8 Mar 2017 1:28 pm

Operator: RJP

Sample : C1703015-001A

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:17 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.65	128	59338	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.04	114	270574	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	278336	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	416764m <i>P</i>	2.39	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	239.00%#

## Target Compounds

					Qvalue
10) Chloroethane	5.63	64	7718	0.13	ppb # 64
28) Methyl Ethyl Ketone	8.93	72	32252	0.96	ppb # 1

Data File : C:\HPCHEM\1\DATA\AO030807.D  
Acq On : 8 Mar 2017 1:28 pm  
Sample : C1703015-001A  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 10 12:13 2017

Vial: 22  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Mar 27 11:00:28 2017  
Response via : Initial Calibration

Abundance

2.5e+07

2.4e+07

2.2e+07

2e+07

1.8e+07

1.6e+07

1.4e+07

1.2e+07

1e+07

8000000

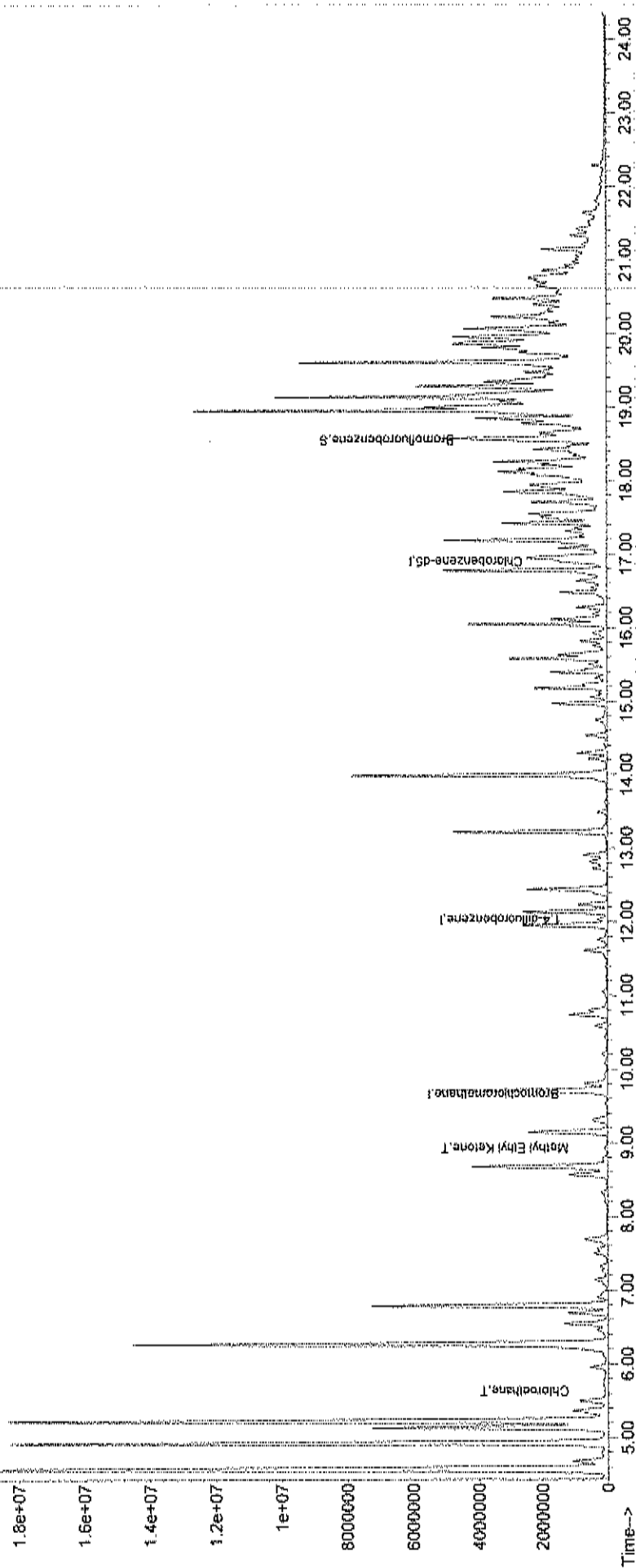
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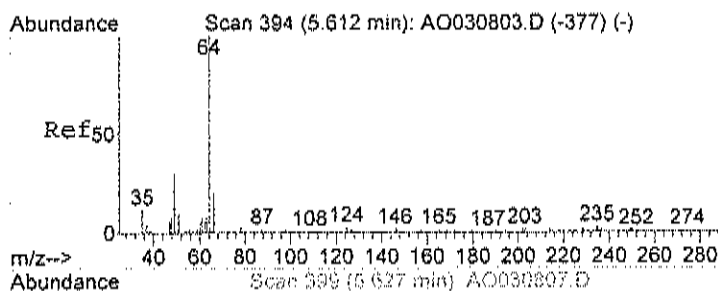
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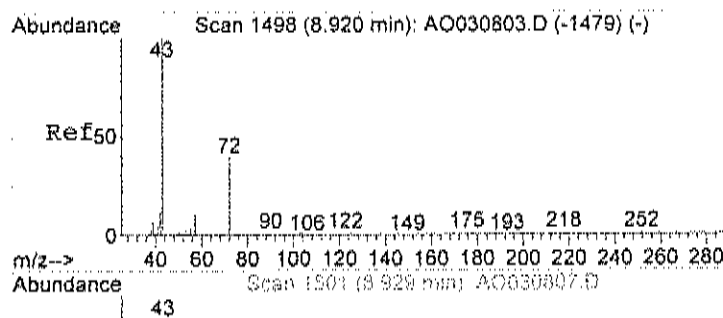
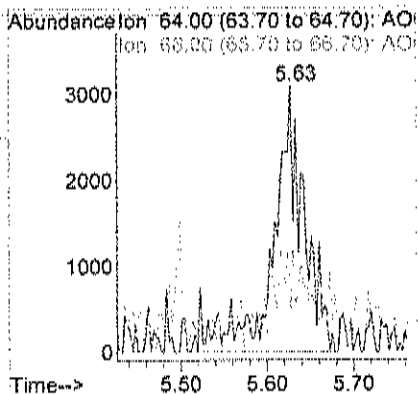
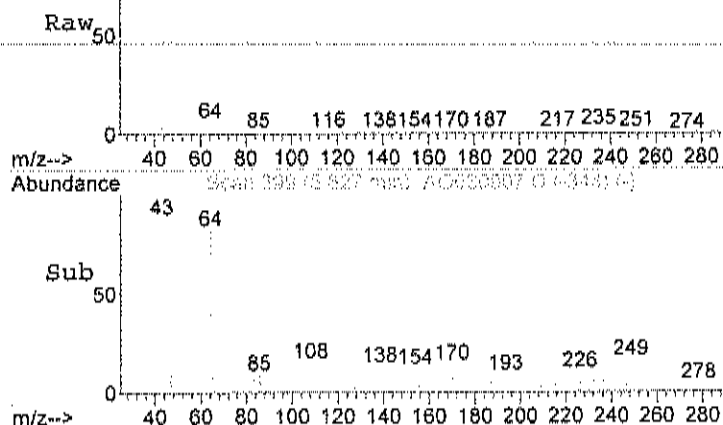
TIC: AO030807.D





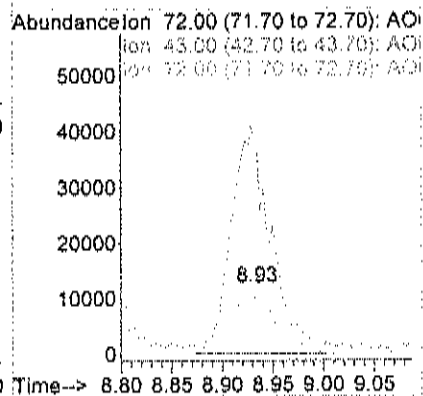
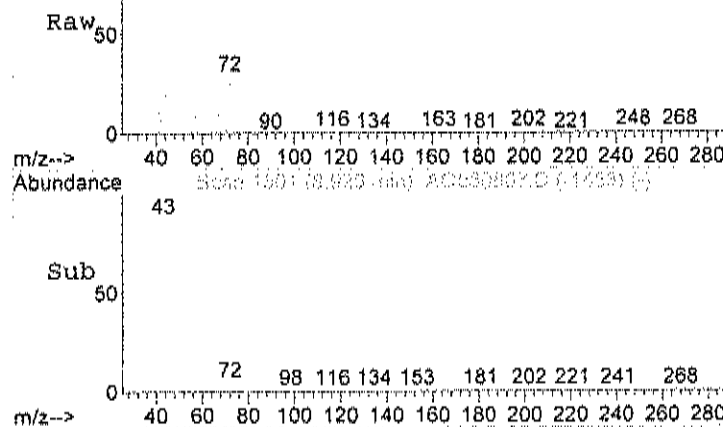
#10  
Chloroethane  
Concen: 0.13 ppb  
RT: 5.63 min Scan# 399  
Delta R.T. 0.02 min  
Lab File: AO030807.D  
Acq: 8 Mar 2017 1:28 pm

Tgt Ion	Ratio	Lower	Upper
64	100		
66	47.5	22.7	34.1#



#28  
Methyl Ethyl Ketone  
Concen: 0.96 ppb  
RT: 8.93 min Scan# 1501  
Delta R.T. -0.01 min  
Lab File: AO030807.D  
Acq: 8 Mar 2017 1:28 pm

Tgt Ion	Ratio	Lower	Upper
72	100		
43	0.0	383.1	423.1#
72	100.0	80.0	120.0



## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-002A

Client Sample ID: SV-2  
 Tag Number: 237.402  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	-1			"Hg		3/7/2017
Lab Vacuum Out	-30			"Hg		3/7/2017
<b>1UG/M3 BY METHOD TO15</b>						
		TO-15				Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 2:09:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	3/8/2017 2:09:00 PM
cis-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 2:09:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 2:09:00 PM
Trichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 2:09:00 PM
Vinyl chloride	< 0.15	0.15		ppbV	1	3/8/2017 2:09:00 PM
Surr: Bromofluorobenzene	235	70-130	S	%REC	1	3/8/2017 2:09:00 PM

**NOTES:**

Surrogate did not meet criteria due to severe matrix interference.

Qualifiers:	**	Quantitation Limit	,	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-002A

Client Sample ID: SV-2  
 Tag Number: 237.402  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY METHOD TO15		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 2:09:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	3/8/2017 2:09:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 2:09:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 2:09:00 PM
Trichloroethene	< 0.81	0.81		ug/m3	1	3/8/2017 2:09:00 PM
Vinyl chloride	< 0.38	0.38		ug/m3	1	3/8/2017 2:09:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte, Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits  
 , Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

Page 2 of 6

Data File : C:\HPCHEM\1\DATA\AO030808.D

Vial: 23

Acq On : 8 Mar 2017 2:09 pm

Operator: RJP

Sample : C1703015-002A

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:18 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.65	128	72867	1.00	ppb	-0.01
35) 1,4-difluorobenzene	12.03	114	317032	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	316562	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.55	95	466729m	2.35	ppb	0.02
Spiked Amount	1.000	Range	70 - 130	Recovery	=	235.00%#

Target Compounds

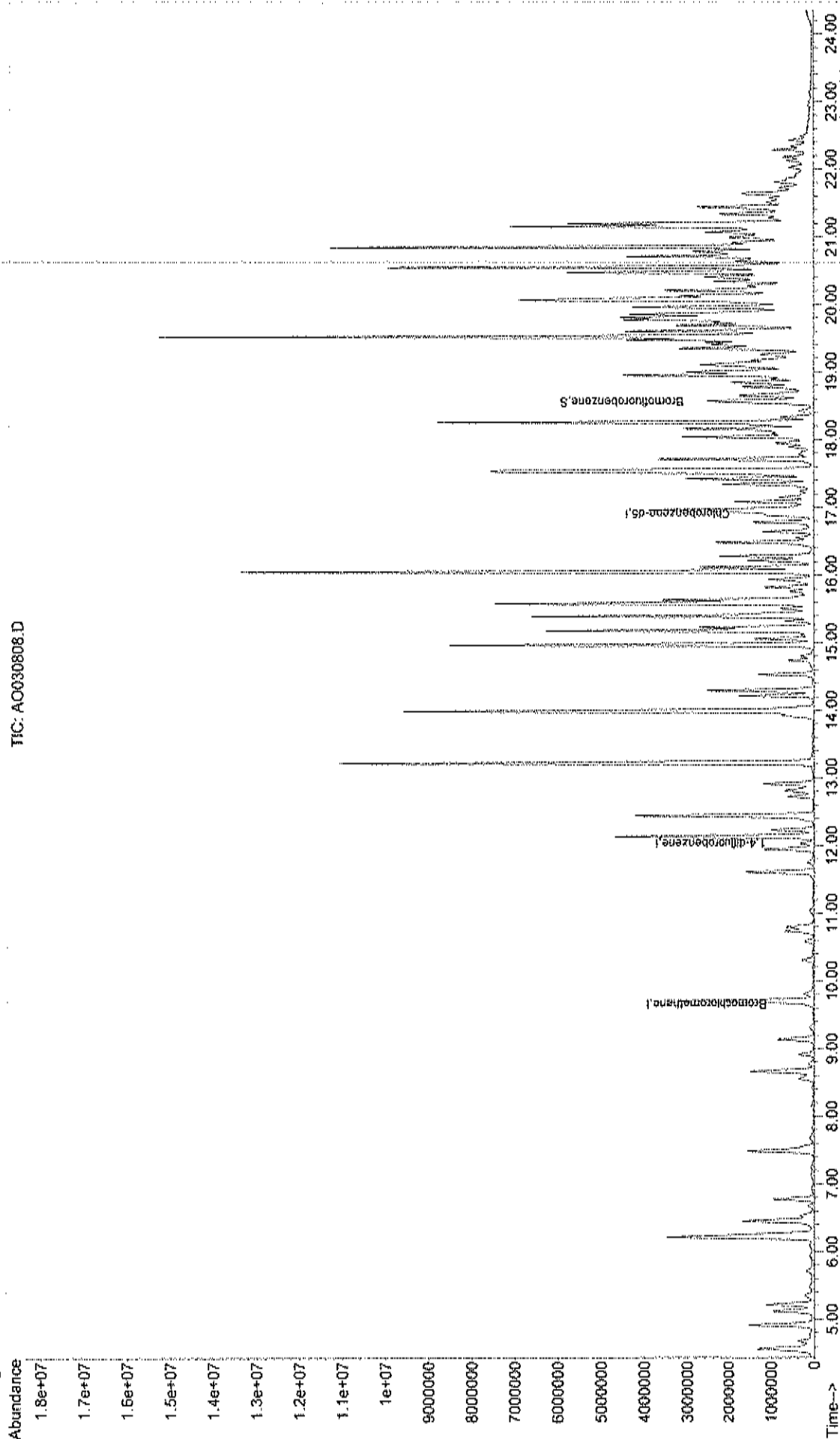
Qvalue

Data File : C:\HPCHEM\1\DATA\AO030808.D  
Acq On : 8 Mar 2017 2:09 pm  
Sample : C1703015-002A  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 10 12:14 2017

Vial: 23  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Mar 27 11:00:28 2017  
Response via : Initial Calibration





## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-003A

Client Sample ID: SV-3  
 Tag Number: 1206.249  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	1			"Hg		3/7/2017
Lab Vacuum Out	-30			"Hg		3/7/2017
<b>1UG/M3 BY METHOD TO15</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 2:48:00 PM
Chloroethane	0.11	0.15	J	ppbV	1	3/8/2017 2:48:00 PM
cis-1,2-Dichloroethene	0.18	0.15		ppbV	1	3/8/2017 2:48:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 2:48:00 PM
Trichloroethene	7.4	1.5		ppbV	10	3/8/2017 11:34:00 PM
Vinyl chloride	< 0.15	0.15		ppbV	1	3/8/2017 2:48:00 PM
Surr: Bromofluorobenzene	97.0	70-130		%REC	10	3/8/2017 11:34:00 PM
Surr: Bromofluorobenzene	166	70-130	S	%REC	1	3/8/2017 2:48:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte, Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

**Centek Laboratories, LLC**

Date: 27-Mar-17

**CLIENT:** LaBella Associates, P.C.  
**Lab Order:** C1703015  
**Project:** 691 and 705 St Paul St  
**Lab ID:** C1703015-003A

**Client Sample ID:** SV-3  
**Tag Number:** 1206.249  
**Collection Date:** 3/3/2017  
**Matrix:** AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>1UG/M3 BY METHOD TO15</b>				<b>TO-15</b>		<b>Analyst: RJP</b>
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 2:48:00 PM
Chloroethane	0.29	0.40	J	ug/m3	1	3/8/2017 2:48:00 PM
cis-1,2-Dichloroethene	0.71	0.59		ug/m3	1	3/8/2017 2:48:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 2:48:00 PM
Trichloroethene	40	8.1		ug/m3	10	3/8/2017 11:34:00 PM
Vinyl chloride	< 0.38	0.38		ug/m3	1	3/8/2017 2:48:00 PM

**NOTES:**

Surrogate did not meet criteria due to severe matrix interference.

<b>Qualifiers:</b>	<b>**</b>	Quantitation Limit	,	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

## Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\AO030809.D  
 Acq On : 8 Mar 2017 2:48 pm  
 Sample : C1703015-003A  
 Misc : A227\_1UG  
 MS Integration Params: RTEINT.P  
 Quant Time: Mar 09 11:10:19 2017

Vial: 24  
 Operator: RJP  
 Inst : MSD #1  
 Multiplr: 1.00

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Mon Mar 06 15:15:36 2017  
 Response via : Initial Calibration  
 DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.65	128	80896	1.00	ppb	-0.01
35) 1,4-difluorobenzene	12.03	114	366728	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	336237	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.55	95	350228m <sup>N</sup>	1.66	ppb	0.01
Spiked Amount	1.000	Range	70 - 130	Recovery	=	166.00%#

## Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
10) Chloroethane	5.62	64	9389	0.11	ppb	# 80
29) cis-1,2-dichloroethene	9.44	61	22541	0.18	ppb	92
44) Trichloroethene	12.88	130	1738106	8.34	ppb	95

Data File : C:\HPCHEM\1\DATA\AO030809.D  
Acq On : 8 Mar 2017 2:48 pm  
Sample : C1703015-003A  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 10 12:15 2017

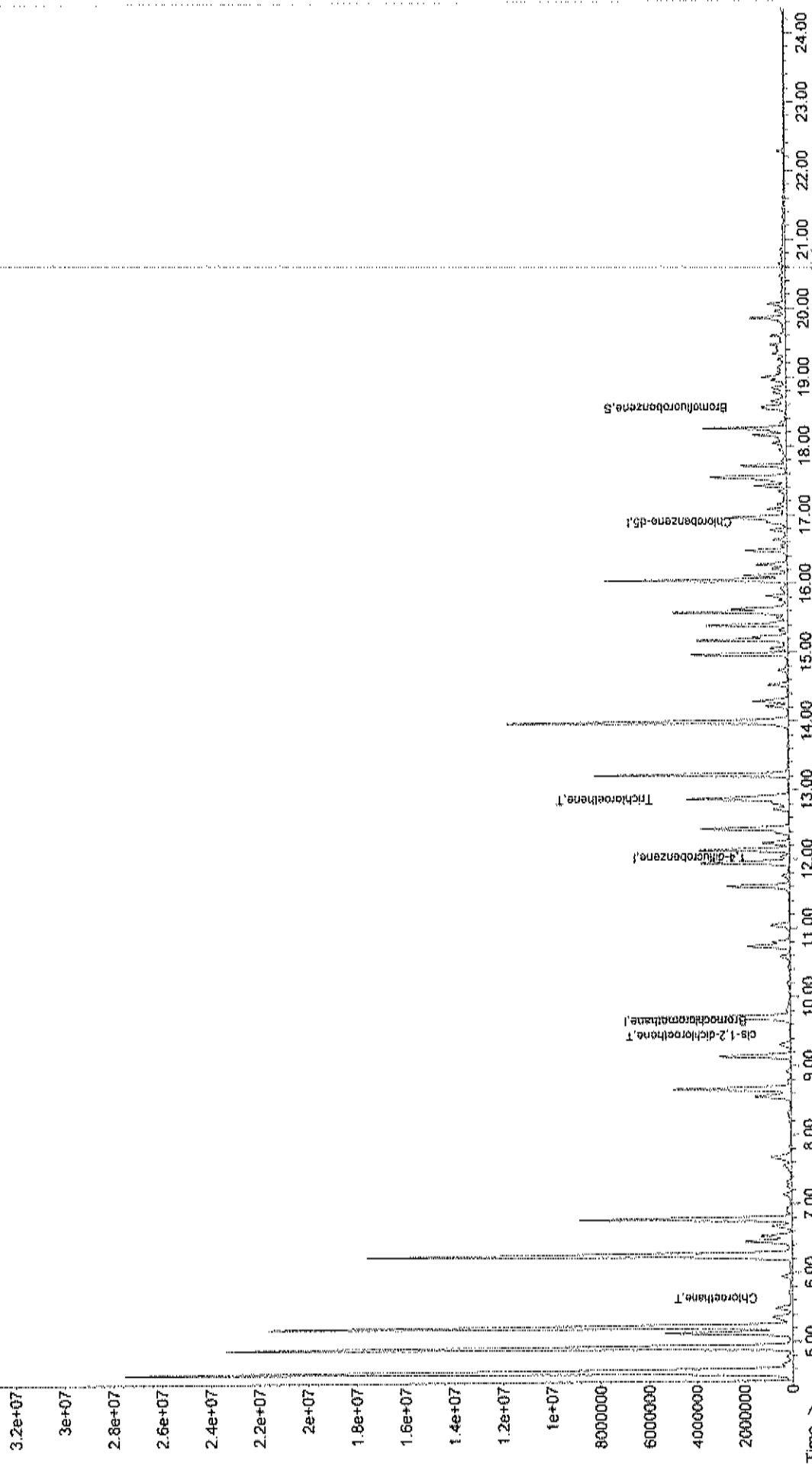
Vial: 24  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

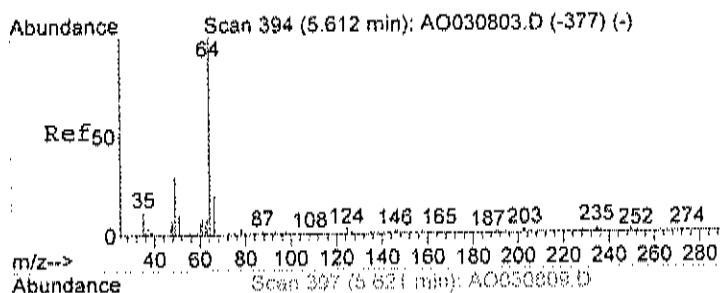
Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Mar 27 11:00:28 2017  
Response via : Initial Calibration

Abundance

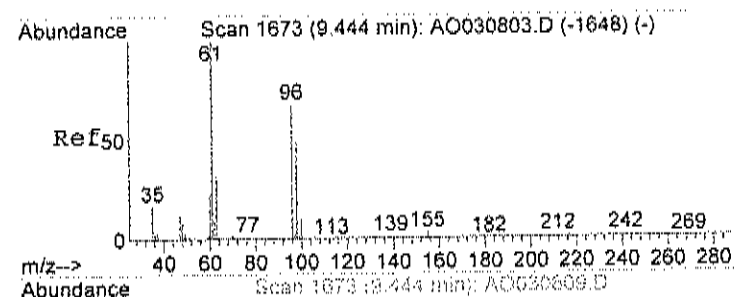
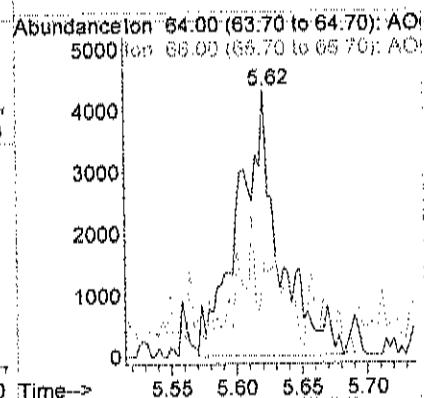
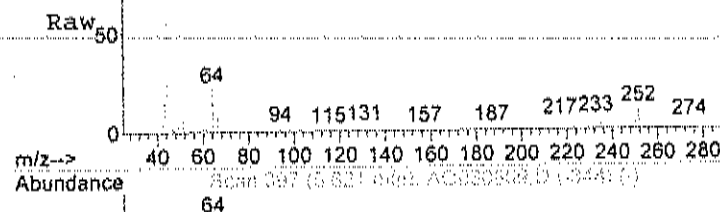
TIC: AO030809.D





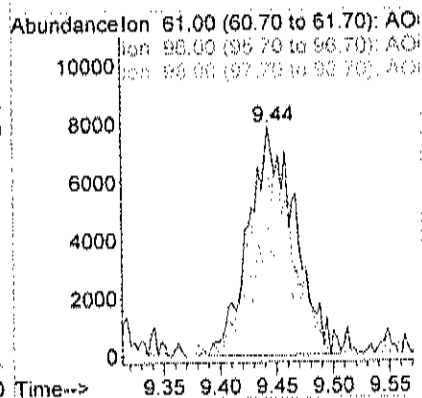
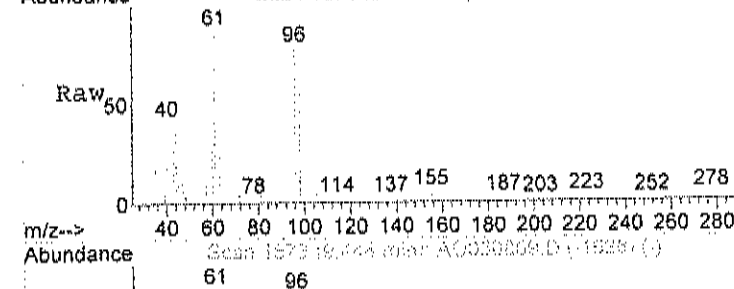
#10  
Chloroethane  
Concen: 0.11 ppb  
RT: 5.62 min Scan# 397  
Delta R.T. 0.01 min  
Lab File: AO030809.D  
Acq: 8 Mar 2017 2:48 pm

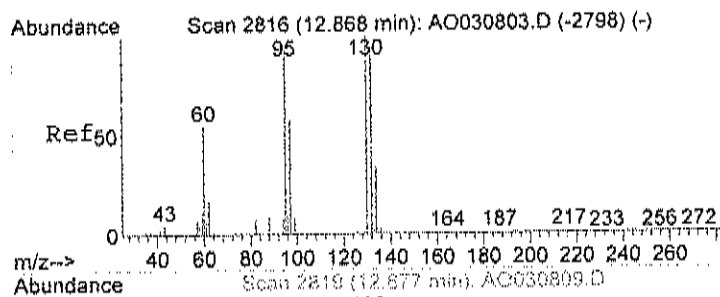
Tgt Ion: 64 Resp: 9389  
Ion Ratio Lower Upper  
64 100  
66 38.9 22.7 34.1#



#29  
cis-1,2-dichloroethene  
Concen: 0.18 ppb  
RT: 9.44 min Scan# 1673  
Delta R.T. -0.02 min  
Lab File: AO030809.D  
Acq: 8 Mar 2017 2:48 pm

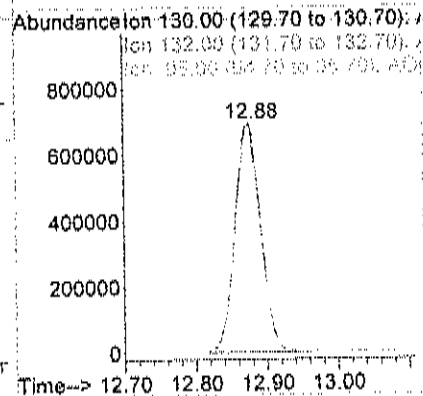
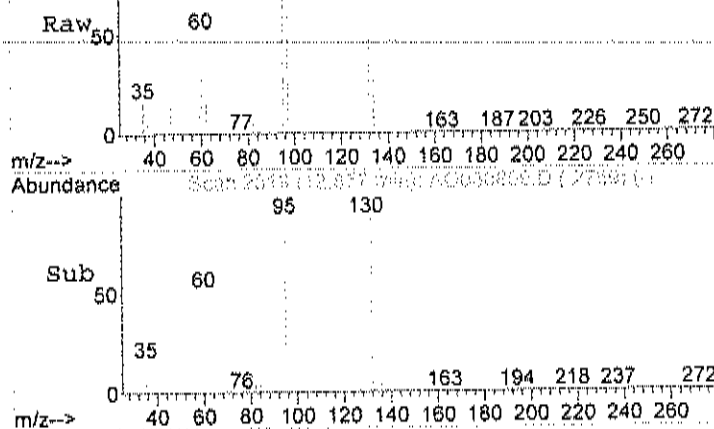
Tgt Ion: 61 Resp: 22541  
Ion Ratio Lower Upper  
61 100  
96 84.9 58.1 98.1  
98 54.4 29.3 69.3





#44  
Trichloroethene  
Concen: 8.34 ppb  
RT: 12.88 min Scan# 2819  
Delta R.T. -0.00 min  
Lab File: AO030809.D  
Acq: 8 Mar 2017 2:48 pm

Tgt Ion: 130 Resp: 1738106  
Ion Ratio Lower Upper  
130 100  
132 98.1 69.9 109.9  
95 97.0 76.3 116.3



Data File : C:\HPCHEM\1\DATA\AO030822.D

Vial: 36

Acq On : 8 Mar 2017 11:34 pm

Operator: RJP

Sample : C1703015-003A 10X

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:32 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.65	128	53827	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.04	114	227536	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	201273	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.53	95	122944	0.97	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	97.00%

## Target Compounds

44) Trichloroethene	12.87	130	95085	0.74	ppb	Qvalue 95
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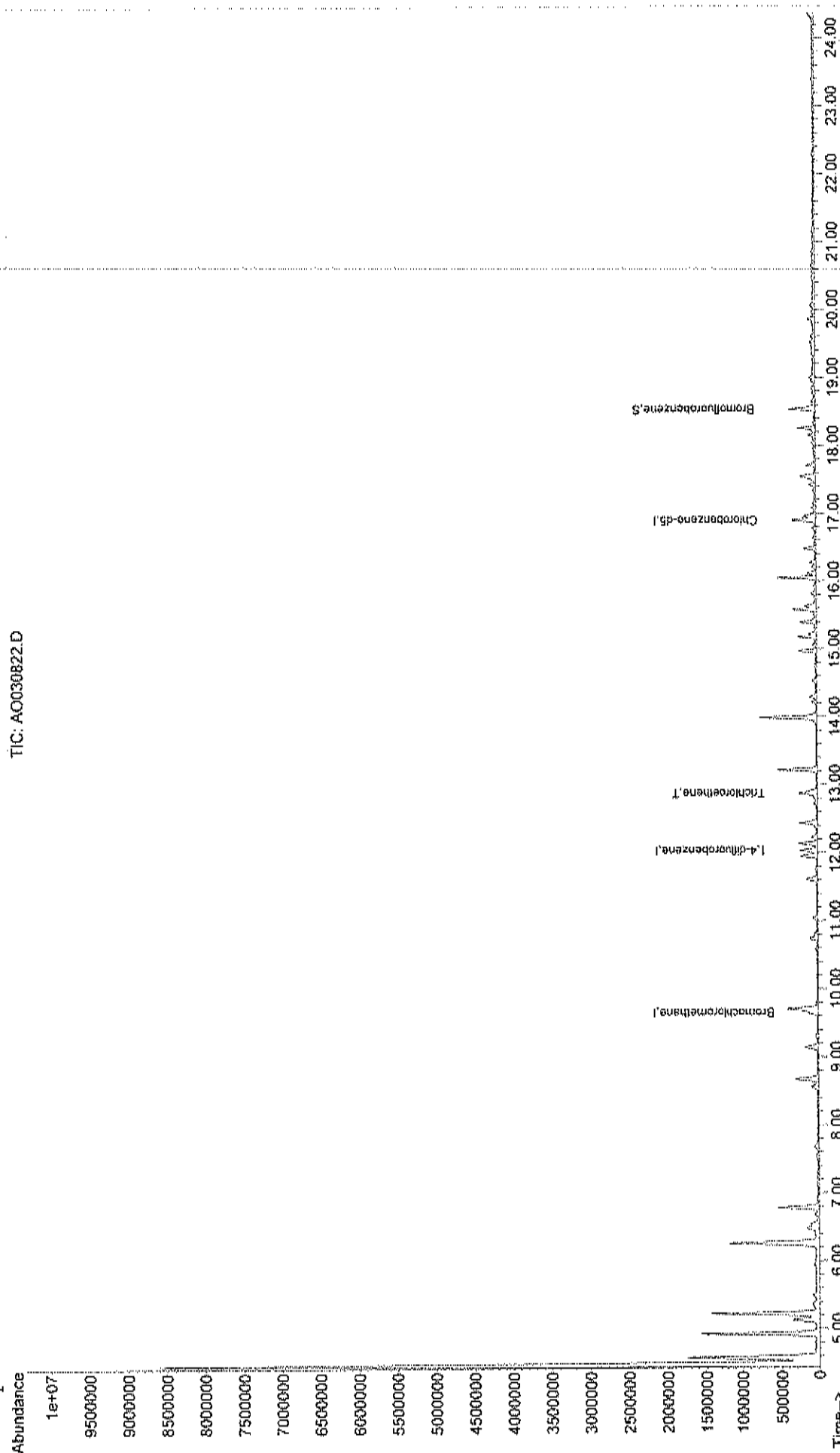
Data File : C:\HPCHEM\1\DATA\AO030822.D  
Acq On : 8 Mar 2017 11:34 pm  
Sample : C1703015-003A 10X  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 10 12:51 2017

Vial: 36  
Operator: RJP  
Inst : MSD #1  
Multiplx: 1.00

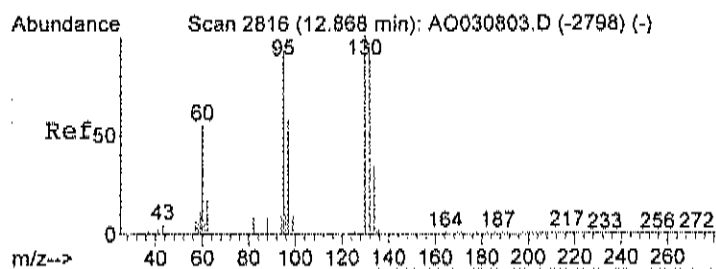
Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integration)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Mar 27 11:00:28 2017  
Response via : Initial Calibration

TIC: AO030822.D

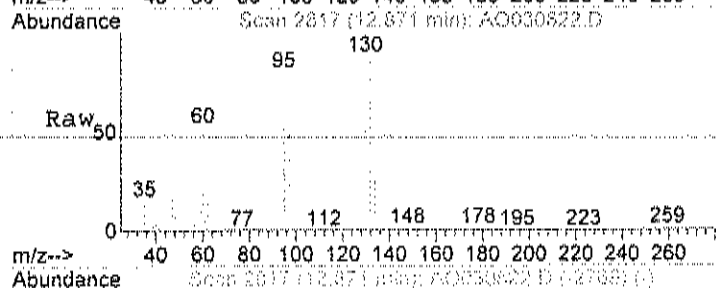






#44  
Trichloroethene  
Concen: 0.74 ppb  
RT: 12.87 min Scan# 2817  
Delta R.T. -0.01 min  
Lab File: AO030822.D  
Acq: 8 Mar 2017 11:34 pm

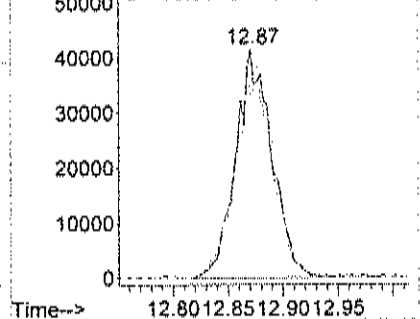
Tgt Ion	130	Resp	95085
Ion Ratio	Lower	Upper	
130	100		
132	96.9	69.9	109.9
95	94.6	76.3	116.3



Abundance Ion 130.00 (129.70 to 130.70):

Ion 132.00 (131.70 to 132.70):

Ion 95.00 (94.70 to 95.70):



## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-004A

Client Sample ID: SV-4  
 Tag Number: 290.299  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
				FLD		Analyst:
Lab Vacuum In	-3			"Hg		3/7/2017
Lab Vacuum Out	-30			"Hg		3/7/2017
<b>1UG/M3 BY METHOD TO15</b>						
				TO-15		Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 4:59:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	3/8/2017 4:59:00 PM
cis-1,2-Dichloroethene	1.0	0.15		ppbV	1	3/8/2017 4:59:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 4:59:00 PM
Trichloroethene	8.9	1.5		ppbV	10	3/9/2017 12:47:00 AM
Vinyl chloride	< 0.15	0.15		ppbV	1	3/8/2017 4:59:00 PM
Surr: Bromofluorobenzene	100	70-130		%REC	10	3/9/2017 12:47:00 AM
Surr: Bromofluorobenzene	190	70-130	S	%REC	1	3/8/2017 4:59:00 PM
<b>NOTES:</b>						
Surrogate did not meet criteria due to severe matrix interference.						

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits  
 , Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-004A

Client Sample ID: SV-4  
 Tag Number: 290.299  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY METHOD TO15		TO-15				Analyst: RJP
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 4:59:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	3/8/2017 4:59:00 PM
cis-1,2-Dichloroethene	4.1	0.59		ug/m3	1	3/8/2017 4:59:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 4:59:00 PM
Trichloroethene	48	8.1		ug/m3	10	3/9/2017 12:47:00 AM
Vinyl chloride	< 0.38	0.38		ug/m3	1	3/8/2017 4:59:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits  
 . Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

Page 4 of 6

Data File : C:\HPCHEM\1\DATA\AO030812.D

Vial: 27

Acq On : 8 Mar 2017 4:59 pm

Operator: RJP

Sample : C1703015-004A

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:22 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.65	128	76299	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.03	114	343779	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	323940	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	386106	1.90	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	190.00%#

## Target Compounds

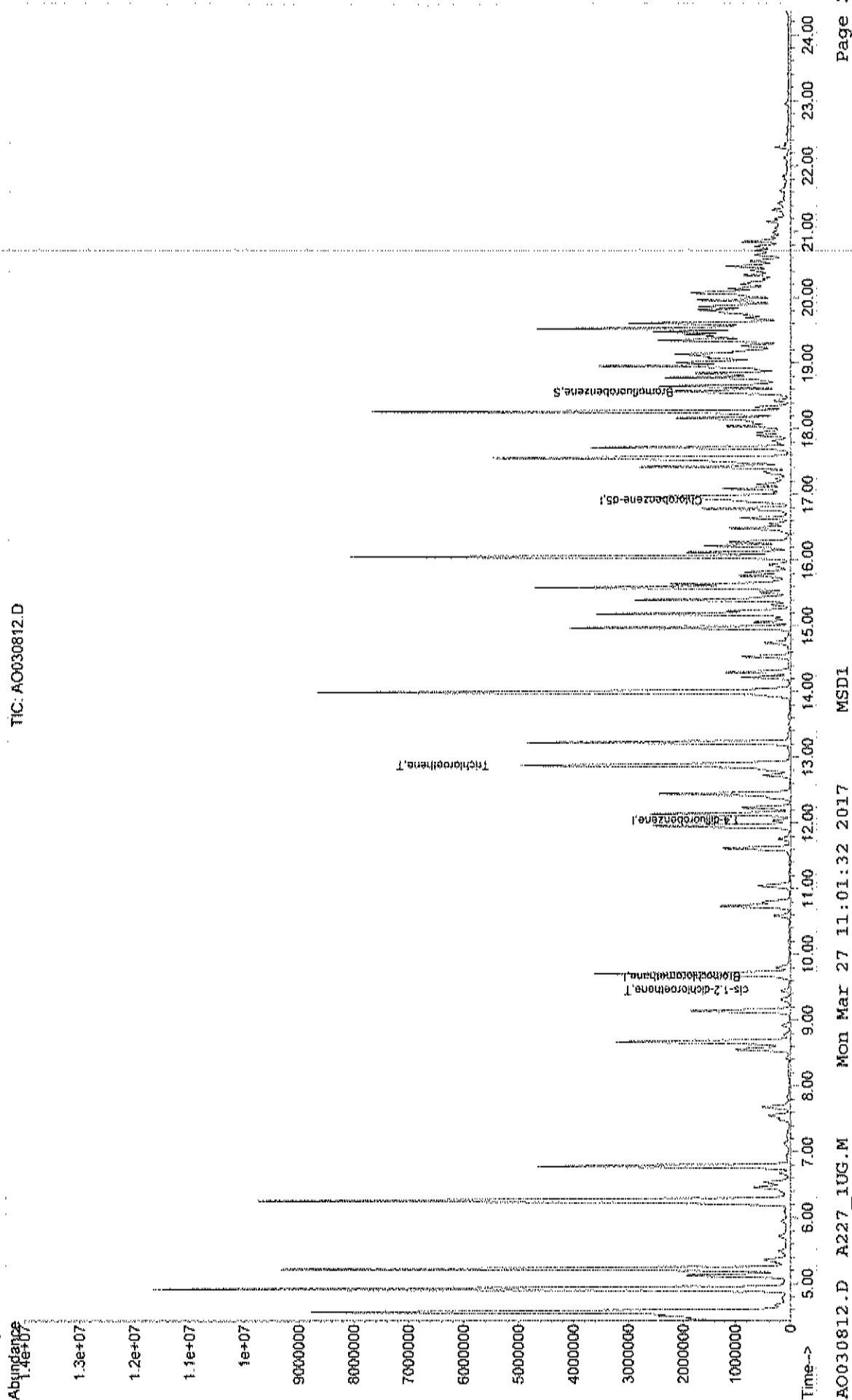
						Qvalue
29) cis-1,2-dichloroethene	9.44	61	126069	1.04	ppb	95
44) Trichloroethene	12.88	130	2043606	10.47	ppb	96

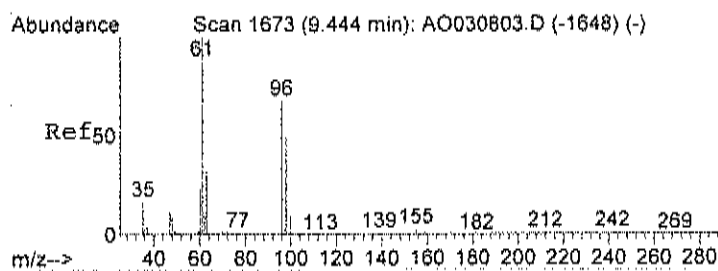
Data File : C:\HPCHEM\1\DATA\AO030812.D  
Acq On : 8 Mar 2017 4:59 pm  
Sample : C1703015-004A  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 10 12:37 2017

Vial: 27  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A227\_1UG.RES

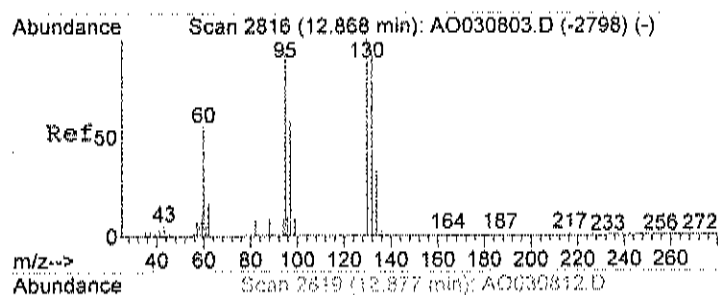
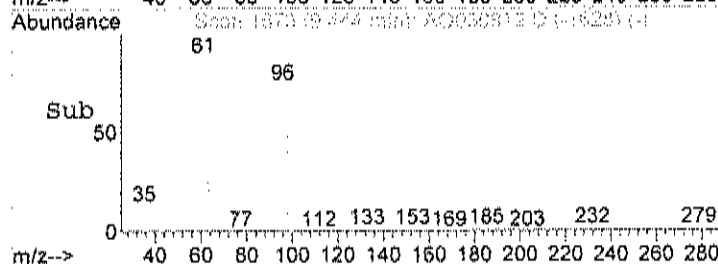
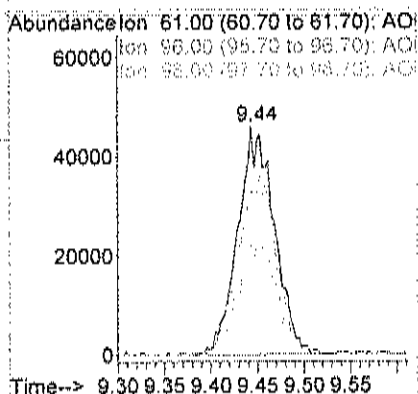
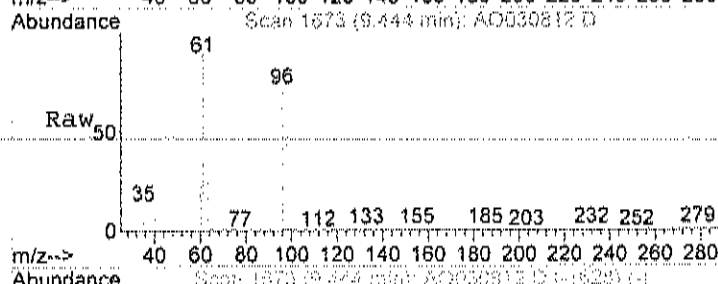
Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Mar 27 11:00:28 2017  
Response via : Initial Calibration





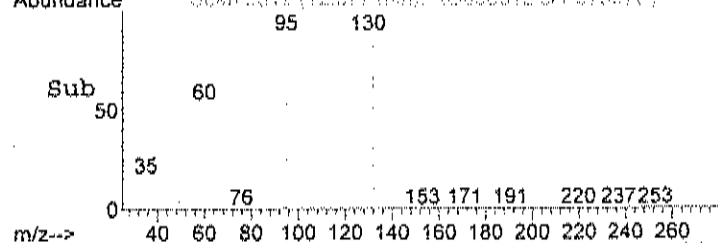
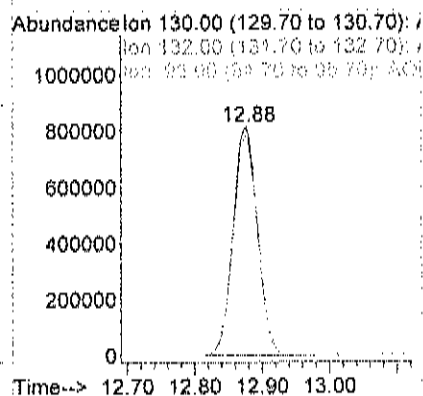
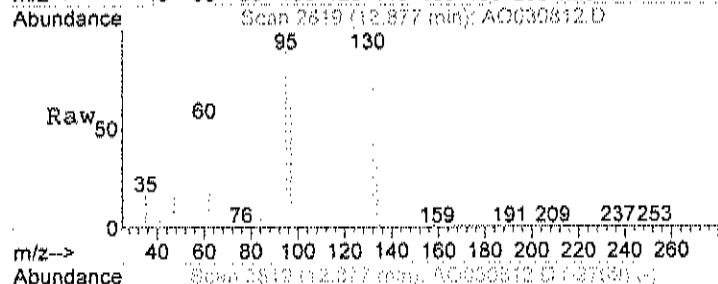
#29  
cis-1,2-dichloroethene  
Concen: 1.04 ppb  
RT: 9.44 min Scan# 1673  
Delta R.T. -0.02 min  
Lab File: AO030812.D  
Acq: 8 Mar 2017 4:59 pm

Tgt Ion:	61	Resp:	126069
Ion	Ratio	Lower	Upper
61	100		
96	82.1	58.1	98.1
98	52.9	29.3	69.3



#44  
Trichloroethene  
Concen: 10.47 ppb  
RT: 12.88 min Scan# 2819  
Delta R.T. -0.00 min  
Lab File: AO030812.D  
Acq: 8 Mar 2017 4:59 pm

Tgt Ion:	130	Resp:	2043606
Ion	Ratio	Lower	Upper
130	100		
132	96.4	69.9	109.9
95	95.1	76.3	116.3



Data File : C:\HPCHEM\1\DATA\AO030824.D

Vial: 38

Acq On : 9 Mar 2017 12:47 am

Operator: RJP

Sample : C1703015-004A 10X

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:34 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	9.65	128	46536	1.00	ppb	-0.01
35) 1,4-difluorobenzene	12.03	114	210199	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	197259	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	123688	1.00	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	100.00%

## Target Compounds

44) Trichloroethene	12.88	130	106647	0.89	ppb	Qvalue 96
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Data File : C:\HPCHEM\1\DATA\AO030824.D  
Acq On : 9 Mar 2017 12:47 am  
Sample : C1703015-004A 10X  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 10 12:52 2017

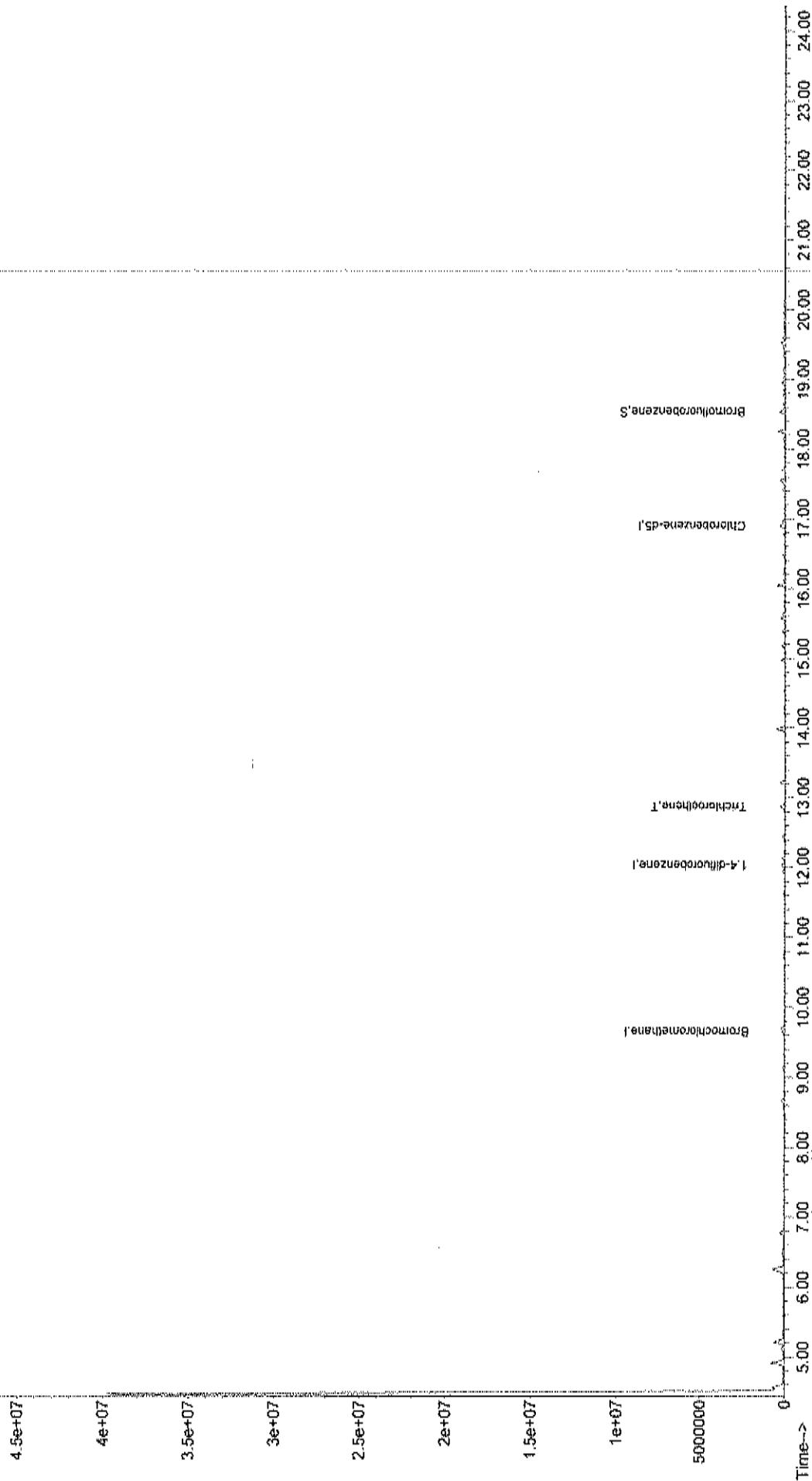
Vial: 38  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Mar 27 11:00:28 2017  
Response via : Initial Calibration

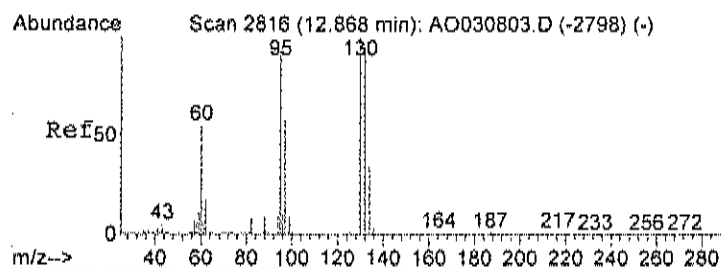
Abundance

TIC: AO030824.D



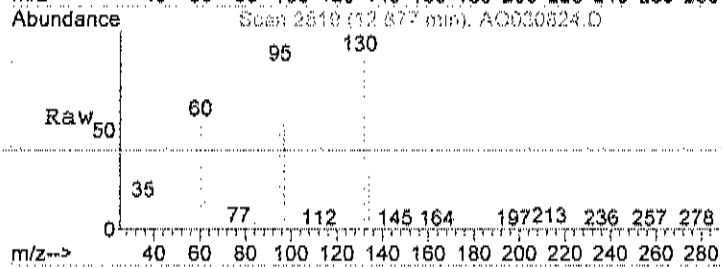
AO030824.D A227\_1UG.M Mon Mar 27 11:01:46 2017 MSD1





#44  
Trichloroethene  
Concen: 0.89 ppb  
RT: 12.88 min Scan# 2819  
Delta R.T. 0.00 min  
Lab File: AO030824.D  
Acq: 9 Mar 2017 12:47 am

Tgt Ion	Ratio	Lower	Upper
130	100		
132	93.9	69.9	109.9
95	92.8	76.3	116.3

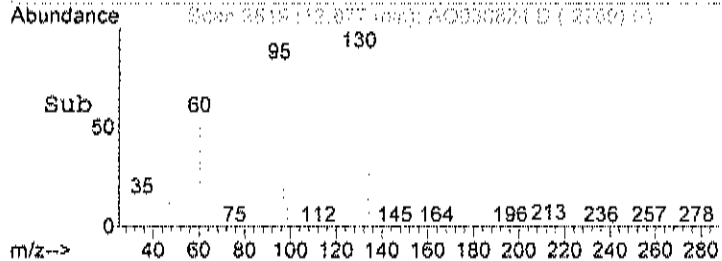
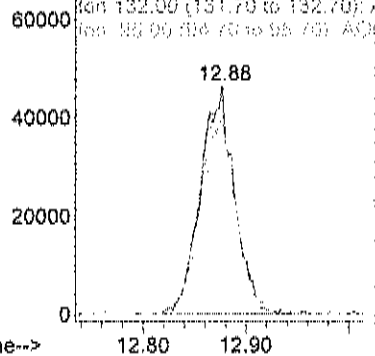


Abundance

Ion 130.00 (129.70 to 130.70)

Ion 132.00 (131.70 to 132.70)

Ion 95.00 (94.70 to 95.70)



## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-005A

Client Sample ID: Duplicate  
 Tag Number: 1184.299  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>		<b>FLD</b>		Analyst:		
Lab Vacuum In	-3			"Hg		3/7/2017
Lab Vacuum Out	-30			"Hg		3/7/2017
<b>1UG/M3 BY METHOD TO15</b>		<b>TO-15</b>		Analyst: RJP		
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 5:39:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	3/8/2017 5:39:00 PM
cis-1,2-Dichloroethene	0.99	0.15		ppbV	1	3/8/2017 5:39:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 5:39:00 PM
Trichloroethene	8.8	1.5		ppbV	10	3/9/2017 2:01:00 AM
Vinyl chloride	< 0.15	0.15		ppbV	1	3/8/2017 5:39:00 PM
Surr: Bromofluorobenzene	98.0	70-130		%REC	10	3/9/2017 2:01:00 AM
Surr: Bromofluorobenzene	187	70-130	S	%REC	1	3/8/2017 5:39:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte, Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-005A

Client Sample ID: Duplicate  
 Tag Number: 1184.299  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY METHOD TO15		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 5:39:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	3/8/2017 5:39:00 PM
cis-1,2-Dichloroethene	3.9	0.59		ug/m3	1	3/8/2017 5:39:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 5:39:00 PM
Trichloroethene	47	8.1		ug/m3	10	3/9/2017 2:01:00 AM
Vinyl chloride	< 0.38	0.38		ug/m3	1	3/8/2017 5:39:00 PM

## NOTES:

Surrogate did not meet criteria due to severe matrix interference.

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

Data File : C:\HPCHEM\1\DATA\AO030813.D

Vial: 28

Acq On : 8 Mar 2017 5:39 pm

Operator: RJP

Sample : C1703015-005A

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:23 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.65	128	77699	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.03	114	348681	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	330439	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	387056	1.87	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	187.00%#

## Target Compounds

						Qvalue
29) cis-1,2-dichloroethene	9.46	61	122307	0.99	ppb	96
44) Trichloroethene	12.88	130	2033182	10.27	ppb	96

Data File : C:\HPCHEM\1\DATA\A0030813.D  
Acq On : 8 Mar 2017 5:39 pm  
Sample : C1703015-005A  
Misc : A227\_IUG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 10 12:38 2017

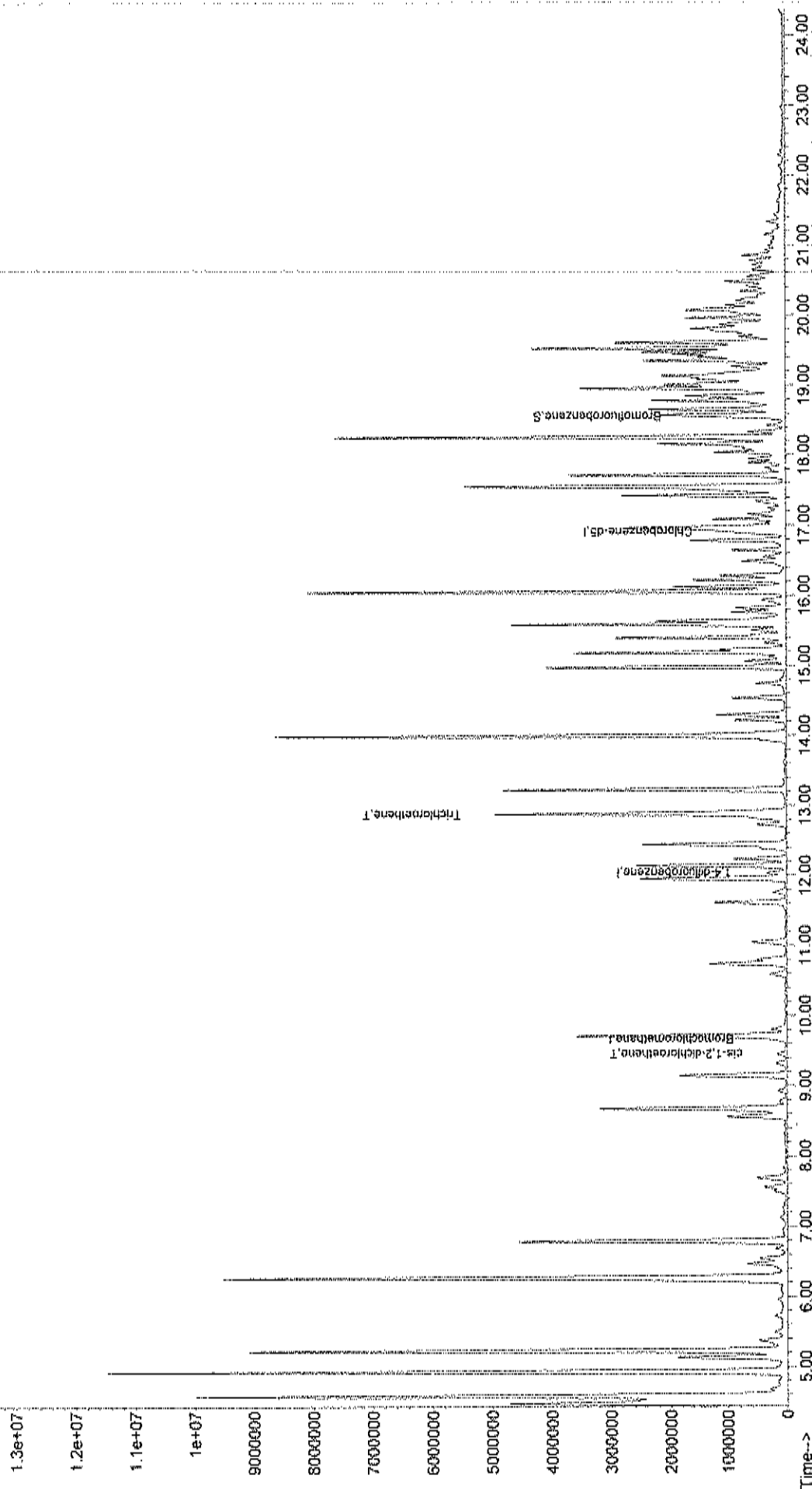
Vial: 28  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

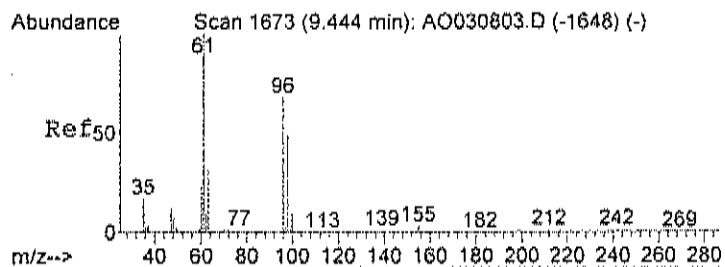
Quant Results File: A227\_IUG.RES

Method : C:\HPCHEM\1\METHODS\A227\_IUG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Mar 27 11:00:28 2017  
Response via : Initial Calibration

Abundance

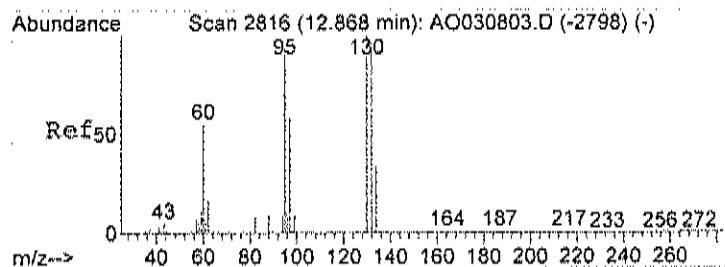
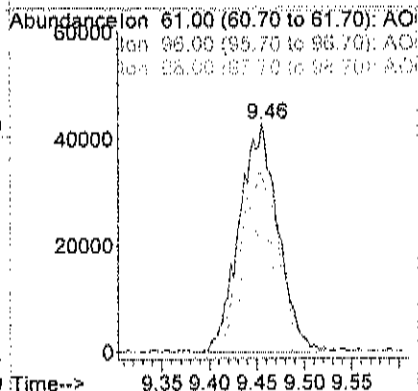
TIC: A0030813.D





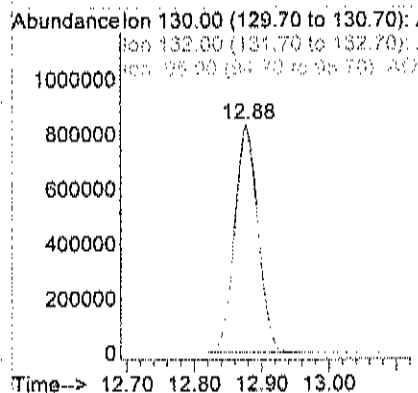
#29  
cis-1,2-dichloroethene  
Concen: 0.99 ppb  
RT: 9.46 min Scan# 1677  
Delta R.T. -0.00 min  
Lab File: AO030813.D  
Acq: 8 Mar 2017 5:39 pm

Tgt Ion	Ratio	Lower	Upper
61	100		
96	79.5	58.1	98.1
98	53.8	29.3	69.3



#44  
Trichloroethene  
Concen: 10.27 ppb  
RT: 12.88 min Scan# 2819  
Delta R.T. -0.00 min  
Lab File: AO030813.D  
Acq: 8 Mar 2017 5:39 pm

Tgt Ion	Ratio	Lower	Upper
130	100		
132	97.0	69.9	109.9
95	95.0	76.3	116.3



## Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\AO030826.D

Vial: 40

Acq On : 9 Mar 2017 2:01 am

Operator: RJP

Sample : C1703015-005A 10X

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:36 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.65	128	46525	1.00	ppb	-0.01
35) 1,4-difluorobenzene	12.03	114	198463	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	190799	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	117295	0.98	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	98.00%

## Target Compounds

44) Trichloroethene	12.87	130	99086	0.88	ppb	Qvalue 96
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Data File : C:\HPCHEM\1\DATA\AO030826.D  
Acq On : 9 Mar 2017 2:01 am  
Sample : C1703015-005A 10X  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 10 12:53 2017

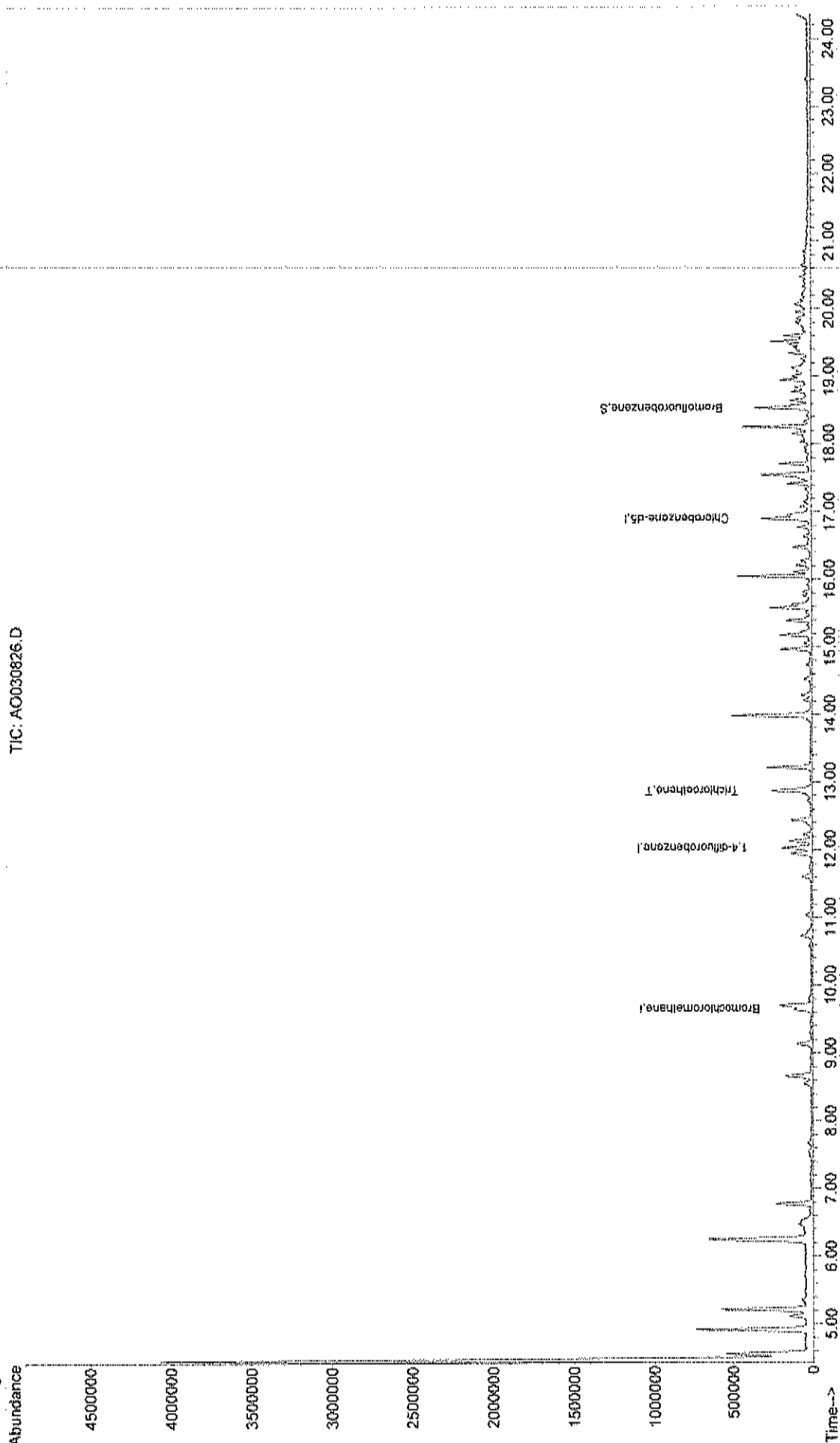
Vial: 40  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A227\_1UG.RES

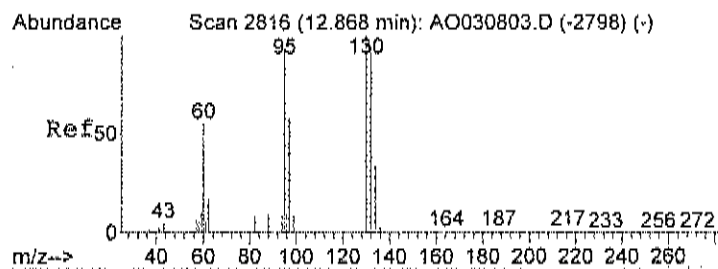
Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Mar 27 11:00:28 2017  
Response via : Initial Calibration

Abundance

TIC: AO030826.D

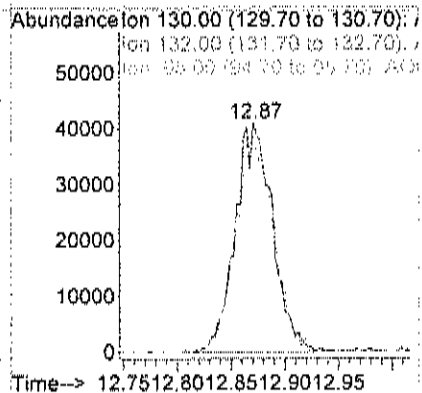






#44  
Trichloroethene  
Concen: 0.88 ppb  
RT: 12.87 min Scan# 2817  
Delta R.T. -0.01 min  
Lab File: AO030826.D  
Acq: 9 Mar 2017 2:01 am

Tgt Ion	130	Resp	99086
Ion	Ratio	Lower	Upper
130	100		
132	97.2	69.9	109.9
95	95.5	76.3	116.3



## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-006A

Client Sample ID: Ambient Air  
 Tag Number: 556.267  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
<b>FIELD PARAMETERS</b>						
			FLD			Analyst:
Lab Vacuum In	1			"Hg		3/7/2017
Lab Vacuum Out	-30			"Hg		3/7/2017
<b>1UG/M3 W/ 0.25UG/M3 CT-TCE-VC</b>						
			TO-15			Analyst: RJP
1,1-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 12:48:00 PM
Chloroethane	< 0.15	0.15		ppbV	1	3/8/2017 12:48:00 PM
cis-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 12:48:00 PM
trans-1,2-Dichloroethene	< 0.15	0.15		ppbV	1	3/8/2017 12:48:00 PM
Trichloroethene	< 0.040	0.040		ppbV	1	3/8/2017 12:48:00 PM
Vinyl chloride	< 0.040	0.040		ppbV	1	3/8/2017 12:48:00 PM
Surr: Bromofluorobenzene	87.0	70-130		%REC	1	3/8/2017 12:48:00 PM

Qualifiers: \*\* Quantitation Limit  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 JN Non-routine analyte. Quantitation estimated.  
 S Spike Recovery outside accepted recovery limits  
 , Results reported are not blank corrected  
 E Estimated Value above quantitation range  
 J Analyte detected below quantitation limit  
 ND Not Detected at the Limit of Detection

## Centek Laboratories, LLC

Date: 27-Mar-17

CLIENT: LaBella Associates, P.C.  
 Lab Order: C1703015  
 Project: 691 and 705 St Paul St  
 Lab ID: C1703015-006A

Client Sample ID: Ambient Air  
 Tag Number: 556.267  
 Collection Date: 3/3/2017  
 Matrix: AIR

Analyses	Result	**Limit	Qual	Units	DF	Date Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC		TO-15		Analyst: RJP		
1,1-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 12:48:00 PM
Chloroethane	< 0.40	0.40		ug/m3	1	3/8/2017 12:48:00 PM
cis-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 12:48:00 PM
trans-1,2-Dichloroethene	< 0.59	0.59		ug/m3	1	3/8/2017 12:48:00 PM
Trichloroethene	< 0.21	0.21		ug/m3	1	3/8/2017 12:48:00 PM
Vinyl chloride	< 0.10	0.10		ug/m3	1	3/8/2017 12:48:00 PM

Qualifiers:	**	Quantitation Limit	.	Results reported are not blank corrected
	B	Analyte detected in the associated Method Blank	E	Estimated Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limit
	JN	Non-routine analyte. Quantitation estimated.	ND	Not Detected at the Limit of Detection
	S	Spike Recovery outside accepted recovery limits		

Data File : C:\HPCHEM\1\DATA\AO030806.D

Vial: 21

Acq On : 8 Mar 2017 12:48 pm

Operator: RJP

Sample : C1703015-006A

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:16 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.65	128	47960	1.00	ppb	-0.01
35) 1,4-difluorobenzene	12.03	114	186316	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	173836	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	94823	0.87	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	87.00%

## Target Compounds

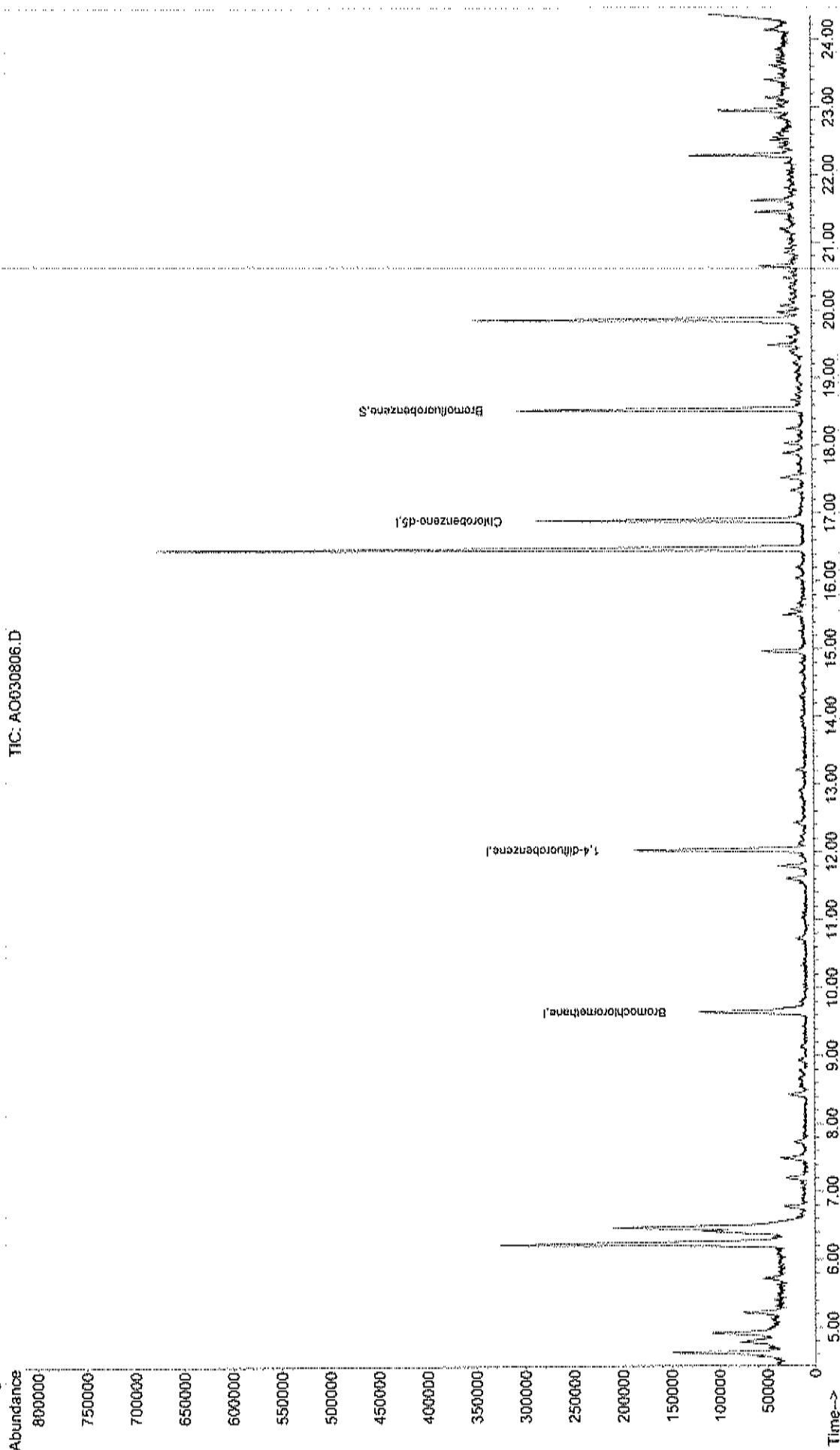
Qvalue

Data File : C:\HPCHEM\1\DATA\AO030806.D  
Acq On : 8 Mar 2017 12:48 pm  
Sample : C1703015-006A  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 10 12:13 2017

Vial: 21  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00  
Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Mar 27 11:00:28 2017  
Response via : Initial Calibration

TIC: AO030806.D



**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**STANDARDS DATA**

**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**INITIAL CALIBRATION**

## Response Factor Report MSD #1

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Mon Feb 27 19:25:53 2017  
 Response via : Initial Calibration

## Calibration Files

0.04 =AO022712.D 0.10 =AO022711.D 0.15 =AO022710.D  
 0.30 =AO022709.D 0.50 =AO022707.D 0.75 =AO022706.D

Compound	0.04	0.10	0.15	0.30	0.50	0.75	Avg	%RSD
1) I Bromochloromethane				ISTD				
2) T Propylene			1.077	1.105	1.080	0.984	1.042	5.29
3) T Freon 12			9.752	9.472	9.687	8.524	9.266	6.10
4) T Chloromethane			1.751	1.749	1.818	1.443	1.659	8.01
5) T Freon 114			8.517	7.919	8.532	6.688	7.921	8.76
6) T Vinyl Chloride	3.341	2.514	2.380	2.148	2.278	1.786	2.301	18.51
7) T Butane			2.756	2.539	2.580	1.994	2.394	10.89
8) T 1,3-butadiene			1.629	1.592	1.691	1.273	1.535	8.91
9) T Bromomethane			3.134	2.822	3.057	2.237	2.831	10.14
10) T Chloroethane			1.080	1.119	1.066	0.854	1.014	8.78
11) T Ethanol			0.839	0.898	0.844	0.579	0.748	14.21
12) T Acrolein			0.771	0.819	0.880	0.592	0.765	11.09
13) T Vinyl Bromide			2.936	2.908	3.017	2.090	2.844	11.03
14) T Freon 11			1.189	1.141	1.175	0.910	1.109	E1 7.97
15) T Acetone			0.963	0.913	0.914	0.735	0.882	7.72
16) T Pentane			1.085	1.027	1.052	0.989	1.010	4.57
17) T Isopropyl alcoh			3.908	3.987	3.323	2.514	3.221	15.59
18) T 1,1-dichloroeth			1.301	1.299	1.276	1.251	1.282	2.86
19) T Freon 113			3.735	3.519	3.538	3.425	3.515	3.18
20) t t-Butyl alcohol			2.463	2.764	2.516	2.361	2.527	6.16
21) T Methylene chlor			1.543	1.399	1.386	1.274	1.335	7.91
22) T Allyl chloride			1.127	1.004	1.070	1.093	1.090	4.30
23) T Carbon disulfid			4.409	4.088	4.065	3.870	3.995	5.01
24) T trans-1,2-dichl			1.946	1.773	1.849	1.747	1.805	4.67
25) T methyl tert-but			2.987	2.995	3.154	3.000	3.166	5.59
26) T 1,1-dichloroeth			2.191	2.383	2.517	2.330	2.386	4.14
27) T Vinyl acetate			2.646	1.955	2.250	2.245	2.361	9.31
28) T Methyl Ethyl Ke			0.522	0.585	0.559	0.561	0.565	4.43
29) T cis-1,2-dichlor			1.554	1.596	1.504	1.534	1.586	3.91
30) T Hexane			1.116	1.103	1.282	1.291	1.306	10.63
31) T Ethyl acetate			2.660	2.650	2.881	2.946	2.971	7.92
32) T Chloroform			4.123	3.930	4.220	4.052	4.078	2.95
33) T Tetrahydrofuran			0.820	0.841	0.811	0.830	0.865	6.76
34) T 1,2-dichloroeth			2.785	2.589	2.684	2.703	2.716	3.35
35) I 1,4-difluorobenzene			ISTD					
36) T 1,1,1-trichloro			1.256	1.165	1.160	1.101	1.135	5.19
37) T Cyclohexane			0.315	0.350	0.352	0.343	0.341	3.81
38) T Carbon tetrachl	2.731	1.619	1.596	1.468	1.467	1.428	1.580	26.26
39) T Benzene			0.955	0.876	0.891	0.875	0.891	3.00
40) T Methyl methacry			0.247	0.283	0.298	0.315	0.312	11.04
41) T 1,4-dioxane			0.163	0.185	0.163	0.165	0.172	4.56
42) T 2,2,4-trimethyl			1.039	1.054	1.153	1.208	1.185	7.86
43) T Heptane			0.320	0.302	0.349	0.377	0.360	9.29
44) T Trichloroethene	0.901	0.552	0.530	0.522	0.507	0.534	0.568	20.69
45) T 1,2-dichloropro			0.334	0.299	0.303	0.300	0.306	4.13
46) T Bromodichlorome			1.167	1.117	1.098	1.099	1.088	3.91
47) T cis-1,3-dichlor			0.508	0.522	0.500	0.520	0.524	2.94
48) T trans-1,3-dichl			0.508	0.480	0.491	0.510	0.513	4.14
49) T 1,1,2-trichloro			0.508	0.512	0.480	0.495	0.488	3.78
50) I Chlorobenzene-d5			ISTD					
51) T Toluene			0.731	0.696	0.720	0.748	0.760	6.06



## Response Factor Report MSD #1

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Mon Feb 27 19:25:53 2017  
 Response via : Initial Calibration

## Calibration Files

0.04 =AO022712.D 0.10 =AO022711.D 0.15 =AO022710.D  
 0.30 =AO022709.D 0.50 =AO022707.D 0.75 =AO022706.D

	Compound	0.04	0.10	0.15	0.30	0.50	0.75	Avg	%RSD
52) T	Methyl Isobutyl				0.639	0.577	0.650	0.683	10.53
53) T	Dibromochlorome			1.456	1.426	1.381	1.463	1.439	2.27
54) T	Methyl Butyl Ke				0.598	0.530	0.565	0.629	12.85
55) T	1,2-dibromoetha			0.876	0.906	0.897	0.935	0.930	3.67
56) T	Tetrachloroethy			0.733	0.649	0.664	0.673	0.691	4.07
57) T	Chlorobenzene			1.201	1.163	1.191	1.257	1.238	3.85
58) T	Ethylbenzene			1.575	1.533	1.604	1.709	1.745	10.00
59) T	m&p-xylene			1.361	1.437	1.526	1.555	1.665	13.45
60) T	Nonane			0.575	0.525	0.631	0.714	0.695	15.40
61) T	Styrene			0.878	0.830	0.888	0.934	0.969	11.14
62) T	Bromoform			1.185	1.256	1.267	1.324	1.307	5.05
63) T	o-xylene			1.302	1.343	1.444	1.649	1.612	13.78
64) T	Cumene			1.936	2.017	2.046	2.238	2.227	9.55
65) S	Bromofluorobenz	0.556	0.553	0.594	0.606	0.637	0.653	0.627	7.46
66) T	1,1,2,2-tetrach			1.230	1.222	1.208	1.298	1.256	2.60
67) T	Propylbenzene			0.530	0.540	0.567	0.628	0.625	11.41
68) T	2-Chlorotoluene			0.495	0.495	0.479	0.533	0.528	7.48
69) T	4-ethyltoluene			1.639	1.780	1.938	2.081	2.062	12.97
70) T	1,3,5-trimethyl			1.480	1.671	1.758	2.015	1.940	14.05
71) T	1,2,4-trimethyl			1.461	1.454	1.406	1.613	1.642	12.06
72) T	1,3-dichloroben			1.145	1.298	1.207	1.305	1.293	6.57
73) T	benzyl chloride			1.018	1.045	1.035	1.114	1.119	7.00
74) T	1,4-dichloroben			0.933	1.095	1.055	1.169	1.136	9.07
75) T	1,2,3-trimethyl			1.465	1.505	1.526	1.664	1.713	12.48
76) T	1,2-dichloroben			1.277	1.263	1.249	1.325	1.297	2.62
77) T	1,2,4-trichloro			0.465	0.433	0.402	0.434	0.468	10.49
78) T	Naphthalene			1.130	1.213	1.033	1.141	1.258	15.12
79) T	Hexachloro-1,3-			1.692	1.457	1.256	1.278	1.359	10.97

Data File : C:\HPCHEM\1\DATA\AO022703.D

Vial: 1

Acq On : 27 Feb 2017 12:01 pm

Operator: RJP

Sample : A1UG\_2.0

Inst : MSD #1

Misc : A120\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 13:34:31 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.66	128	34488	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.04	114	147413	1.00	ppb	0.00
50) Chlorobenzene-d5	16.91	117	118307	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	80069	1.01	ppb	0.00
Spiked Amount	1.000	Range	70 ~ 130	Recovery	=	101.00%

## Target Compounds

						Qvalue
2) Propylene	4.56	41	67283m	2.01	ppb	
3) Freon 12	4.63	85	637506	1.94	ppb	98
4) Chloromethane	4.80	50	113936m	1.90	ppb	
5) Freon 114	4.90	85	558550	1.90	ppb	98
6) Vinyl Chloride	5.03	62	151882	2.04	ppb	91
7) Butane	5.23	43	159751	1.86	ppb	89
8) 1,3-butadiene	5.17	39	107879m	1.87	ppb	
9) Bromomethane	5.45	94	194464	1.83	ppb	95
10) Chloroethane	5.63	64	67257	1.91	ppb	89
11) Ethanol	5.73	45	47585m	2.00	ppb	
12) Acrolein	6.09	56	52048	1.83	ppb	91
13) Vinyl Bromide	5.98	106	202045	1.91	ppb	96
14) Freon 11	6.42	101	776050	1.88	ppb	98
15) Acetone	6.23	58	61266	1.98	ppb	# 76
16) Pentane	6.79	42	68027	1.94	ppb	# 41
17) Isopropyl alcohol	6.47	45	204229	1.87	ppb	# 100
18) 1,1-dichloroethene	7.10	96	89203	1.96	ppb	99
19) Freon 113	7.50	101	238988	1.99	ppb	96
20) t-Butyl alcohol	7.12	59	182177	2.01	ppb	# 91
21) Methylene chloride	7.21	84	87576	1.93	ppb	93
22) Allyl chloride	7.34	41	78216	2.02	ppb	94
23) Carbon disulfide	7.56	76	261881	1.90	ppb	82
24) trans-1,2-dichloroethene	8.25	61	123974	2.01	ppb	93
25) methyl tert-butyl ether	8.55	73	236087	2.07	ppb	91
26) 1,1-dichloroethane	8.48	63	167678	1.96	ppb	96
27) Vinyl acetate	8.63	43	178454m	2.19	ppb	
28) Methyl Ethyl Ketone	8.93	72	40336	2.07	ppb	# 1
29) cis-1,2-dichloroethene	9.46	61	115901	2.07	ppb	93
30) Hexane	9.71	57	102300	2.08	ppb	88
31) Ethyl acetate	9.72	43	225431	2.06	ppb	98
32) Chloroform	9.81	83	284156	1.92	ppb	96
33) Tetrahydrofuran	10.32	42	65779	2.13	ppb	# 70
34) 1,2-dichloroethane	10.73	62	190579	1.99	ppb	96
36) 1,1,1-trichloroethane	11.05	97	326280	1.94	ppb	97
37) Cyclohexane	11.96	56	104270	2.12	ppb	88
38) Carbon tetrachloride	11.80	117	402678	1.93	ppb	98
39) Benzene	11.62	78	260440	1.99	ppb	94
40) Methyl methacrylate	13.10	41	103894m	2.12	ppb	
41) 1,4-dioxane	12.89	58	51011	2.11	ppb	# 33
42) 2,2,4-trimethylpentane	12.92	57	374982	2.06	ppb	88
43) Heptane	13.23	43	117429	2.17	ppb	85
44) Trichloroethene	12.88	130	157508	2.00	ppb	92
45) 1,2-dichloropropane	12.60	63	89213	2.05	ppb	99

(#)=qualifier out of range (m)=manual integration

AO022703.D A227\_1UG.M

Mon Mar 06 09:22:53 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO022703.D

Vial: 1

Acq On : 27 Feb 2017 12:01 pm

Operator: RJP

Sample : A1UG\_2.0

Inst : MSD #1

Misc : A120\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 13:34:31 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

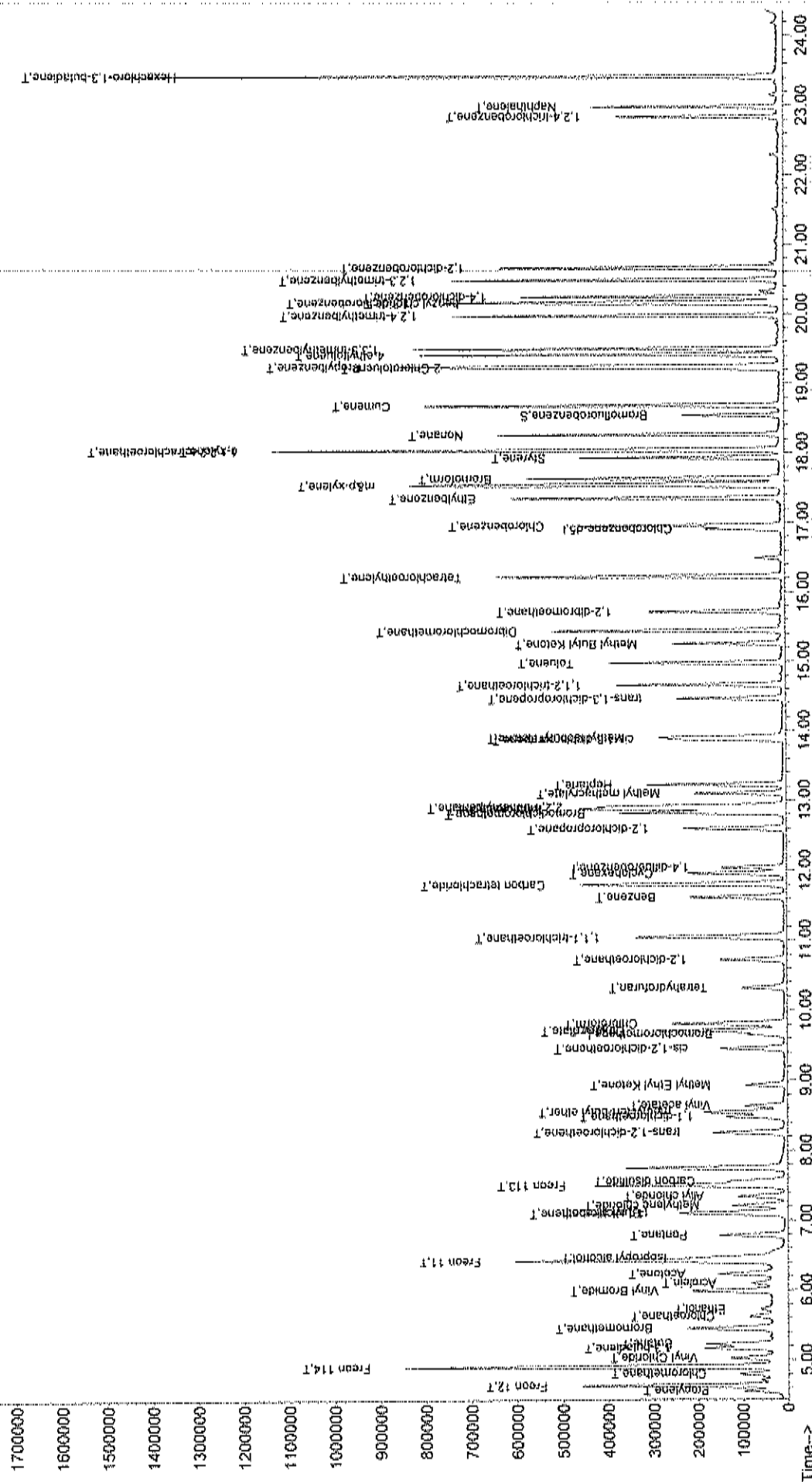
Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.83	83	313826	1.95	ppb	98
47) cis-1,3-dichloropropene	13.88	75	161710	2.10	ppb	95
48) trans-1,3-dichloropropene	14.47	75	160601	2.11	ppb	95
49) 1,1,2-trichloroethane	14.66	97	140491	1.98	ppb	97
51) Toluene	14.98	92	198866	2.09	ppb	87
52) Methyl Isobutyl Ketone	13.91	43	185015	1.95	ppb	96
53) Dibromochloromethane	15.44	129	352272	2.00	ppb	93
54) Methyl Butyl Ketone	15.26	43	176699	1.90	ppb	98
55) 1,2-dibromoethane	15.72	107	231834	1.96	ppb	97
56) Tetrachloroethylene	16.22	164	170159	2.04	ppb	97
57) Chlorobenzene	16.95	112	307845	1.98	ppb	97
58) Ethylbenzene	17.35	91	482120	2.21	ppb	98
59) m&p-xylene	17.54	91	936602	4.26	ppb	99
60) Nonane	18.26	43	197031	2.13	ppb	95
61) Styrene	17.93	104	269222	2.27	ppb	92
62) Bromoform	17.63	173	327396	1.96	ppb	99
63) o-xylene	18.05	91	452697	2.16	ppb	95
64) Cumene	18.68	105	604876	2.15	ppb	99
66) 1,1,2,2-tetrachloroethane	18.04	83	304924	1.95	ppb	95
67) Propylbenzene	19.25	120	170201	2.12	ppb	# 1
68) 2-Chlorotoluene	19.22	126	142423	2.22	ppb	# 1
69) 4-ethyltoluene	19.41	105	576004	2.16	ppb	98
70) 1,3,5-trimethylbenzene	19.49	105	536902	2.07	ppb	98
71) 1,2,4-trimethylbenzene	19.97	105	464766	2.20	ppb	93
72) 1,3-dichlorobenzene	20.16	146	336426	2.04	ppb	98
73) benzyl chloride	20.14	91	289019	1.97	ppb	89
74) 1,4-dichlorobenzene	20.25	146	295119	2.02	ppb	96
75) 1,2,3-trimethylbenzene	20.49	105	489968	2.13	ppb	95
76) 1,2-dichlorobenzene	20.66	146	318561	1.95	ppb	98
77) 1,2,4-trichlorobenzene	22.83	180	129235	1.98	ppb	95
78) Naphthalene	22.97	128	376207	2.05	ppb	94
79) Hexachloro-1,3-butadiene	23.41	225	314561	1.80	ppb	98

-----  
 (#) = qualifier out of range (m) = manual integration (+) = signals summed  
 AO022703.D A227\_1UG.M Mon Mar 06 09:22:54 2017 MSD1



Data File : C:\HPCHEM\1\DATA\AO022704.D

Vial: 2

Acq On : 27 Feb 2017 12:42 pm

Operator: RJP

Sample : A1UG\_1.5

Inst : MSD #1

Misc : A120\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 13:37:02 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	9.65	128	36100	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.04	114	151427	1.00	ppb	0.00
50) Chlorobenzene-d5	16.91	117	120601	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	81420	1.00	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	100.00%

## Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Propylene	4.55	41	55806	1.59	ppb	89
3) Freon 12	4.63	85	447890	1.30	ppb	98
4) Chloromethane	4.80	50	80196m $\beta$	1.27	ppb	
5) Freon 114	4.90	85	382883	1.24	ppb	94
6) Vinyl Chloride	5.02	62	101642	1.30	ppb	89
7) Butane	5.22	43	113187	1.26	ppb	90
8) 1,3-butadiene	5.17	39	78428m $\beta$	1.30	ppb	
9) Bromomethane	5.45	94	143623	1.29	ppb	94
10) Chloroethane	5.62	64	50578	1.37	ppb	# 84
11) Ethanol	5.73	45	36233	1.45	ppb	# 84
12) Acrolein	6.10	56	40470	1.36	ppb	86
13) Vinyl Bromide	5.99	106	155521	1.41	ppb	94
14) Freon 11	6.42	101	578839	1.34	ppb	100
15) Acetone	6.23	58	45684	1.41	ppb	# 62
16) Pentane	6.78	42	52005	1.42	ppb	# 36
17) Isopropyl alcohol	6.48	45	159740	1.40	ppb	# 100
18) 1,1-dichloroethene	7.09	96	68749	1.44	ppb	96
19) Freon 113	7.50	101	181865	1.44	ppb	97
20) t-Butyl alcohol	7.12	59	129743	1.37	ppb	# 87
21) Methylene chloride	7.22	84	65431	1.38	ppb	92
22) Allyl chloride	7.34	41	59816	1.48	ppb	92
23) Carbon disulfide	7.56	76	206115	1.43	ppb	83
24) trans-1,2-dichloroethene	8.26	61	94649	1.47	ppb	94
25) methyl tert-butyl ether	8.55	73	175245	1.47	ppb	90
26) 1,1-dichloroethane	8.48	63	127822	1.43	ppb	95
27) Vinyl acetate	8.64	43	130963m $\beta$	1.54	ppb	
28) Methyl Ethyl Ketone	8.93	72	30056	1.48	ppb	# 83
29) cis-1,2-dichloroethene	9.46	61	84779	1.45	ppb	92
30) Hexane	9.72	57	76403	1.48	ppb	87
31) Ethyl acetate	9.71	43	167003	1.46	ppb	98
32) Chloroform	9.81	83	211706	1.36	ppb	98
33) Tetrahydrofuran	10.33	42	48492	1.50	ppb	# 70
34) 1,2-dichloroethane	10.72	62	145890	1.46	ppb	98
36) 1,1,1-trichloroethane	11.05	97	242355	1.40	ppb	96
37) Cyclohexane	11.96	56	77508	1.53	ppb	88
38) Carbon tetrachloride	11.80	117	304694	1.42	ppb	100
39) Benzene	11.62	78	198866	1.48	ppb	94
40) Methyl methacrylate	13.10	41	74232	1.47	ppb	# 82
41) 1,4-dioxane	12.90	58	39753	1.60	ppb	# 30
42) 2,2,4-trimethylpentane	12.92	57	281936	1.51	ppb	88
43) Heptane	13.23	43	86145	1.55	ppb	86
44) Trichloroethene	12.88	130	119387	1.48	ppb	97
45) 1,2-dichloropropane	12.60	63	66632	1.49	ppb	98

(#)=qualifier out of range (m)=manual integration

AO022704.D A227\_1UG.M

Mon Mar 06 09:22:57 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO022704.D

Vial: 2

Acq On : 27 Feb 2017 12:42 pm

Operator: RJP

Sample : A1UG\_1.5

Inst : MSD #1

Misc : A120\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 13:37:02 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

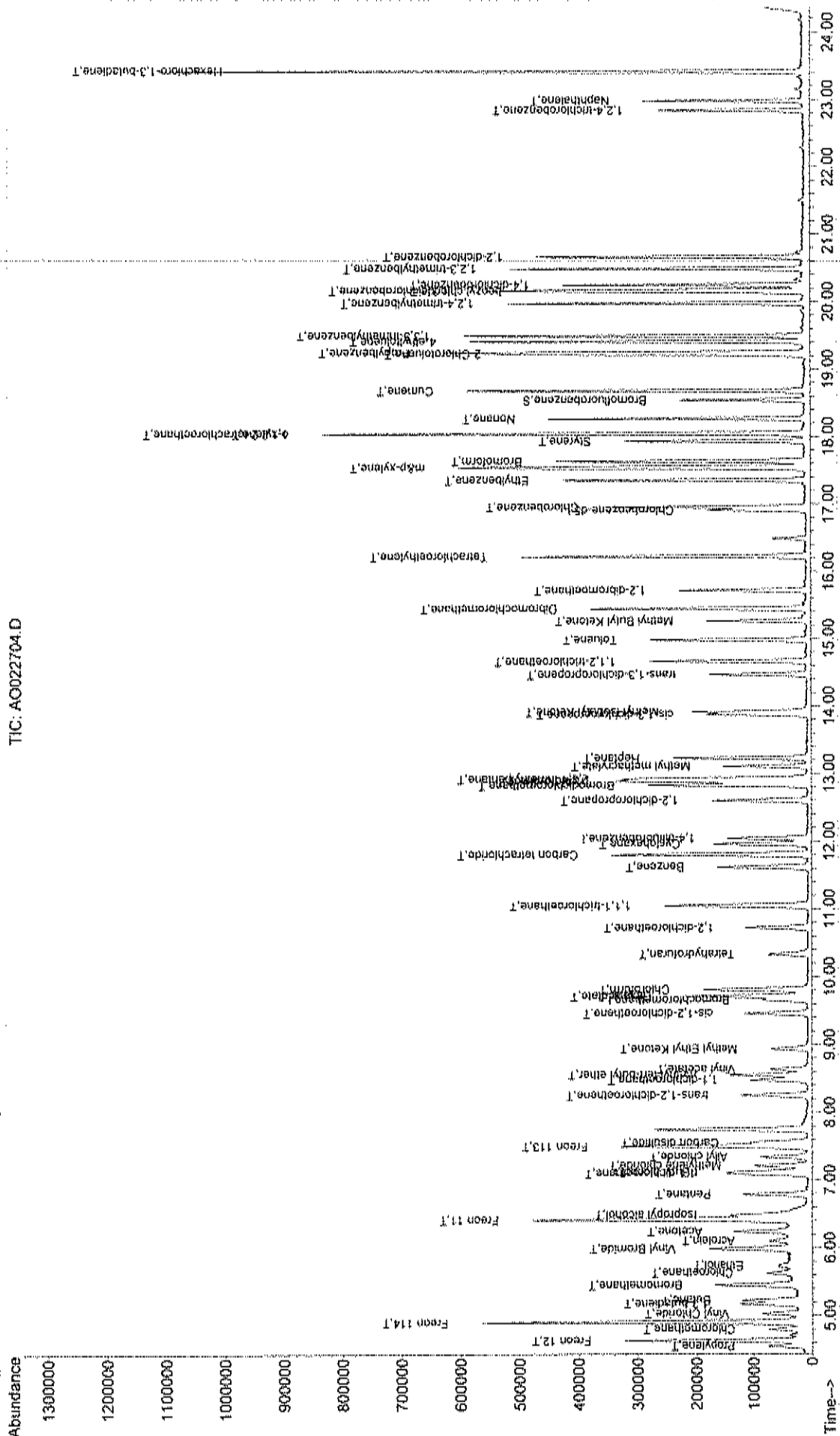
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.83	83	233846	1.41	ppb	96
47) cis-1,3-dichloropropene	13.87	75	119377	1.51	ppb	96
48) trans-1,3-dichloropropene	14.47	75	116355	1.49	ppb	96
49) 1,1,2-trichloroethane	14.67	97	104823	1.44	ppb	100
51) Toluene	14.98	92	143255	1.48	ppb	86
52) Methyl Isobutyl Ketone	13.91	43	136808	1.41	ppb	96
53) Dibromochloromethane	15.45	129	260970	1.46	ppb	93
54) Methyl Butyl Ketone	15.26	43	127927	1.35	ppb	98
55) 1,2-dibromoethane	15.72	107	171510	1.42	ppb	96
56) Tetrachloroethylene	16.22	164	126330	1.49	ppb	97
57) Chlorobenzene	16.95	112	230234	1.46	ppb	96
58) Ethylbenzene	17.35	91	344400	1.55	ppb	98
59) m&p-xylene	17.54	91	674977	3.01	ppb	100
60) Nonane	18.26	43	139459	1.48	ppb	96
61) Styrene	17.93	104	197650	1.63	ppb	88
62) Bromoform	17.63	173	244288	1.43	ppb	98
63) o-xylene	18.05	91	319499	1.49	ppb	96
64) Cumene	18.68	105	436193	1.52	ppb	99
66) 1,1,2,2-tetrachloroethane	18.04	83	230628	1.44	ppb	97
67) Propylbenzene	19.24	120	124849	1.52	ppb #	1
68) 2-Chlorotoluene	19.22	126	98971	1.52	ppb #	1
69) 4-ethyltoluene	19.41	105	417512	1.54	ppb	99
70) 1,3,5-trimethylbenzene	19.49	105	388020	1.47	ppb	97
71) 1,2,4-trimethylbenzene	19.97	105	329611	1.53	ppb	94
72) 1,3-dichlorobenzene	20.16	146	244983m	1.46	ppb	
73) benzyl chloride	20.14	91	215675	1.44	ppb	92
74) 1,4-dichlorobenzene	20.25	146	218869m	1.47	ppb	
75) 1,2,3-trimethylbenzene	20.49	105	346422	1.48	ppb	94
76) 1,2-dichlorobenzene	20.66	146	240399	1.44	ppb	97
77) 1,2,4-trichlorobenzene	22.83	180	95449	1.44	ppb	96
78) Naphthalene	22.97	128	259809	1.39	ppb	95
79) Hexachloro-1,3-butadiene	23.41	225	234742	1.32	ppb	100

Data File : C:\HPCHEM\1\DATA\AO022704.D  
Acq On : 27 Feb 2017 12:42 pm  
Sample : A1UG\_1.5  
Misc : A120\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Feb 27 14:26 2017

Vial: 2  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00  
Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Feb 27 19:25:53 2017  
Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

TIC: AO022704.D



Data File : C:\HPCHEM\1\DATA\AO022705.D

Vial: 3

Acq On : 27 Feb 2017 1:22 pm

Operator: RJP

Sample : A1UG\_1.25

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 14:21:04 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev.(Min)
1) Bromochloromethane	9.66	128	33434	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.04	114	145107	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	118466	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	77962	0.98	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	98.00%

## Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Propylene	4.55	41	45886	1.42	ppb	82
3) Freon 12	4.63	85	405753	1.27	ppb	97
4) Chloromethane	4.79	50	71614m	1.23	ppb	
5) Freon 114	4.89	85	355820	1.25	ppb	98
6) Vinyl Chloride	5.02	62	96809	1.34	ppb	90
7) Butane	5.22	43	106701	1.28	ppb	93
8) 1,3-butadiene	5.17	39	68251m	1.22	ppb	
9) Bromomethane	5.44	94	127291	1.24	ppb	94
10) Chloroethane	5.63	64	45045	1.32	ppb	89
11) Ethanol	5.74	45	30805	1.33	ppb	82
12) Acrolein	6.09	56	34273	1.25	ppb	88
13) Vinyl Bromide	5.99	106	129727	1.27	ppb	96
14) Freon 11	6.42	101	478226	1.19	ppb	99
15) Acetone	6.23	58	38074	1.27	ppb	# 67
16) Pentane	6.78	42	43027	1.26	ppb	# 40
17) Isopropyl alcohol	6.47	45	129586	1.22	ppb	# 100
18) 1,1-dichloroethene	7.10	96	56144	1.27	ppb	97
19) Freon 113	7.50	101	149371	1.28	ppb	97
20) t-Butyl alcohol	7.11	59	112735	1.28	ppb	# 90
21) Methylene chloride	7.21	84	55538	1.26	ppb	94
22) Allyl chloride	7.33	41	47631	1.27	ppb	93
23) Carbon disulfide	7.56	76	166949	1.25	ppb	83
24) trans-1,2-dichloroethene	8.25	61	78916	1.32	ppb	94
25) methyl tert-butyl ether	8.55	73	142379	1.29	ppb	88
26) 1,1-dichloroethane	8.48	63	103193	1.25	ppb	93
27) Vinyl acetate	8.64	43	102981m	1.31	ppb	
28) Methyl Ethyl Ketone	8.94	72	25140	1.33	ppb	# 1
29) cis-1,2-dichloroethene	9.45	61	69788	1.29	ppb	94
30) Hexane	9.72	57	59417	1.24	ppb	90
31) Ethyl acetate	9.72	43	135614	1.28	ppb	99
32) Chloroform	9.81	83	176976	1.23	ppb	100
33) Tetrahydrofuran	10.33	42	39561	1.32	ppb	76
34) 1,2-dichloroethane	10.72	62	120237	1.30	ppb	97
36) 1,1,1-trichloroethane	11.05	97	199154	1.20	ppb	98
37) Cyclohexane	11.95	56	62283	1.28	ppb	88
38) Carbon tetrachloride	11.80	117	247922	1.21	ppb	97
39) Benzene	11.61	78	160434	1.24	ppb	94
40) Methyl methacrylate	13.10	41	61746	1.28	ppb	# 84
41) 1,4-dioxane	12.91	58	32323	1.36	ppb	# 29
42) 2,2,4-trimethylpentane	12.92	57	228477	1.28	ppb	88
43) Heptane	13.22	43	69030	1.29	ppb	81
44) Trichloroethene	12.88	130	98431	1.27	ppb	96
45) 1,2-dichloropropane	12.59	63	54586	1.28	ppb	99

(#)=qualifier out of range (m)=manual integration

AO022705.D A227\_1UG.M

Mon Mar 06 09:23:01 2017

MSD1

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Data File : C:\HPCHEM\1\DATA\AO022705.D

Vial: 3

Acq On : 27 Feb 2017 1:22 pm

Operator: RJP

Sample : A1UG\_1.25

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 14:21:04 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

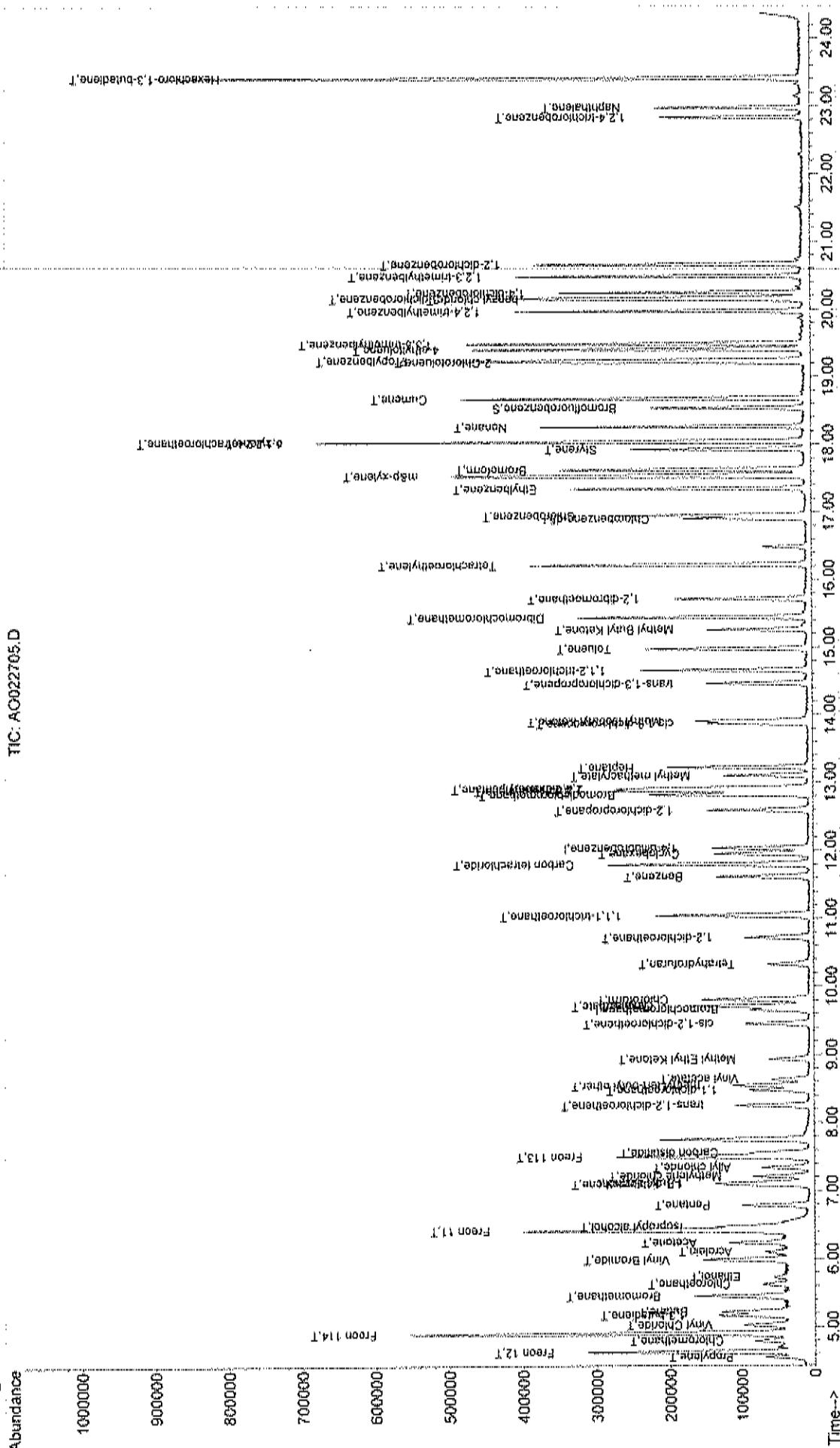
DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.83	83	191263	1.21	ppb	96
47) cis-1,3-dichloropropene	13.87	75	96035	1.27	ppb	95
48) trans-1,3-dichloropropene	14.46	75	96628	1.29	ppb	99
49) 1,1,2-trichloroethane	14.66	97	85391	1.23	ppb	99
51) Toluene	14.97	92	115953	1.22	ppb	87
52) Methyl Isobutyl Ketone	13.91	43	105820	1.11	ppb	96
53) Dibromochloromethane	15.44	129	209627	1.19	ppb	92
54) Methyl Butyl Ketone	15.26	43	99894	1.07	ppb	96
55) 1,2-dibromoethane	15.72	107	140106	1.18	ppb	98
56) Tetrachloroethylene	16.23	164	101954	1.22	ppb	98
57) Chlorobenzene	16.95	112	188047	1.21	ppb	97
58) Ethylbenzene	17.35	91	270529	1.24	ppb	97
59) m&p-xylene	17.54	91	539522	2.45	ppb	100
60) Nonane	18.26	43	113508	1.23	ppb	97
61) Styrene	17.93	104	150564	1.27	ppb	89
62) Bromoform	17.63	173	197222	1.18	ppb	98
63) o-xylene	18.04	91	258358	1.23	ppb	95
64) Cumene	18.68	105	344068	1.22	ppb	98
66) 1,1,2,2-tetrachloroethane	18.04	83	186037	1.19	ppb	94
67) Propylbenzene	19.24	120	98861	1.23	ppb	# 1
68) 2-Chlorotoluene	19.21	126	81240	1.27	ppb	# 1
69) 4-ethyltoluene	19.41	105	329649	1.24	ppb	99
70) 1,3,5-trimethylbenzene	19.49	105	313638	1.21	ppb	99
71) 1,2,4-trimethylbenzene	19.97	105	260242	1.23	ppb	94
72) 1,3-dichlorobenzene	20.16	146	192000m	1.16	ppb	
73) benzyl chloride	20.14	91	174719	1.19	ppb	91
74) 1,4-dichlorobenzene	20.24	146	178085m	1.22	ppb	
75) 1,2,3-trimethylbenzene	20.49	105	265866	1.15	ppb	92
76) 1,2-dichlorobenzene	20.66	146	191377	1.17	ppb	98
77) 1,2,4-trichlorobenzene	22.83	180	71785	1.10	ppb	93
78) Naphthalene	22.97	128	204556m	1.11	ppb	
79) Hexachloro-1,3-butadiene	23.41	225	191140	1.09	ppb	99

Data File : C:\HPCHEM\1\DATA\AO022705.D  
 Vial: 3  
 Acq On : 27 Feb 2017 1:22 pm  
 Operator: RJP  
 Sample : ALUG\_1.25  
 Inst : MSD #1  
 Misc : A227\_1UG  
 Multiplr: 1.00  
 MS Integration Params: RTEINT.P  
 Quant Time: Feb 27 14:27 2017  
 Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
 Title : TO-15 VOA Standards for 5 point calibration  
 Last Update : Mon Feb 27 19:25:53 2017  
 Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

TIC: AO022705.D



Data File : C:\HPCHEM\1\DATA\AO022706.D

Vial: 4

Acq On : 27 Feb 2017 2:01 pm

Operator: RJP

Sample : A1UG\_0.75

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 14:27:59 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.66	128	34548	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.04	114	139180	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	114597	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	74856	0.97	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	97.00%

## Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Propylene	4.56	41	25499	0.76	ppb	77
3) Freon 12	4.64	85	220863	0.67	ppb	98
4) Chloromethane	4.81	50	37377m	0.62	ppb	
5) Freon 114	4.90	85	173292	0.59	ppb	93
6) Vinyl Chloride	5.03	62	46269	0.62	ppb	93
7) Butane	5.23	43	51654	0.60	ppb	91
8) 1,3-butadiene	5.18	39	32992m	0.57	ppb	
9) Bromomethane	5.45	94	57960	0.54	ppb	94
10) Chloroethane	5.63	64	22140	0.63	ppb	90
11) Ethanol	5.74	45	14995	0.63	ppb	82
12) Acrolein	6.11	56	15338	0.54	ppb	89
13) Vinyl Bromide	5.99	106	54158	0.51	ppb	91
14) Freon 11	6.42	101	235743	0.57	ppb	99
15) Acetone	6.25	58	19048	0.61	ppb	# 64
16) Pentane	6.80	42	25624	0.73	ppb	# 51
17) Isopropyl alcohol	6.49	45	65135	0.60	ppb	# 100
18) 1,1-dichloroethene	7.10	96	32415	0.71	ppb	99
19) Freon 113	7.50	101	88734	0.74	ppb	96
20) t-Butyl alcohol	7.14	59	61182	0.67	ppb	# 86
21) Methylene chloride	7.21	84	33001	0.72	ppb	93
22) Allyl chloride	7.34	41	28333	0.73	ppb	93
23) Carbon disulfide	7.56	76	100285	0.73	ppb	81
24) trans-1,2-dichloroethene	8.26	61	45269	0.73	ppb	92
25) methyl tert-butyl ether	8.55	73	77736	0.68	ppb	91
26) 1,1-dichloroethane	8.48	63	60381	0.71	ppb	94
27) Vinyl acetate	8.64	43	58164m	0.71	ppb	
28) Methyl Ethyl Ketone	8.95	72	14528	0.75	ppb	# 76
29) cis-1,2-dichloroethene	9.45	61	39739	0.71	ppb	90
30) Hexane	9.72	57	33440	0.68	ppb	# 86
31) Ethyl acetate	9.72	43	76325	0.70	ppb	98
32) Chloroform	9.82	83	104984	0.71	ppb	100
33) Tetrahydrofuran	10.35	42	21496	0.69	ppb	# 71
34) 1,2-dichloroethane	10.73	62	70040	0.73	ppb	94
36) 1,1,1-trichloroethane	11.05	97	114950	0.72	ppb	98
37) Cyclohexane	11.96	56	35814	0.77	ppb	88
38) Carbon tetrachloride	11.80	117	149024	0.76	ppb	99
39) Benzene	11.62	78	91372	0.74	ppb	94
40) Methyl methacrylate	13.11	41	32920	0.71	ppb	# 80
41) 1,4-dioxane	12.92	58	17273	0.76	ppb	# 17
42) 2,2,4-trimethylpentane	12.92	57	126105	0.73	ppb	87
43) Heptane	13.23	43	39307	0.77	ppb	84
44) Trichloroethene	12.88	130	55754	0.75	ppb	97
45) 1,2-dichloropropane	12.61	63	31317	0.76	ppb	99

(#)=qualifier out of range (m)=manual integration

AO022706.D A227\_1UG.M

Mon Mar 06 09:23:05 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO022706.D

Vial: 4

Acq On : 27 Feb 2017 2:01 pm

Operator: RJP

Sample : A1UG\_0.75

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 14:27:59 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.83	83	114702	0.75	ppb	99
47) cis-1,3-dichloropropene	13.88	75	54268	0.75	ppb	96
48) trans-1,3-dichloropropene	14.46	75	53258	0.74	ppb	94
49) 1,1,2-trichloroethane	14.66	97	51718	0.77	ppb	100
51) Toluene	14.98	92	64300	0.70	ppb	87
52) Methyl Isobutyl Ketone	13.92	43	55839	0.61	ppb	98
53) Dibromochloromethane	15.44	129	125722	0.74	ppb	94
54) Methyl Butyl Ketone	15.26	43	48541	0.54	ppb	99
55) 1,2-dibromoethane	15.72	107	80352	0.70	ppb	96
56) Tetrachloroethylene	16.23	164	57805	0.72	ppb	94
57) Chlorobenzene	16.95	112	108029	0.72	ppb	96
58) Ethylbenzene	17.35	91	146857	0.70	ppb	96
59) m&p-xylene	17.55	91	267284m	1.25	ppb	
60) Nonane	18.26	43	61387	0.69	ppb	97
61) Styrene	17.92	104	80315	0.70	ppb	87
62) Bromoform	17.63	173	113759	0.70	ppb	97
63) o-xylene	18.05	91	141768	0.70	ppb	93
64) Cumene	18.68	105	192362	0.70	ppb	97
66) 1,1,2,2-tetrachloroethane	18.04	83	111551	0.74	ppb	98
67) Propylbenzene	19.25	120	54005	0.69	ppb	# 1
68) 2-Chlorotoluene	19.22	126	45788	0.74	ppb	# 1
69) 4-ethyltoluene	19.41	105	178881	0.69	ppb	100
70) 1,3,5-trimethylbenzene	19.49	105	173200	0.69	ppb	99
71) 1,2,4-trimethylbenzene	19.97	105	138646	0.68	ppb	94
72) 1,3-dichlorobenzene	20.16	146	112138m	0.70	ppb	
73) benzyl chloride	20.14	91	95731	0.67	ppb	91
74) 1,4-dichlorobenzene	20.24	146	100495m	0.71	ppb	
75) 1,2,3-trimethylbenzene	20.48	105	143014	0.64	ppb	93
76) 1,2-dichlorobenzene	20.66	146	113886	0.72	ppb	97
77) 1,2,4-trichlorobenzene	22.83	180	37319	0.59	ppb	94
78) Naphthalene	22.98	128	98076	0.55	ppb	96
79) Hexachloro-1,3-butadiene	23.41	225	109840	0.65	ppb	99

[illegible]

Data File : C:\HPCHEM\1\DATA\AO022707.D

Vial: 5

Acq On : 27 Feb 2017 2:38 pm

Operator: RJP

Sample : A1UG\_0.50

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 15:23:11 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.66	128	32086	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.04	114	132674	1.00	ppb	0.00
50) Chlorobenzene-d5	16.91	117	112670	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	71762	0.95	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	95.00%

## Target Compounds

						Qvalue
2) Propylene	4.56	41	17331	0.56	ppb	79
3) Freon 12	4.64	85	155415	0.51	ppb	97
4) Chloromethane	4.80	50	29173m	0.52	ppb	
5) Freon 114	4.90	85	136873	0.50	ppb	96
6) Vinyl Chloride	5.04	62	36547	0.53	ppb	87
7) Butane	5.23	43	41383	0.52	ppb	91
8) 1,3-butadiene	5.17	39	27126m	0.51	ppb	
9) Bromomethane	5.46	94	49045	0.50	ppb	91
10) Chloroethane	5.62	64	17099	0.52	ppb	# 74
11) Ethanol	5.74	45	13536	0.61	ppb	# 75
12) Acrolein	6.10	56	14121	0.53	ppb	84
13) Vinyl Bromide	5.99	106	48399	0.49	ppb	91
14) Freon 11	6.42	101	188460	0.49	ppb	99
15) Acetone	6.24	58	14669	0.51	ppb	# 64
16) Pentane	6.79	42	16870	0.52	ppb	# 47
17) Isopropyl alcohol	6.48	45	53316	0.52	ppb	# 100
18) 1,1-dichloroethene	7.10	96	20467	0.48	ppb	98
19) Freon 113	7.50	101	56765	0.51	ppb	97
20) t-Butyl alcohol	7.14	59	40359	0.48	ppb	# 87
21) Methylene chloride	7.21	84	22237	0.53	ppb	89
22) Allyl chloride	7.34	41	17163	0.48	ppb	89
23) Carbon disulfide	7.56	76	65215	0.51	ppb	81
24) trans-1,2-dichloroethene	8.26	61	29663	0.52	ppb	95
25) methyl tert-butyl ether	8.56	73	50598	0.48	ppb	92
26) 1,1-dichloroethane	8.48	63	40382	0.51	ppb	94
27) Vinyl acetate	8.63	43	36091m	0.48	ppb	
28) Methyl Ethyl Ketone	8.94	72	8973	0.50	ppb	# 1
29) cis-1,2-dichloroethene	9.47	61	24130	0.46	ppb	89
30) Hexane	9.71	57	20572	0.45	ppb	88
31) Ethyl acetate	9.72	43	46227	0.45	ppb	98
32) Chloroform	9.81	83	67700	0.49	ppb	98
33) Tetrahydrofuran	10.34	42	13012	0.45	ppb	# 71
34) 1,2-dichloroethane	10.73	62	43067	0.48	ppb	93
36) 1,1,1-trichloroethane	11.05	97	76976	0.51	ppb	98
37) Cyclohexane	11.96	56	23337	0.53	ppb	90
38) Carbon tetrachloride	11.80	117	97284	0.52	ppb	99
39) Benzene	11.62	78	59123	0.50	ppb	95
40) Methyl methacrylate	13.10	41	19757	0.45	ppb	# 83
41) 1,4-dioxane	12.90	58	10792	0.50	ppb	# 22
42) 2,2,4-trimethylpentane	12.92	57	76483	0.47	ppb	87
43) Heptane	13.22	43	23168	0.47	ppb	88
44) Trichloroethene	12.88	130	33635	0.47	ppb	91
45) 1,2-dichloropropane	12.60	63	20090	0.51	ppb	97

(#)=qualifier out of range (m)=manual integration

AO022707.D A227\_1UG.M

Mon Mar 06 09:23:09 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO022707.D

Vial: 5

Acq On : 27 Feb 2017 2:38 pm

Operator: RJP

Sample : A1UG\_0.50

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 15:23:11 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.83	83	72852	0.50	ppb	96
47) cis-1,3-dichloropropene	13.87	75	33191	0.48	ppb	94
48) trans-1,3-dichloropropene	14.46	75	32559	0.47	ppb	94
49) 1,1,2-trichloroethane	14.66	97	31820	0.50	ppb	93
51) Toluene	14.98	92	40547	0.45	ppb	89
52) Methyl Isobutyl Ketone	13.91	43	32500	0.36	ppb	97
53) Dibromochloromethane	15.45	129	77782	0.46	ppb	92
54) Methyl Butyl Ketone	15.26	43	29867	0.34	ppb	98
55) 1,2-dibromoethane	15.71	107	50516	0.45	ppb	94
56) Tetrachloroethylene	16.23	164	37382	0.47	ppb	93
57) Chlorobenzene	16.95	112	67067	0.45	ppb	97
58) Ethylbenzene	17.35	91	90343	0.44	ppb	96
59) m&p-xylene	17.54	91	171931	0.82	ppb	98
60) Nonane	18.26	43	35556	0.40	ppb	98
61) Styrene	17.92	104	50031	0.44	ppb	87
62) Bromoform	17.63	173	71380	0.45	ppb	98
63) o-xylene	18.05	91	81347	0.41	ppb	96
64) Cumene	18.68	105	115251	0.43	ppb	98
66) 1,1,2,2-tetrachloroethane	18.04	83	68065	0.46	ppb	94
67) Propylbenzene	19.25	120	31968	0.42	ppb #	1
68) 2-Chlorotoluene	19.22	126	26957	0.44	ppb #	1
69) 4-ethyltoluene	19.41	105	109155	0.43	ppb	100
70) 1,3,5-trimethylbenzene	19.49	105	99045	0.40	ppb	94
71) 1,2,4-trimethylbenzene	19.97	105	79227	0.39	ppb	88
72) 1,3-dichlorobenzene	20.16	146	67990	0.43	ppb	98
73) benzyl chloride	20.13	91	58313	0.42	ppb	88
74) 1,4-dichlorobenzene	20.24	146	59437	0.43	ppb	95
75) 1,2,3-trimethylbenzene	20.49	105	85990	0.39	ppb	91
76) 1,2-dichlorobenzene	20.66	146	70376	0.45	ppb	95
77) 1,2,4-trichlorobenzene	22.83	180	22620	0.36	ppb	92
78) Naphthalene	22.97	128	58170	0.33	ppb	95
79) Hexachloro-1,3-butadiene	23.41	225	70735	0.42	ppb	99

-----  
 (#) = qualifier out of range (m) = manual integration (+) = signals summed  
 AO022707.D A227\_1UG.M Mon Mar 06 09:23:09 2017 MSD1





Data File : C:\HPCHEM\1\DATA\AO022708.D

Vial: 6

Acq On : 27 Feb 2017 3:18 pm

Operator: RJP

Sample : A1UG\_1.0

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 19:08:49 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	9.66	128	33193	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.05	114	135563	1.00	ppb	0.01
50) Chlorobenzene-d5	16.91	117	112581	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	74287	0.98	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	98.00%

## Target Compounds

						Qvalue
2) Propylene	4.57	41	32616	1.01	ppb	82
3) Freon 12	4.63	85	314355	0.99	ppb	97
4) Chloromethane	4.80	50	55242m	0.95	ppb	
5) Freon 114	4.91	85	266565	0.94	ppb	98
6) Vinyl Chloride	5.03	62	71834	1.00	ppb	88
7) Butane	5.23	43	77054	0.93	ppb	87
8) 1,3-butadiene	5.17	39	47992m	0.87	ppb	
9) Bromomethane	5.45	94	95659	0.94	ppb	92
10) Chloroethane	5.63	64	33347	0.99	ppb	93
11) Ethanol	5.73	45	24047	1.05	ppb	87
12) Acrolein	6.09	56	24461	0.90	ppb	97
13) Vinyl Bromide	5.99	106	96108	0.95	ppb	95
14) Freon 11	6.42	101	371968	0.93	ppb	98
15) Acetone	6.24	58	29555	0.99	ppb	# 67
16) Pentane	6.79	42	31576	0.93	ppb	# 39
17) Isopropyl alcohol	6.47	45	100359	0.95	ppb	# 100
18) 1,1-dichloroethene	7.10	96	40530	0.93	ppb	98
19) Freon 113	7.50	101	116271	1.00	ppb	97
20) t-Butyl alcohol	7.12	59	78945	0.91	ppb	# 88
21) Methylene chloride	7.21	84	42167	0.96	ppb	95
22) Allyl chloride	7.35	41	34889	0.94	ppb	94
23) Carbon disulfide	7.57	76	130323	0.98	ppb	81
24) trans-1,2-dichloroethene	8.26	61	56081	0.95	ppb	92
25) methyl tert-butyl ether	8.55	73	103821	0.95	ppb	90
26) 1,1-dichloroethane	8.48	63	79911	0.97	ppb	98
27) Vinyl acetate	8.63	43	77184m	0.99	ppb	
28) Methyl Ethyl Ketone	8.94	72	18231	0.97	ppb	# 1
29) cis-1,2-dichloroethene	9.46	61	52575	0.98	ppb	93
30) Hexane	9.71	57	44374	0.94	ppb	87
31) Ethyl acetate	9.72	43	100553	0.96	ppb	98
32) Chloroform	9.81	83	133894	0.94	ppb	97
33) Tetrahydrofuran	10.34	42	27337	0.92	ppb	# 65
34) 1,2-dichloroethane	10.73	62	87348	0.95	ppb	95
36) 1,1,1-trichloroethane	11.05	97	152512	0.98	ppb	97
37) Cyclohexane	11.96	56	44782	0.99	ppb	88
38) Carbon tetrachloride	11.80	117	192671	1.00	ppb	99
39) Benzene	11.62	78	119767	0.99	ppb	95
40) Methyl methacrylate	13.10	41	44765	0.99	ppb	# 83
41) 1,4-dioxane	12.91	58	23899	1.08	ppb	# 34
42) 2,2,4-trimethylpentane	12.92	57	169708	1.01	ppb	89
43) Heptane	13.23	43	50744	1.02	ppb	84
44) Trichloroethene	12.88	130	72013	1.00	ppb	96
45) 1,2-dichloropropane	12.60	63	42501	1.06	ppb	99

(#)= qualifier out of range (m)= manual integration

AO022708.D A227\_1UG.M

Mon Mar 06 09:23:13 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO022708.D

Vial: 6

Acq On : 27 Feb 2017 3:18 pm

Operator: RJP

Sample : A1UG\_1.0

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 19:08:49 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

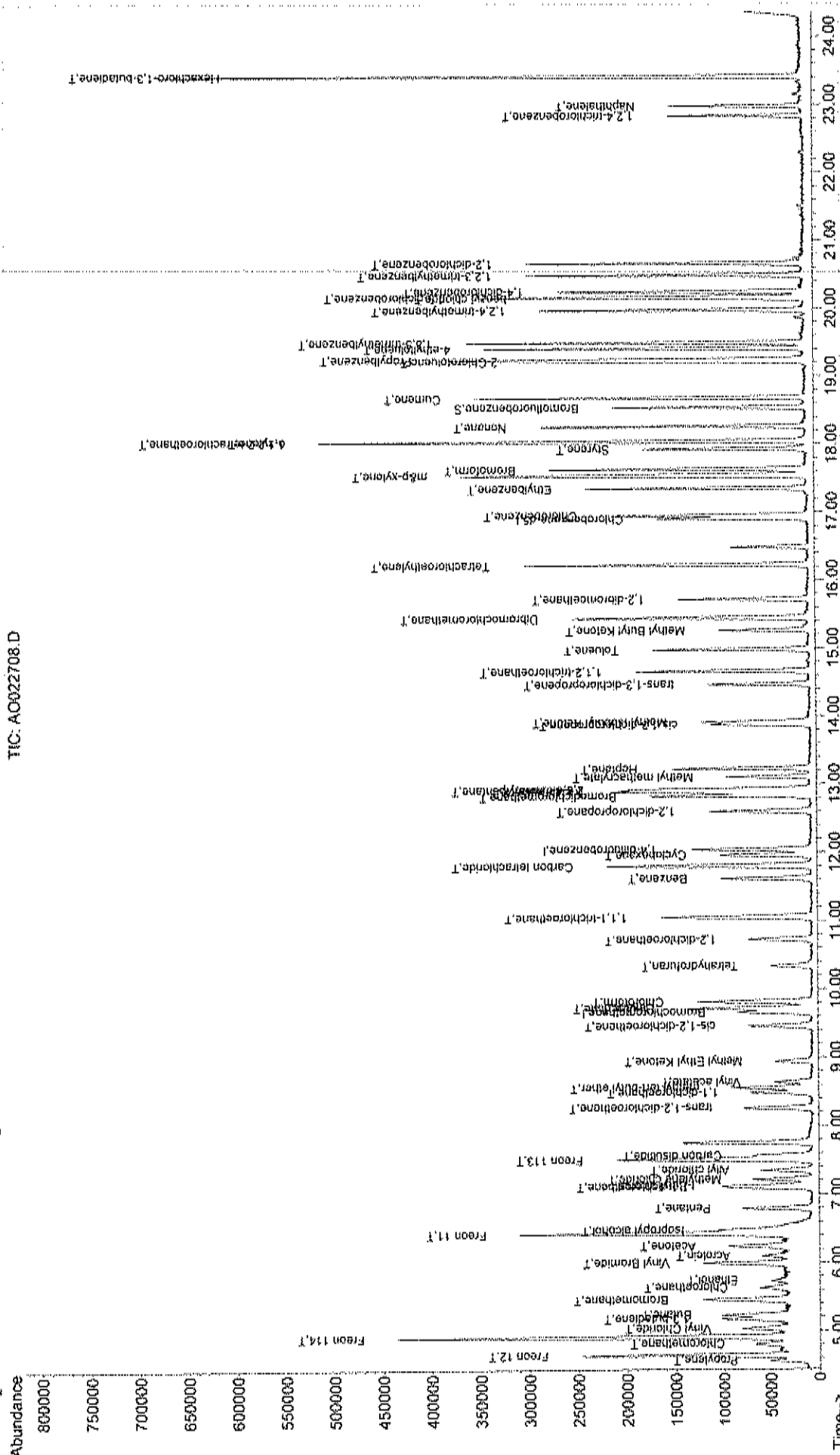
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.82	83	145355	0.98	ppb	97
47) cis-1,3-dichloropropene	13.87	75	72816	1.03	ppb	97
48) trans-1,3-dichloropropene	14.47	75	71040	1.01	ppb	98
49) 1,1,2-trichloroethane	14.67	97	67705	1.04	ppb	95
51) Toluene	14.97	92	86837	0.96	ppb	88
52) Methyl Isobutyl Ketone	13.92	43	74181	0.82	ppb	97
53) Dibromochloromethane	15.45	129	162151	0.97	ppb	92
54) Methyl Butyl Ketone	15.26	43	65439	0.74	ppb	99
55) 1,2-dibromoethane	15.72	107	107224	0.95	ppb	97
56) Tetrachloroethylene	16.22	164	78753	0.99	ppb	100
57) Chlorobenzene	16.95	112	140548	0.95	ppb	98
58) Ethylbenzene	17.35	91	199134	0.96	ppb	98
59) m&p-xylene	17.54	91	399403	1.91	ppb	100
60) Nonane	18.26	43	84171	0.96	ppb	98
61) Styrene	17.92	104	109876	0.97	ppb	87
62) Bromoform	17.63	173	153180	0.96	ppb	98
63) o-xylene	18.04	91	195397	0.98	ppb	94
64) Cumene	18.68	105	257295	0.96	ppb	99
66) 1,1,2,2-tetrachloroethane	18.03	83	142956	0.96	ppb	96
67) Propylbenzene	19.24	120	73966	0.97	ppb	# 1
68) 2-Chlorotoluene	19.21	126	59127	0.97	ppb	# 1
69) 4-ethyltoluene	19.41	105	234955	0.93	ppb	94
70) 1,3,5-trimethylbenzene	19.49	105	232316	0.94	ppb	98
71) 1,2,4-trimethylbenzene	19.98	105	186326	0.93	ppb	91
72) 1,3-dichlorobenzene	20.16	146	148198	0.94	ppb	# 35
73) benzyl chloride	20.14	91	129063	0.92	ppb	88
74) 1,4-dichlorobenzene	20.24	146	132228	0.95	ppb	# 25
75) 1,2,3-trimethylbenzene	20.49	105	198131	0.91	ppb	93
76) 1,2-dichlorobenzene	20.66	146	145697	0.94	ppb	96
77) 1,2,4-trichlorobenzene	22.83	180	51354	0.83	ppb	96
78) Naphthalene	22.97	128	128684	0.74	ppb	94
79) Hexachloro-1,3-butadiene	23.41	225	142832	0.86	ppb	98

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 (#) = qualifier out of range (m) = manual integration (+) = signals summed  
 AO022708.D A227\_1UG.M Mon Mar 06 09:23:13 2017 MSD1

Data File : C:\HPCHEM\1\DATA\AO022708.D  
Acq On : 27 Feb 2017 3:18 pm  
Sample : A1UG\_1.0  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Feb 27 19:11 2017  
Quant Results File: A227\_1UG.RBS

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Feb 27 19:25:53 2017  
Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

TIC: AO022708.D



Data File : C:\HPCHEM\1\DATA\AO022709.D

Vial: 7

Acq On : 27 Feb 2017 3:55 pm

Operator: RJP

Sample : A1UG\_0,30

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 19:09:15 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	9.66	128	31550	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.04	114	124545	1.00	ppb	0.00
50) Chlorobenzene-d5	16.91	117	107408	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	65066	0.90	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	90.00%

## Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Propylene	4.56	41	10462	0.34	ppb	80
3) Freon 12	4.64	85	89651	0.30	ppb	99
4) Chloromethane	4.80	50	16559	0.30	ppb	68
5) Freon 114	4.90	85	74954	0.28	ppb	97
6) Vinyl Chloride	5.03	62	20335	0.30	ppb	96
7) Butane	5.23	43	24032	0.31	ppb	94
8) 1,3-butadiene	5.18	39	15070m	0.29	ppb	
9) Bromomethane	5.45	94	26711	0.27	ppb	94
10) Chloroethane	5.63	64	10587m	0.33	ppb	
11) Ethanol	5.73	45	8502	0.39	ppb	# 77
12) Acrolein	6.11	56	7749	0.30	ppb	89
13) Vinyl Bromide	5.99	106	27527	0.28	ppb	96
14) Freon 11	6.42	101	107968	0.29	ppb	99
15) Acetone	6.25	58	8644	0.30	ppb	# 52
16) Pentane	6.79	42	9719	0.30	ppb	# 45
17) Isopropyl alcohol	6.49	45	37741	0.38	ppb	# 100
18) 1,1-dichloroethene	7.10	96	12295	0.30	ppb	95
19) Freon 113	7.51	101	33311	0.30	ppb	98
20) t-Butyl alcohol	7.14	59	26163	0.32	ppb	# 87
21) Methylene chloride	7.21	84	13237m	0.32	ppb	
22) Allyl chloride	7.33	41	9501	0.27	ppb	86
23) Carbon disulfide	7.57	76	38692	0.31	ppb	79
24) trans-1,2-dichloroethene	8.25	61	16783	0.30	ppb	92
25) methyl tert-butyl ether	8.56	73	28349	0.27	ppb	94
26) 1,1-dichloroethane	8.48	63	22556	0.29	ppb	92
27) Vinyl acetate	8.65	43	18505m	0.25	ppb	
28) Methyl Ethyl Ketone	8.94	72	5537	0.31	ppb	# 1
29) cis-1,2-dichloroethene	9.46	61	15110	0.30	ppb	96
30) Hexane	9.72	57	10444	0.23	ppb	# 85
31) Ethyl acetate	9.72	43	25081	0.25	ppb	98
32) Chloroform	9.81	83	37202	0.27	ppb	95
33) Tetrahydrofuran	10.34	42	7963	0.28	ppb	# 61
34) 1,2-dichloroethane	10.73	62	24501	0.28	ppb	88
36) 1,1,1-trichloroethane	11.05	97	43540	0.31	ppb	98
37) Cyclohexane	11.95	56	13093	0.31	ppb	89
38) Carbon tetrachloride	11.80	117	54833	0.31	ppb	99
39) Benzene	11.62	78	32717	0.30	ppb	95
40) Methyl methacrylate	13.11	41	10561	0.25	ppb	# 75
41) 1,4-dioxane	12.92	58	6900	0.34	ppb	# 23
42) 2,2,4-trimethylpentane	12.92	57	39374	0.26	ppb	78
43) Heptane	13.23	43	11286	0.25	ppb	# 76
44) Trichloroethene	12.89	130	19504	0.29	ppb	94
45) 1,2-dichloropropane	12.60	63	11186	0.30	ppb	94

(#)=qualifier out of range (m)=manual integration

AO022709.D A227\_1UG.M

Mon Mar 06 09:23:17 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO022709.D

Vial: 7

Acq On : 27 Feb 2017 3:55 pm

Operator: RJP

Sample : A1UG\_0,30

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 19:09:15 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

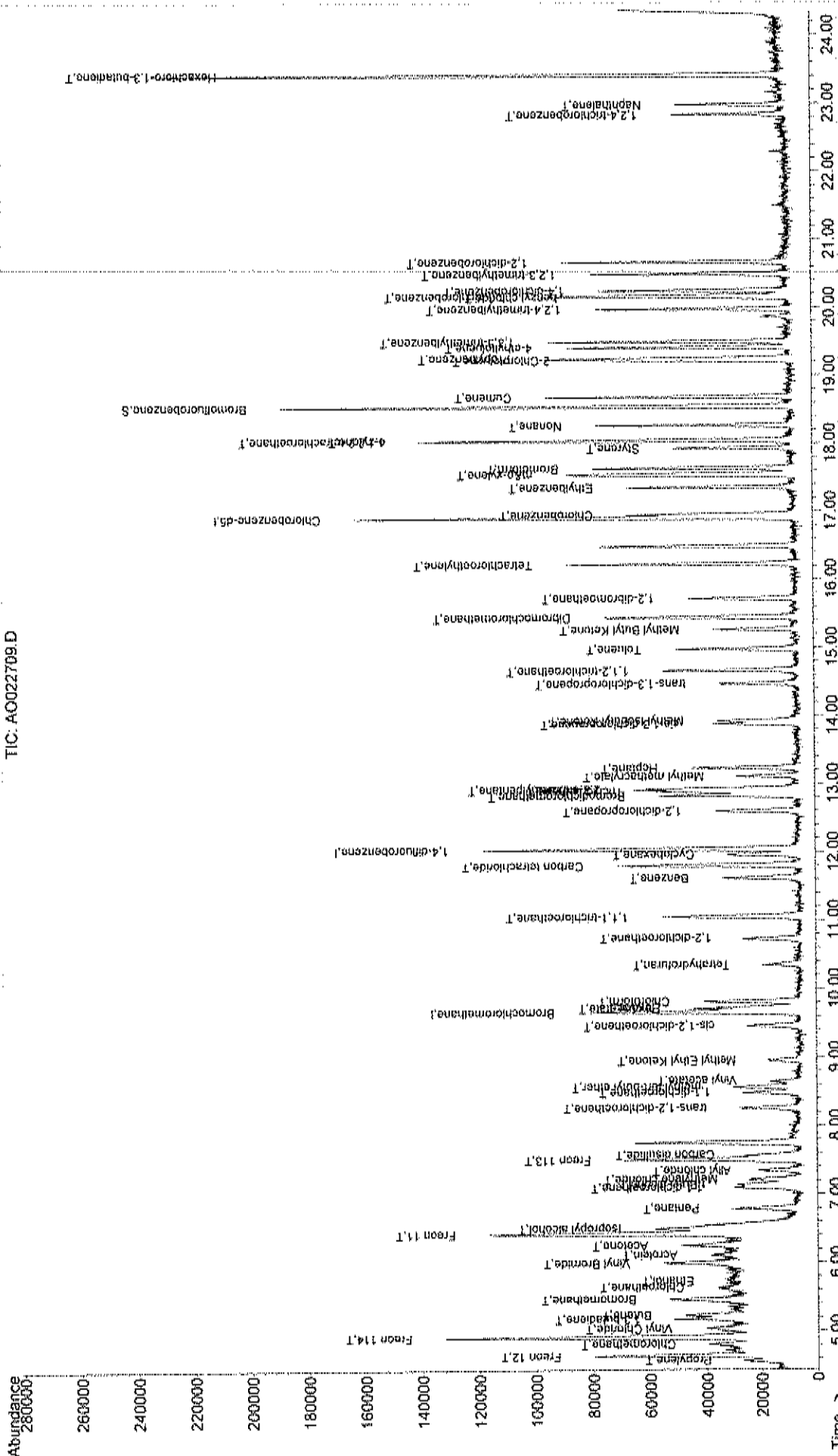
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.83	83	41743	0.31	ppb	93
47) cis-1,3-dichloropropene	13.88	75	19493	0.30	ppb	93
48) trans-1,3-dichloropropene	14.46	75	17923	0.28	ppb	91
49) 1,1,2-trichloroethane	14.66	97	19136	0.32	ppb	99
51) Toluene	14.98	92	22427	0.26	ppb	87
52) Methyl Isobutyl Ketone	13.92	43	20602	0.24	ppb	96
53) Dibromochloromethane	15.44	129	45949	0.29	ppb	94
54) Methyl Butyl Ketone	15.26	43	19265	0.23	ppb	98
55) 1,2-dibromoethane	15.71	107	29196	0.27	ppb	99
56) Tetrachloroethylene	16.23	164	20923	0.28	ppb	91
57) Chlorobenzene	16.95	112	37466	0.27	ppb	93
58) Ethylbenzene	17.35	91	49406	0.25	ppb	99
59) m&p-xylene	17.54	91	92629m	0.46	ppb	
60) Nonane	18.26	43	16910	0.20	ppb	89
61) Styrene	17.92	104	26733	0.25	ppb	84
62) Bromoform	17.63	173	40478	0.27	ppb	98
63) o-xylene	18.04	91	43261	0.23	ppb	96
64) Cumene	18.68	105	64995	0.25	ppb	98
66) 1,1,2,2-tetrachloroethane	18.04	83	39374	0.28	ppb	95
67) Propylbenzene	19.25	120	17405	0.24	ppb	# 1
68) 2-Chlorotoluene	19.22	126	15954	0.27	ppb	# 1
69) 4-ethyltoluene	19.41	105	57355	0.24	ppb	99
70) 1,3,5-trimethylbenzene	19.49	105	53840	0.23	ppb	93
71) 1,2,4-trimethylbenzene	19.97	105	46850	0.24	ppb	88
72) 1,3-dichlorobenzene	20.16	146	41823	0.28	ppb	# 60
73) benzyl chloride	20.14	91	33683	0.25	ppb	92
74) 1,4-dichlorobenzene	20.25	146	35292m	0.27	ppb	
75) 1,2,3-trimethylbenzene	20.49	105	48489	0.23	ppb	90
76) 1,2-dichlorobenzene	20.66	146	40713	0.27	ppb	100
77) 1,2,4-trichlorobenzene	22.83	180	13945	0.24	ppb	87
78) Naphthalene	22.97	128	39073	0.23	ppb	97
79) Hexachloro-1,3-butadiene	23.41	225	46959	0.30	ppb	99

-----  
 (#) = qualifier out of range (m) = manual integration (+) = signals summed  
 AO022709.D A227\_1UG.M Mon Mar 06 09:23:17 2017 MSD1

Data File : C:\HPCHEM\1\DATA\AO022709.D  
Acq On : 27 Feb 2017 3:55 pm  
Sample : A1UG\_0,30  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Feb 27 19:13 2017  
Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Feb 27 19:25:53 2017  
Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

TIC: AO022709.D



Data File : C:\HPCHEM\1\DATA\AO022710.D

Vial: 8

Acq On : 27 Feb 2017 4:31 pm

Operator: RJP

Sample : A1UG\_0.15

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 19:09:36 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	9.66	128	29609	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.04	114	114280	1.00	ppb	0.00
50) Chlorobenzene-d5	16.91	117	101915	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	60579	0.88	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	88.00%

## Target Compounds

						Qvalue
2) Propylene	4.56	41	4783	0.17	ppb	89
3) Freon 12	4.63	85	43314	0.15	ppb	99
4) Chloromethane	4.80	50	7779m	0.15	ppb	
5) Freon 114	4.90	85	37829	0.15	ppb	96
6) Vinyl Chloride	5.03	62	10571	0.17	ppb	84
7) Butane	5.22	43	12241	0.17	ppb	92
8) 1,3-butadiene	5.17	39	7233m	0.15	ppb	
9) Bromomethane	5.45	94	13921	0.15	ppb	92
10) Chloroethane	5.62	64	4796	0.16	ppb	# 68
11) Ethanol	5.74	45	3726m	0.18	ppb	
12) Acrolein	6.10	56	3425	0.14	ppb	79
13) Vinyl Bromide	5.99	106	13042	0.14	ppb	89
14) Freon 11	6.42	101	52808	0.15	ppb	99
15) Acetone	6.25	58	4275	0.16	ppb	# 63
16) Pentane	6.78	42	4819	0.16	ppb	# 48
17) Isopropyl alcohol	6.49	45	17359	0.19	ppb	# 100
18) 1,1-dichloroethene	7.10	96	5776	0.15	ppb	92
19) Freon 113	7.50	101	16587	0.16	ppb	95
20) t-Butyl alcohol	7.15	59	10937	0.14	ppb	# 88
21) Methylene chloride	7.22	84	6851	0.18	ppb	93
22) Allyl chloride	7.34	41	5007m	0.15	ppb	
23) Carbon disulfide	7.56	76	19583	0.17	ppb	87
24) trans-1,2-dichloroethene	8.25	61	8641	0.16	ppb	90
25) methyl tert-butyl ether	8.57	73	13267	0.14	ppb	94
26) 1,1-dichloroethane	8.49	63	9732	0.13	ppb	88
27) Vinyl acetate	8.63	43	11754	0.17	ppb	83
28) Methyl Ethyl Ketone	8.93	72	2318	0.14	ppb	# 80
29) cis-1,2-dichloroethene	9.46	61	6901	0.14	ppb	86
30) Hexane	9.72	57	4955	0.12	ppb	# 83
31) Ethyl acetate	9.72	43	11816	0.13	ppb	91
32) Chloroform	9.81	83	18312	0.14	ppb	93
33) Tetrahydrofuran	10.35	42	3640m	0.14	ppb	
34) 1,2-dichloroethane	10.72	62	12370	0.15	ppb	93
36) 1,1,1-trichloroethane	11.05	97	21536	0.16	ppb	96
37) Cyclohexane	11.97	56	5394	0.14	ppb	# 76
38) Carbon tetrachloride	11.81	117	27365	0.17	ppb	98
39) Benzene	11.62	78	16374	0.16	ppb	94
40) Methyl methacrylate	13.10	41	4234	0.11	ppb	# 64
41) 1,4-dioxane	12.91	58	2799	0.15	ppb	# 14
42) 2,2,4-trimethylpentane	12.93	57	17819	0.13	ppb	84
43) Heptane	13.24	43	5482	0.13	ppb	# 73
44) Trichloroethene	12.88	130	9089	0.15	ppb	86
45) 1,2-dichloropropane	12.60	63	5722	0.17	ppb	95

(# ) = qualifier out of range (m) = manual integration

AO022710.D A227\_1UG.M

Mon Mar 06 09:23:21 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO022710.D

Vial: 8

Acq On : 27 Feb 2017 4:31 pm

Operator: RJP

Sample : A1UG\_0.15

Inst : MSD #1

Misc : A227\_1UG

Multiplier: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 19:09:36 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.83	83	20004	0.16	ppb	98
47) cis-1,3-dichloropropene	13.87	75	8700	0.15	ppb	90
48) trans-1,3-dichloropropene	14.46	75	8707	0.15	ppb	98
49) 1,1,2-trichloroethane	14.67	97	8712	0.16	ppb	99
51) Toluene	14.98	92	11171	0.14	ppb	94
53) Dibromochloromethane	15.44	129	22261	0.15	ppb	100
55) 1,2-dibromoethane	15.72	107	13395	0.13	ppb	94
56) Tetrachloroethylene	16.23	164	11207	0.16	ppb	95
57) Chlorobenzene	16.95	112	18365	0.14	ppb	96
58) Ethylbenzene	17.35	91	24082	0.13	ppb	95
59) m&p-xylene	17.55	91	41618	0.22	ppb	98
60) Nonane	18.26	43	8797m	0.11	ppb	
61) Styrene	17.93	104	13421	0.13	ppb	81
62) Bromoform	17.63	173	18122	0.13	ppb	94
63) o-xylene	18.05	91	19911	0.11	ppb	99
64) Cumene	18.68	105	29600	0.12	ppb	95
66) 1,1,2,2-tetrachloroethane	18.04	83	18805	0.14	ppb	94
67) Propylbenzene	19.25	120	8095	0.12	ppb	# 1
68) 2-Chlorotoluene	19.22	126	7562	0.14	ppb	# 1
69) 4-ethyltoluene	19.41	105	25058	0.11	ppb	98
70) 1,3,5-trimethylbenzene	19.49	105	22630	0.10	ppb	87
71) 1,2,4-trimethylbenzene	19.97	105	22333	0.12	ppb	81
72) 1,3-dichlorobenzene	20.17	146	17500	0.12	ppb	99
73) benzyl chloride	20.14	91	15570	0.12	ppb	87
74) 1,4-dichlorobenzene	20.24	146	14264	0.11	ppb	97
75) 1,2,3-trimethylbenzene	20.48	105	22396	0.11	ppb	# 83
76) 1,2-dichlorobenzene	20.66	146	19520	0.14	ppb	96
77) 1,2,4-trichlorobenzene	22.83	180	7102	0.13	ppb	98
78) Naphthalene	22.97	128	17274	0.11	ppb	91
79) Hexachloro-1,3-butadiene	23.41	225	25872	0.17	ppb	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed  
 AO022710.D A227\_1UG.M Mon Mar 06 09:23:21 2017 MSD1

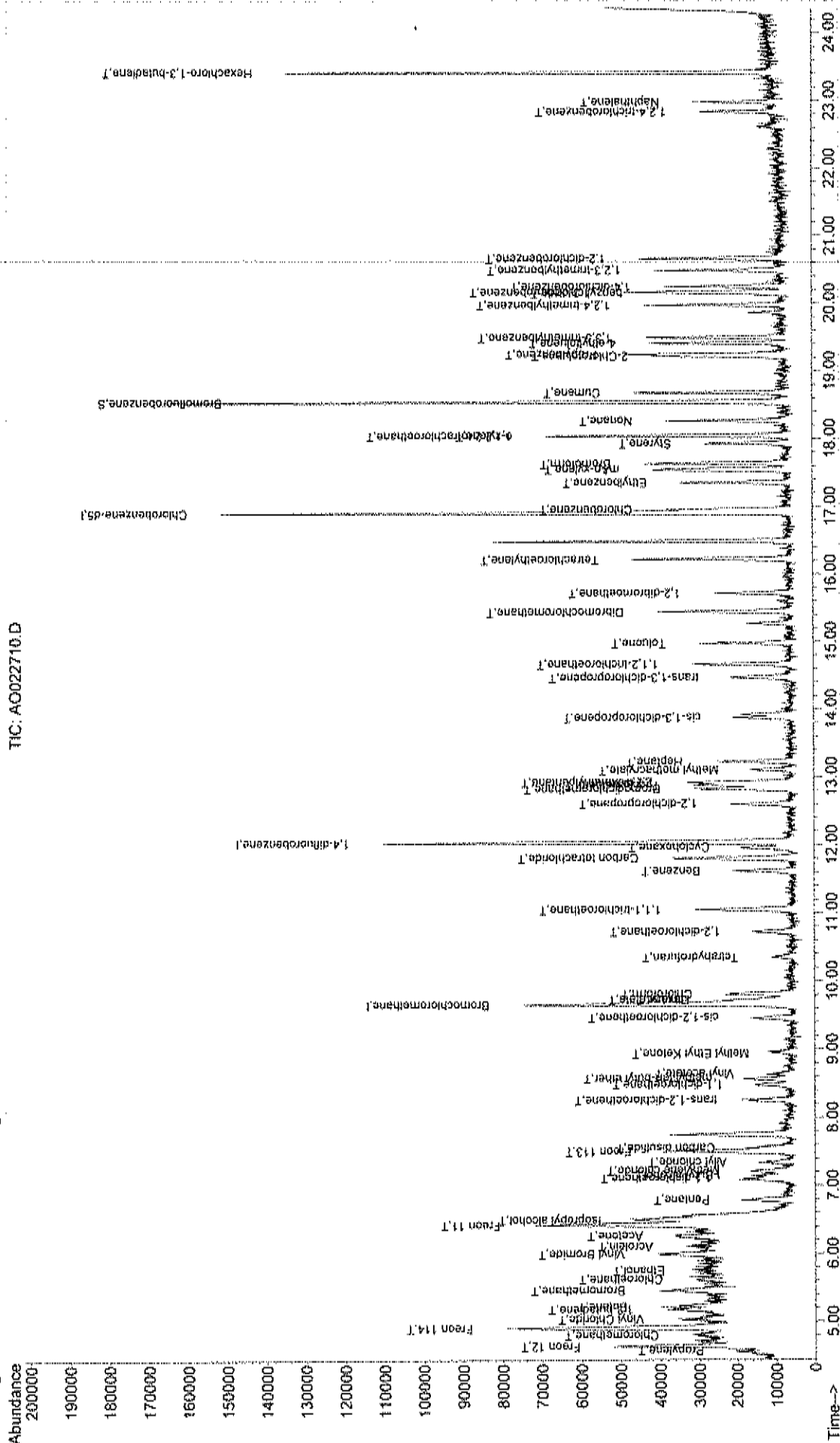


Data File : C:\HPCHEM\1\DATA\A0022710.D  
Acq On : 27 Feb 2017 4:31 pm  
Sample : A1UG 0.15  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Feb 27 19:15 2017

Vial: 8  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00  
Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Feb 27 19:25:53 2017  
Response via : Continuing Cal File: C:\HPCHEM\1\DATA\A0022702.D

THC: A0022710.D



Data File : C:\HPCHEM\1\DATA\AO022711.D

Vial: 9

Acq On : 27 Feb 2017 5:08 pm

Operator: RJP

Sample : A1UG\_0.10

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 19:10:00 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	9.66	128	29938	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.04	114	109904	1.00	ppb	0.00
50) Chlorobenzene-d5	16.91	117	98325	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	54341	0.82	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	82.00%

## Target Compounds

					Qvalue
6) Vinyl Chloride	5.04	62	7526	0.12	ppb 84
38) Carbon tetrachloride	11.80	117	17794	0.11	ppb 97
44) Trichloroethene	12.89	130	6065	0.10	ppb 87

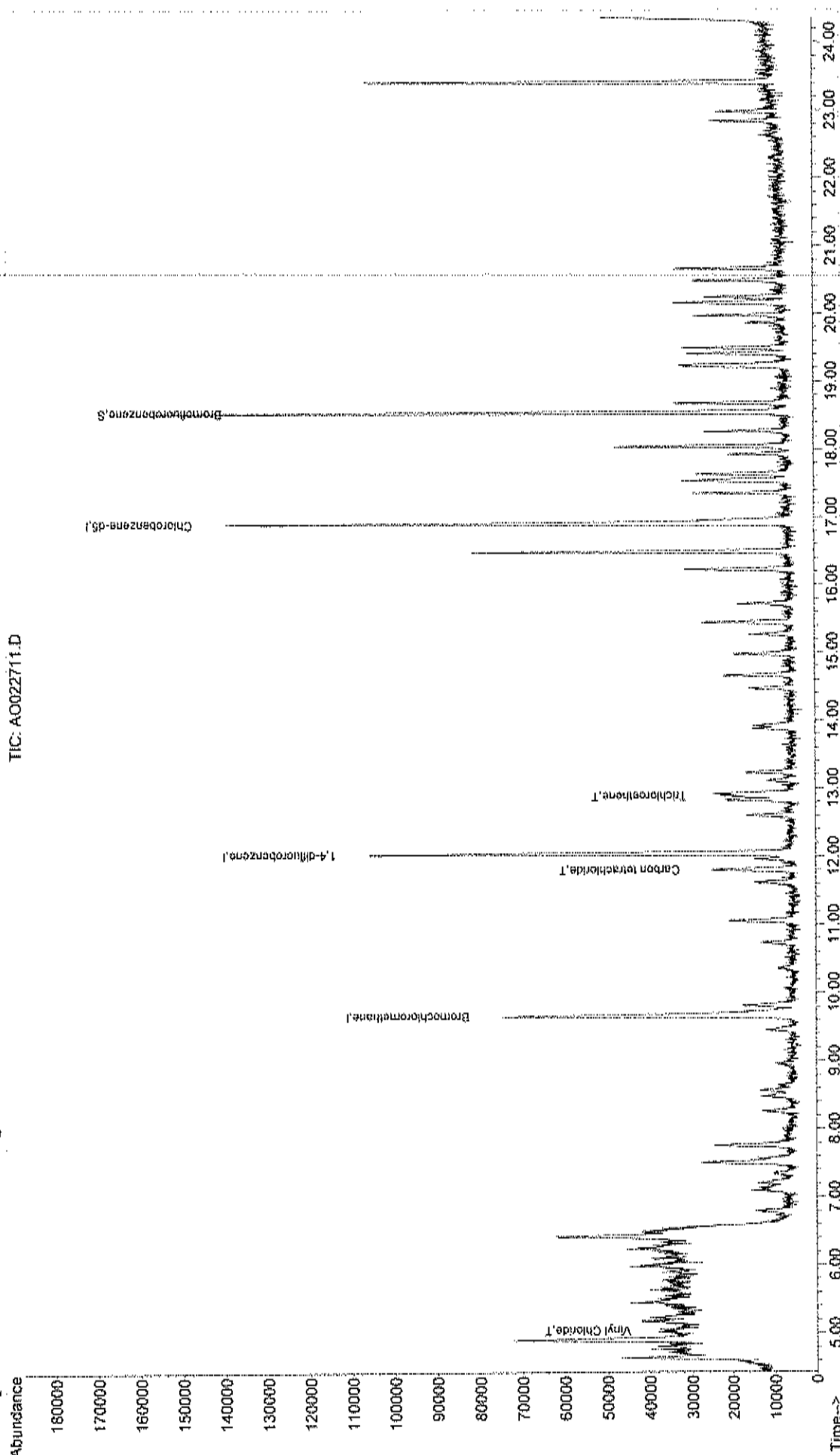
Quantitation Report (X4) AO022711.D

Data File : C:\HPCHEM\1\DATA\AO022711.D  
Acq On : 27 Feb 2017 5:08 pm  
Sample : A1UG\_0.10  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Feb 27 19:16 2017 Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Feb 27 19:25:53 2017  
Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

Vial: 9  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

TIC: AO022711.D



Data File : C:\HPCHEM\1\DATA\AO022712.D

Vial: 10

Acq On : 27 Feb 2017 5:45 pm

Operator: RJP

Sample : A1UG\_0.04

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 27 19:10:27 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Feb 27 13:34:10 2017

Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.66	128	29129	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.05	114	103388	1.00	ppb	0.01
50) Chlorobenzene-d5	16.91	117	93432	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	51957	0.83	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	83.00%

## Target Compounds

6) Vinyl Chloride	5.03	62	3893m	0.06	ppb	Qvalue
38) Carbon tetrachloride	11.80	117	11296m	0.08	ppb	
44) Trichloroethene	12.88	130	3725m	0.07	ppb	

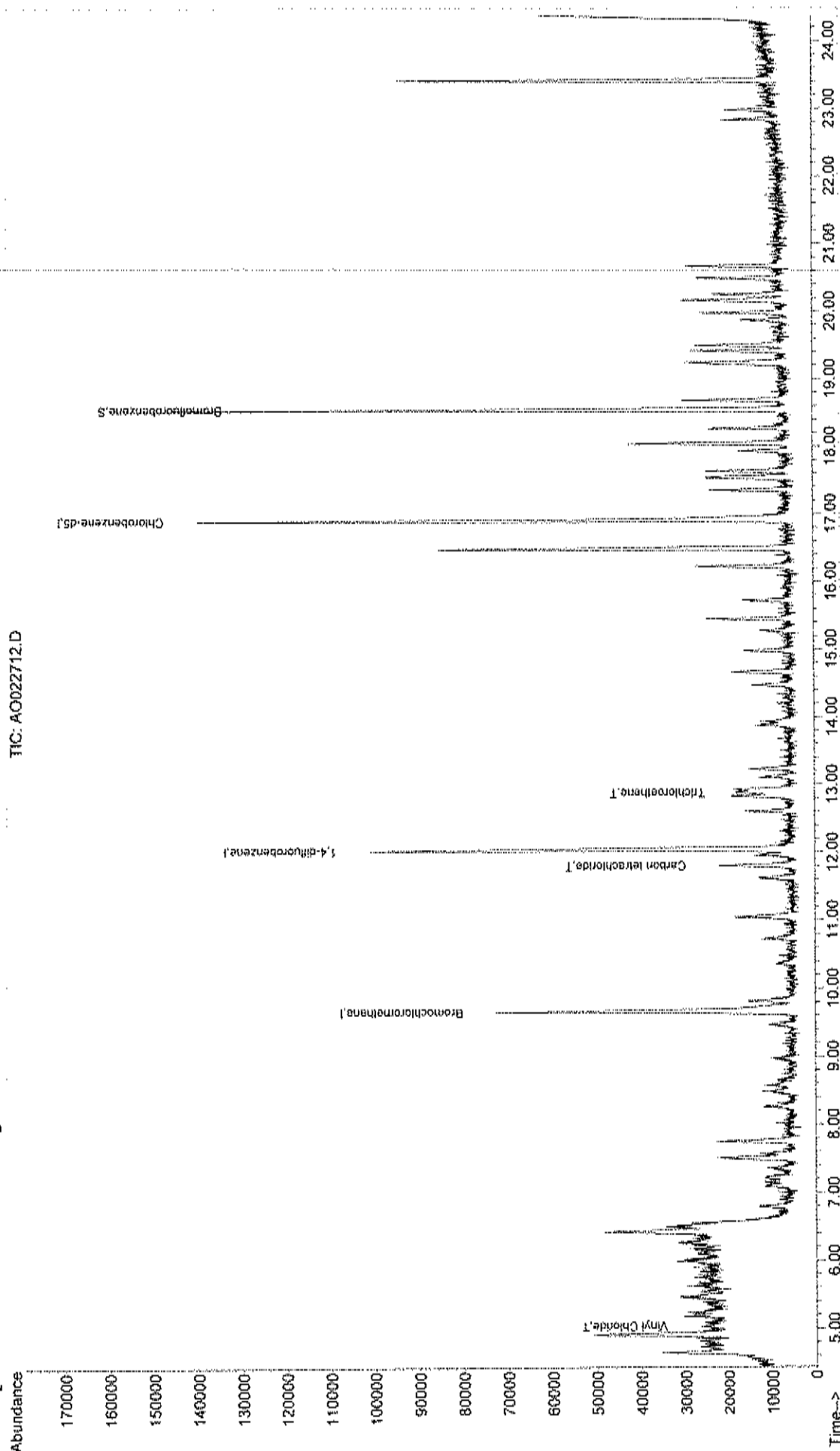
Data File : C:\HPCHEM\1\DATA\AO022712.D  
Acq On : 27 Feb 2017 5:45 pm  
Sample : A1UG 0.04  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Feb 27 19:25 2017

Vial: 10  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Feb 27 19:25:53 2017  
Response via : Continuing Cal File: C:\HPCHEM\1\DATA\AO022702.D

THC: AO022712.D



**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**CALIBRATION VERIFICATION**

## Evaluate Continuing Calibration Report

Data File : C:\HPCHEM\1\DATA\AO030803.D

Vial: 3

Acq On : 8 Mar 2017 10:33 am

Operator: RJP

Sample : A1UG\_1.0

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 27 11:00:28 2017

Response via : Multiple Level Calibration

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

Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane	1.000	1.000	0.0	155#	0.00
2 T	Propylene	1.042	0.908	12.9	144	0.00
3 T	Freon 12	9.266	10.061	-8.6	165#	0.00
4 T	Chloromethane	1.659	1.378	16.9	129	0.00
5 T	Freon 114	7.921	7.623	3.8	148	0.00
6 T	Vinyl Chloride	2.301	1.895	17.6	136	0.00
7 T	Butane	2.394	2.055	14.2	138	0.00
8 T	1,3-butadiene	1.535	1.410	8.1	152#	0.00
9 T	Bromomethane	2.831	2.442	13.7	132	0.00
10 T	Chloroethane	1.014	0.804	20.7	124	0.00
11 T	Ethanol	0.748	0.556	25.7	119	0.00
12 T	Acrolein	0.765	0.583	23.8	123	0.00
13 T	Vinyl Bromide	2.844	2.497	12.2	134	0.00
14 T	Freon 11	11.092	11.621	-4.8	161#	0.00
15 T	Acetone	0.882	0.734	16.8	128	0.00
16 T	Pentane	1.010	0.835	17.3	137	0.00
17 T	Isopropyl alcohol	3.221	2.645	17.9	136	0.00
18 T	1,1-dichloroethene	1.282	1.389	-8.3	177#	0.00
19 T	Freon 113	3.515	3.854	-9.6	171#	0.00
20 t	t-Butyl alcohol	2.527	2.816	-11.4	184#	0.00
21 T	Methylene chloride	1.335	1.268	5.0	155#	0.00
22 T	Allyl chloride	1.090	1.109	-1.7	164#	0.00
23 T	Carbon disulfide	3.995	3.725	6.8	148	0.00
24 T	trans-1,2-dichloroethene	1.805	1.947	-7.9	179#	0.00
25 T	methyl tert-butyl ether	3.166	3.784	-19.5	188#	0.00
26 T	1,1-dichloroethane	2.386	2.480	-3.9	160#	0.00
27 T	Vinyl acetate	2.361	3.637	-54.0#	243#	0.00
28 T	Methyl Ethyl Ketone	0.565	0.600	-6.2	170#	0.00
29 T	cis-1,2-dichloroethene	1.586	1.793	-13.1	176#	0.00
30 T	Hexane	1.306	1.346	-3.1	156#	0.00
31 T	Ethyl acetate	2.971	2.955	0.5	152#	0.00
32 T	Chloroform	4.078	4.790	-17.5	185#	0.00
33 T	Tetrahydrofuran	0.865	0.815	5.8	154#	0.00
34 T	1,2-dichloroethane	2.716	3.358	-23.6	198#	0.00
35 I	1,4-difluorobenzene	1.000	1.000	0.0	172#	0.00
36 T	1,1,1-trichloroethane	1.135	1.381	-21.7	211#	0.00
37 T	Cyclohexane	0.341	0.312	8.5	162#	0.00
38 T	Carbon tetrachloride	1.580	1.763	-11.6	213#	0.00
39 T	Benzene	0.891	0.884	0.8	172#	0.00
40 T	Methyl methacrylate	0.312	0.340	-9.0	177#	0.00
41 T	1,4-dioxane	0.172	0.137	20.3	134	0.00
42 T	2,2,4-trimethylpentane	1.185	1.052	11.2	144	0.00
43 T	Heptane	0.360	0.322	10.6	148	0.00
44 T	Trichloroethene	0.568	0.593	-4.4	192#	0.00
45 T	1,2-dichloropropane	0.306	0.250	18.3	137	0.00
46 T	Bromodichloromethane	1.088	1.203	-10.6	193#	0.00
47 T	cis-1,3-dichloropropene	0.524	0.528	-0.8	169#	0.00
48 T	trans-1,3-dichloropropene	0.513	0.594	-15.8	195#	0.00
49 T	1,1,2-trichloroethane	0.488	0.457	6.4	157#	0.00

-----  
(#) = Out of Range

AO030803.D A227\_1UG.M

Mon Mar 27 11:03:46 2017

MSD1

Page 1

## Evaluate Continuing Calibration Report

Data File : C:\HPCHEM\1\DATA\AO030803.D

Vial: 3

Acq On : 8 Mar 2017 10:33 am

Operator: RJP

Sample : A1UG\_1.0

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 27 11:00:28 2017

Response via : Multiple Level Calibration

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

Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
51 T	Toluene	0.760	0.831	-9.3	188#	0.00
52 T	Methyl Isobutyl Ketone	0.683	0.530	22.4	141	0.00
53 T	Dibromochloromethane	1.439	1.560	-8.4	189#	0.00
54 T	Methyl Butyl Ketone	0.629	0.478	24.0	144	0.00
55 T	1,2-dibromoethane	0.930	0.924	0.6	170#	0.00
56 T	Tetrachloroethylene	0.691	0.815	-17.9	204#	0.00
57 T	Chlorobenzene	1.238	1.346	-8.7	189#	0.00
58 T	Ethylbenzene	1.745	2.006	-15.0	198#	0.00
59 T	m&p-xylene	1.665	1.964	-18.0	194#	0.00
60 T	Nonane	0.695	0.624	10.2	146	0.00
61 T	Styrene	0.969	1.148	-18.5	206#	0.00
62 T	Bromoform	1.307	1.428	-9.3	184#	0.00
63 T	o-xylene	1.612	1.859	-15.3	187#	0.00
64 T	Cumene	2.227	2.617	-17.5	200#	0.00
65 S	Bromofluorobenzene	0.627	0.662	-5.6	175#	0.00
66 T	1,1,2,2-tetrachloroethane	1.256	1.031	17.9	142	0.00
67 T	Propylbenzene	0.625	0.728	-16.5	194#	0.00
68 T	2-Chlorotoluene	0.528	0.633	-19.9	211#	0.00
69 T	4-ethyltoluene	2.062	2.438	-18.2	204#	0.00
70 T	1,3,5-trimethylbenzene	1.940	2.323	-19.7	197#	0.00
71 T	1,2,4-trimethylbenzene	1.642	2.082	-26.8	220#	0.00
72 T	1,3-dichlorobenzene	1.293	1.536	-18.8	204#	0.00
73 T	benzyl chloride	1.119	1.365	-22.0	208#	0.00
74 T	1,4-dichlorobenzene	1.136	1.366	-20.2	203#	0.00
75 T	1,2,3-trimethylbenzene	1.713	2.104	-22.8	209#	0.00
76 T	1,2-dichlorobenzene	1.297	1.504	-16.0	203#	0.00
77 T	1,2,4-trichlorobenzene	0.468	0.562	-20.1	215#	0.00
78 T	Naphthalene	1.258	1.266	-0.6	194#	0.00
79 T	Hexachloro-1,3-butadiene	1.359	1.375	-1.2	190#	0.00



Data File : C:\HPCHEM\1\DATA\AO030803.D

Vial: 3

Acq On : 8 Mar 2017 10:33 am

Operator: RJP

Sample : A1UG\_1.0

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:13 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	9.64	128	51607	1.00	ppb	-0.02
35) 1,4-difluorobenzene	12.03	114	232876	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	196939	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.53	95	130288	1.06	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	106.00%

## Target Compounds

						Qvalue
2) Propylene	4.54	41	46864	0.87	ppb	67
3) Freon 12	4.62	85	519234	1.09	ppb	97
4) Chloromethane	4.78	50	71090m	0.83	ppb	
5) Freon 114	4.89	85	393406	0.96	ppb	98
6) Vinyl Chloride	5.02	62	97776	0.82	ppb	96
7) Butane	5.21	43	106056	0.86	ppb	# 86
8) 1,3-butadiene	5.16	39	72754m	0.92	ppb	
9) Bromomethane	5.44	94	126042	0.86	ppb	94
10) Chloroethane	5.61	64	41479	0.79	ppb	# 81
11) Ethanol	5.72	45	28685	0.74	ppb	84
12) Acrolein	6.08	56	30071	0.76	ppb	94
13) Vinyl Bromide	5.98	106	128853	0.88	ppb	94
14) Freon 11	6.40	101	599733	1.05	ppb	99
15) Acetone	6.23	58	37904	0.83	ppb	# 32
16) Pentane	6.78	42	43107	0.83	ppb	# 18
17) Isopropyl alcohol	6.46	45	136507	0.82	ppb	# 100
18) 1,1-dichloroethene	7.08	96	71696	1.08	ppb	97
19) Freon 113	7.49	101	198900	1.10	ppb	98
20) t-Butyl alcohol	7.11	59	145313	1.11	ppb	# 93
21) Methylene chloride	7.20	84	65431	0.95	ppb	94
22) Allyl chloride	7.33	41	57216	1.02	ppb	88
23) Carbon disulfide	7.55	76	192233	0.93	ppb	78
24) trans-1,2-dichloroethene	8.24	61	100476	1.08	ppb	93
25) methyl tert-butyl ether	8.54	73	195286m	1.20	ppb	
26) 1,1-dichloroethane	8.47	63	127980	1.04	ppb	93
27) Vinyl acetate	8.62	43	187691	1.54	ppb	96
28) Methyl Ethyl Ketone	8.92	72	30974	1.06	ppb	# 1
29) cis-1,2-dichloroethene	9.44	61	92536	1.13	ppb	95
30) Hexane	9.70	57	69443	1.03	ppb	# 82
31) Ethyl acetate	9.71	43	152501	0.99	ppb	99
32) Chloroform	9.80	83	247201	1.17	ppb	99
33) Tetrahydrofuran	10.32	42	42068	0.94	ppb	# 51
34) 1,2-dichloroethane	10.71	62	173309m	1.24	ppb	
36) 1,1,1-trichloroethane	11.03	97	321582	1.22	ppb	98
37) Cyclohexane	11.95	56	72657	0.91	ppb	# 72
38) Carbon tetrachloride	11.79	117	410561	1.12	ppb	98
39) Benzene	11.61	78	205947	0.99	ppb	92
40) Methyl methacrylate	13.10	41	79152	1.09	ppb	# 88
41) 1,4-dioxane	12.90	58	31932	0.80	ppb	# 22
42) 2,2,4-trimethylpentane	12.91	57	244875	0.89	ppb	# 77
43) Heptane	13.22	43	74984	0.89	ppb	# 76
44) Trichloroethene	12.87	130	138089	1.04	ppb	95
45) 1,2-dichloropropane	12.59	63	58298	0.82	ppb	97

(#)=qualifier out of range (m)=manual integration

AO030803.D A227\_1UG.M

Mon Mar 27 11:03:52 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO030803.D

Vial: 3

Acq On : 8 Mar 2017 10:33 am

Operator: RJP

Sample : A1UG\_1.0

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:13 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

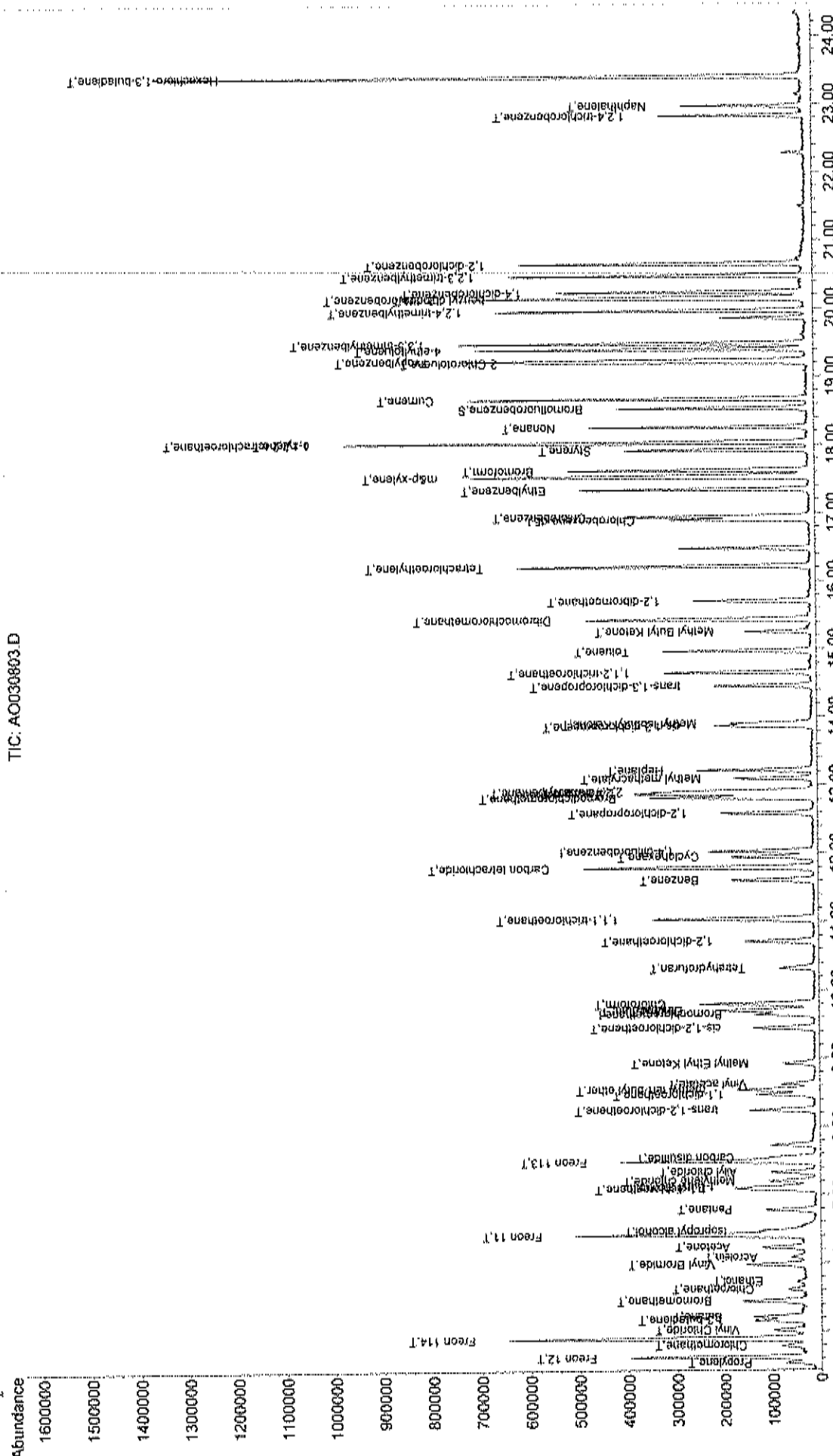
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.81	83	280105	1.11	ppb	96
47) cis-1,3-dichloropropene	13.87	75	122992	1.01	ppb	91
48) trans-1,3-dichloropropene	14.45	75	138346	1.16	ppb	97
49) 1,1,2-trichloroethane	14.66	97	106404	0.94	ppb	98
51) Toluene	14.97	92	163558	1.09	ppb	89
52) Methyl Isobutyl Ketone	13.90	43	104327	0.78	ppb	94
53) Dibromochloromethane	15.44	129	307235	1.08	ppb	93
54) Methyl Butyl Ketone	15.26	43	94051	0.76	ppb	95
55) 1,2-dibromoethane	15.70	107	181895	0.99	ppb	97
56) Tetrachloroethylene	16.22	164	160582	1.18	ppb	99
57) Chlorobenzene	16.94	112	265103	1.09	ppb	94
58) Ethylbenzene	17.34	91	395136	1.15	ppb	96
59) m&p-xylene	17.53	91	773530	2.36	ppb	99
60) Nonane	18.26	43	122957	0.90	ppb	93
61) Styrene	17.92	104	226068	1.18	ppb	90
62) Bromoform	17.62	173	281276	1.09	ppb	98
63) o-xylene	18.04	91	366065	1.15	ppb	97
64) Cumene	18.67	105	515313	1.18	ppb	96
66) 1,1,2,2-tetrachloroethane	18.03	83	203066	0.82	ppb	96
67) Propylbenzene	19.24	120	143415	1.17	ppb #	1
68) 2-Chlorotoluene	19.20	126	124731	1.20	ppb #	1
69) 4-ethyltoluene	19.40	105	480060	1.18	ppb	96
70) 1,3,5-trimethylbenzene	19.48	105	457565	1.20	ppb	96
71) 1,2,4-trimethylbenzene	19.96	105	410061m	1.27	ppb	
72) 1,3-dichlorobenzene	20.16	146	302405	1.19	ppb #	38
73) benzyl chloride	20.13	91	268754	1.22	ppb	91
74) 1,4-dichlorobenzene	20.24	146	268946	1.20	ppb #	25
75) 1,2,3-trimethylbenzene	20.48	105	414336	1.23	ppb	92
76) 1,2-dichlorobenzene	20.65	146	296198	1.16	ppb	94
77) 1,2,4-trichlorobenzene	22.82	180	110615	1.20	ppb	97
78) Naphthalene	22.97	128	249307	1.01	ppb	94
79) Hexachloro-1,3-butadiene	23.40	225	270885	1.01	ppb	100

(#) = qualifier out of range (m) = manual integration (+) = signals summed  
 AO030803.D A227\_1UG.M Mon Mar 27 11:03:52 2017 MSD1

Data File : C:\HPCHEM\1\DATA\AO030803.D  
Acq On : 8 Mar 2017 10:33 am  
Sample : A1UG\_1.0  
Misc : A227\_1UG  
MMS Integration Params: RTEINT.P  
Quant Time: Mar 9 15:45 2017  
Vial: 3  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00  
Quant Results File: A227\_1UG.RES

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Method      : C:\HPCHEM\1\METHODS\A227 1UG.M (RTE Integrator)
Title       : TO-15 VOA Standards for 5 point calibration
Last Update : Mon Mar 27 11:00:28 2017
Response via : Initial Calibration
```

TIC: A0030803.D



**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**RAW DATA**

BFB

Data File : C:\HPCHEM\1\DATA\AO022701.D

Vial: 1

Acq On : 27 Feb 2017 10:36 am

Operator: RJP

Sample : BFB1UG

Inst : MSD #1

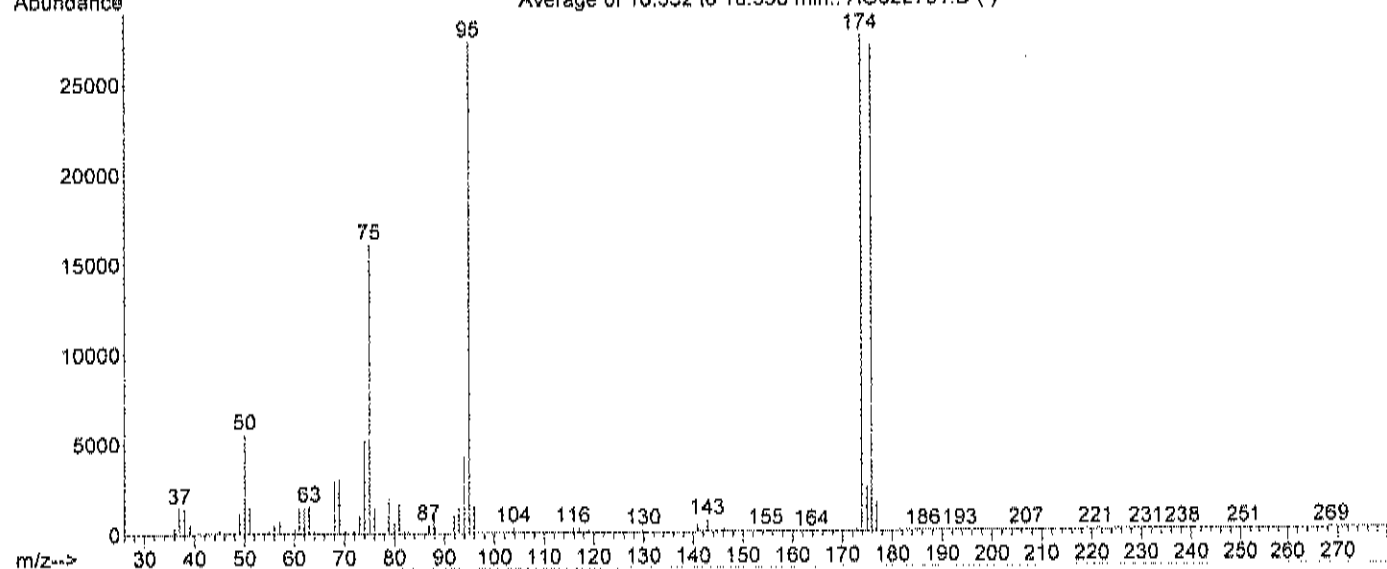
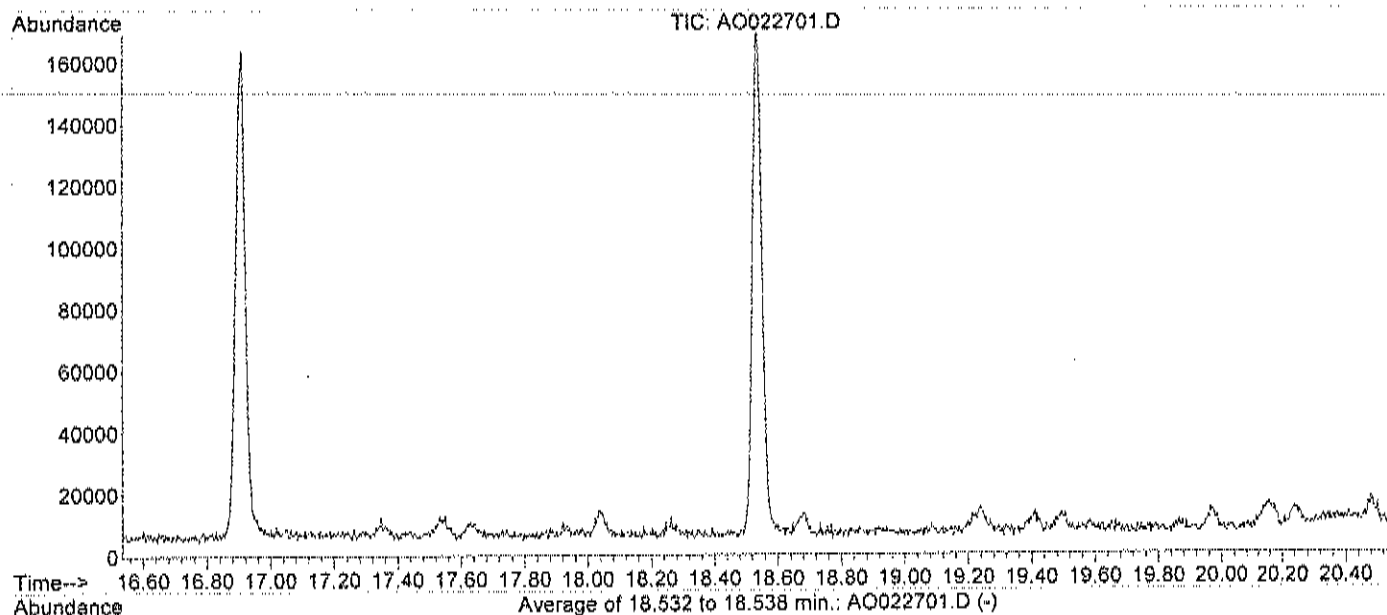
Misc : A120\_1UG - 200cc

Multiplr: 1.00

MS Integration Params: RTEINT.P

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration



Spectrum Information: Average of 18.532 to 18.538 min.

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	8	40	20.4	5563	PASS
75	95	30	66	58.8	16074	PASS
95	95	100	100	100.0	27336	PASS
96	95	5	9	5.4	1473	PASS
173	174	0.00	2	0.4	108	PASS
174	95	50	120	100.9	27589	PASS
175	174	4	9	8.8	2430	PASS
176	174	95	101	98.0	27042	PASS
177	176	5	9	6.1	1638	PASS

BFB

Data File : C:\HPCHEM\1\DATA\AO030801.D

Vial: 1

Acq On : 8 Mar 2017 9:00 am

Operator: RJP

Sample : BFB1UG

Inst : MSD #1

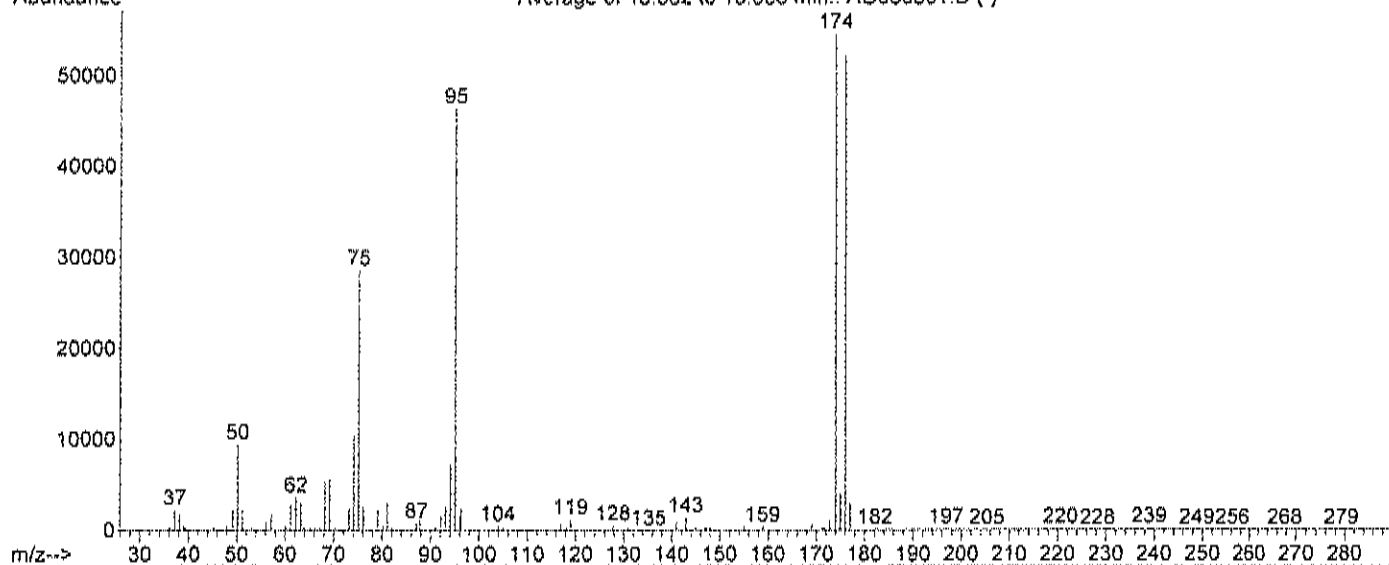
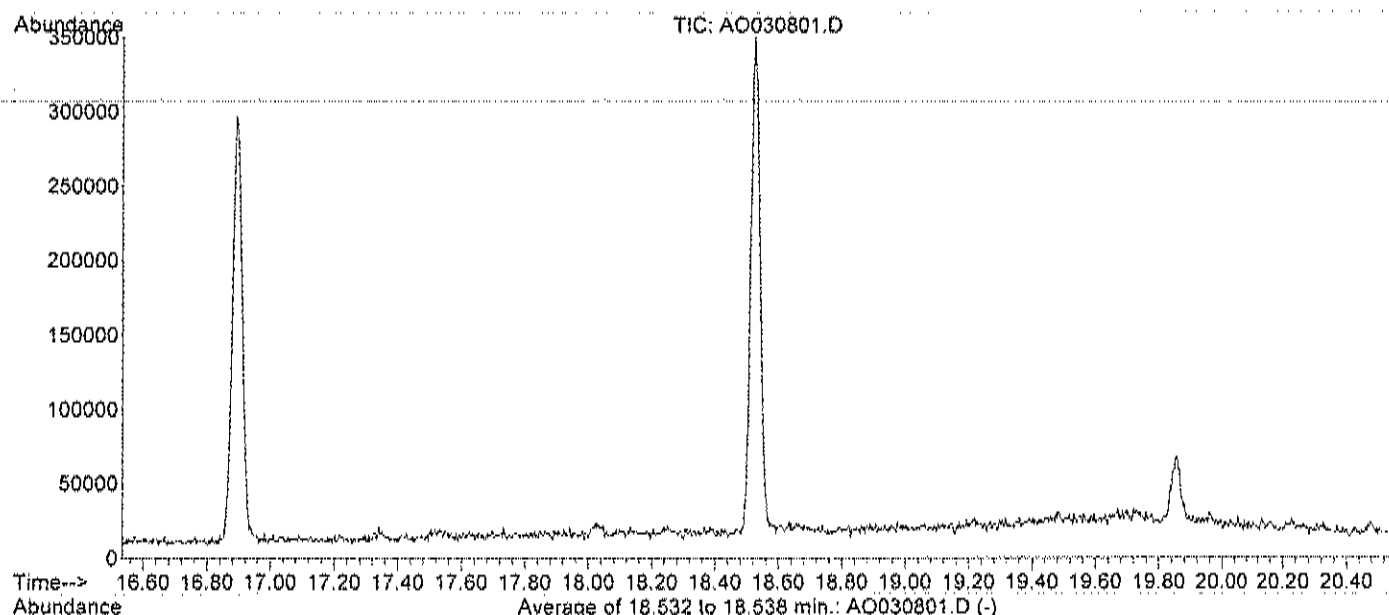
Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration



Spectrum Information: Average of 18.532 to 18.538 min.

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	8	40	20.4	9459	PASS
75	95	30	66	61.6	28567	PASS
95	95	100	100	100.0	46394	PASS
96	95	5	9	5.2	2408	PASS
173	174	0.00	2	2.0	1068	PASS
174	95	50	120	117.6	54570	PASS
175	174	4	9	7.5	4117	PASS
176	174	95	101	95.8	52253	PASS
177	176	5	9	5.6	2913	PASS

**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**RAW QC DATA**

Date: 27-Mar-17

## CENTEK LABORATORIES, LLC

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.

Work Order: C1703015

Project: 691 and 705 St Paul St

TestCode: 0.25CT-TCE-VC

Sample ID	AMB1UG-030817	SampType: MBLK	TestCode: 0.25CT-TCE-	Units: ppbv	Prep Date:	RunNo: 12014					
Client ID:	ZZZZZ	Batch ID: R12014	TestNo: TO-15		Analysis Date: 3/8/2017	SeqNo: 140534					
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	< 0.15	0.15									
Chloroethane	< 0.15	0.15									
cis-1,2-Dichloroethene	< 0.15	0.15									
trans-1,2-Dichloroethene	< 0.15	0.15									
Trichloroethene	< 0.040	0.040									
Vinyl chloride	< 0.040	0.040									

## Qualifiers:

. Results reported are not blank corrected  
 J Analyte detected below quantitation limit  
 S Spike Recovery outside accepted recovery limits

E Estimated Value above quantitation range  
 ND Not Detected at the Limit of Detection

E Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits

Page 1 of 1



Data File : C:\HPCHEM\1\DATA\AO030805.D

Vial: 5

Acq On : 8 Mar 2017 12:00 pm

Operator: RJP

Sample : AMB1UG-030817

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:15 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	9.64	128	51456	1.00	ppb	-0.01
35) 1,4-difluorobenzene	12.03	114	199472	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	169731	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.53	95	90186	0.85	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	85.00%

Target Compounds

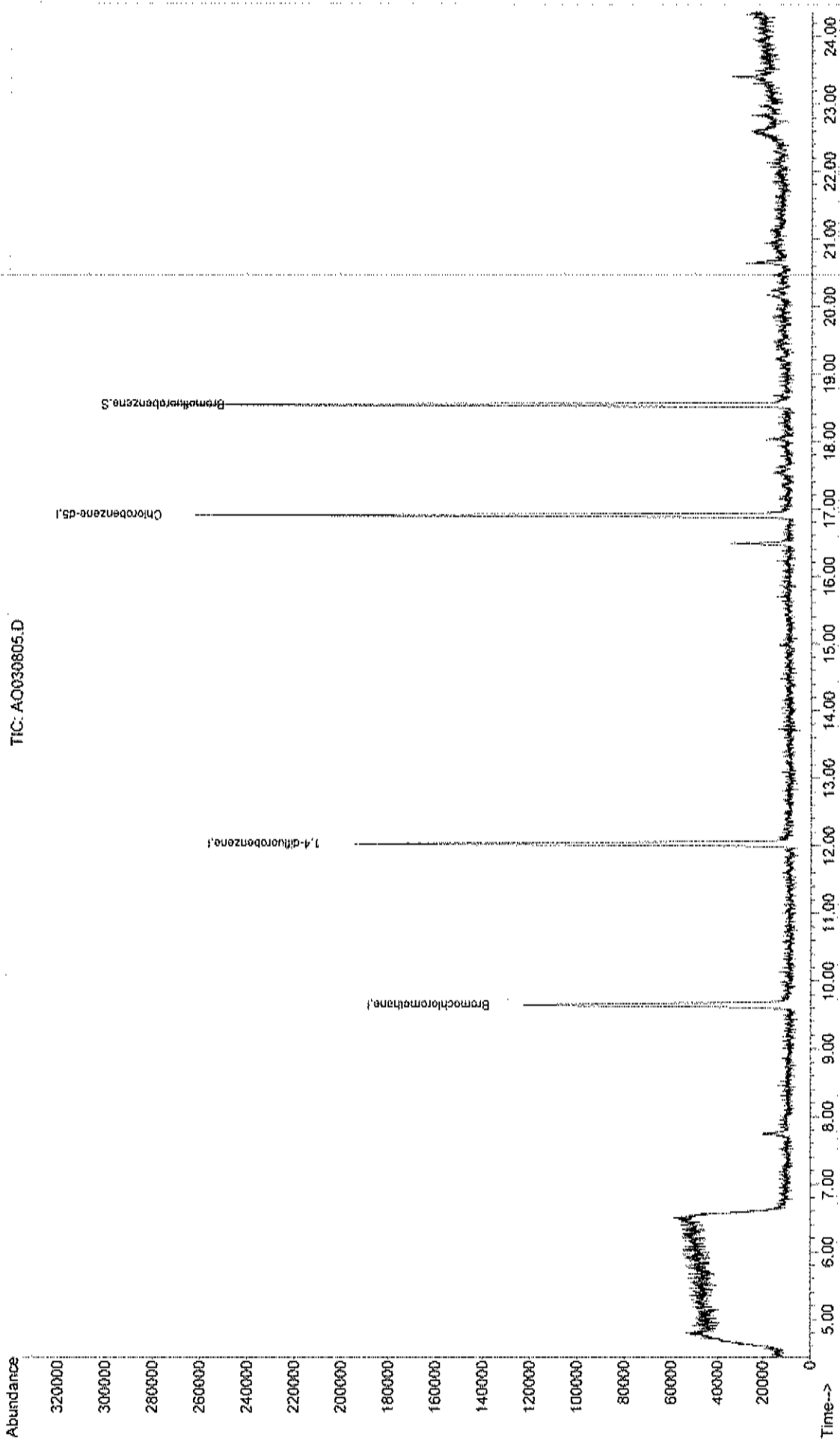
Qvalue

Data File : C:\HPCHEM\1\DATA\A0030805.D  
Acq On : 8 Mar 2017 12:00 pm  
Sample : AMB1UG-030817  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 9 12:10 2017

Vial: 5  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Mar 27 11:00:28 2017  
Response via : Initial Calibration



Date: 27-Mar-17

## CENTEK LABORATORIES, LLC

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.

Work Order: C1703015

Project: 691 and 705 St Paul St

TestCode: 1ugM3\_TO15

Sample ID	C1703015-003A	MS	Sample Type:	MS	TestCode: 1ugM3_TO15	Units: ppbV	Prep Date:	RunNo: 12014			
Client ID:	SV-3	Batch ID: R12014	TestNo: TO-15	Analysis Date: 3/8/2017	SeqNo: 140546						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	1.090	0.15	1	0	109	70	130				
Chloroethane	0.8100	0.15	1	0.11	70.0	70	130				
cis-1,2-Dichloroethene	1.400	0.15	1	0.18	122	70	130				
trans-1,2-Dichloroethene	1.140	0.15	1	0	114	70	130				
Trichloroethene	9.890	0.15	1	8.34	155	70	130				S
Vinyl chloride	0.7200	0.15	1	0	72.0	70	130				

Sample ID	C1703015-003A MS	SampType: MSD	TestCode: 1ugM3_TO15	Units: ppbV	Prep Date:	RunNo: 12014					
Client ID: SV-3	Batch ID: R12014	TestNo: TO-15	Analysis Date: 3/8/2017	SeqNo: 140547							
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	1.160	0.15	1	0	116	70	130	1.09	6.22	30	
Chloroethane	0.8600	0.15	1	0.11	75.0	70	130	0.81	5.99	30	
cis-1,2-Dichloroethene	1.440	0.15	1	0.18	126	70	130	1.4	2.82	30	
trans-1,2-Dichloroethene	1.180	0.15	1	0	118	70	130	1.14	3.45	30	
Trichloroethene	9.550	0.15	1	8.34	121	70	130	9.89	3.50	30	
Vinyl chloride	0.7500	0.15	1	0	75.0	70	130	0.72	4.08	30	

Qualifiers:	Results reported are not blank corrected	E	Estimated Value above quantitation range	H	Holding times for preparation or analysis exceeded
J	Analyte detected below quantitation limit	ND	Not Detected at the Limit of Detection	R	RPD outside accepted recovery limits
S	Spike Recovery outside accepted recovery limits				

Page 1 of 1

Data File : C:\HPCHEM\1\DATA\AO030810.D

Vial: 25

Acq On : 8 Mar 2017 3:32 pm

Operator: RJP

Sample : C1703015-003A MS

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:20 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.65	128	80202	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.04	114	356830	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	344072	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	367606	1.70	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	170.00%#

## Target Compounds

						Qvalue
6) Vinyl Chloride	5.02	62	133363	0.72	ppb	89
10) Chloroethane	5.62	64	66208	0.81	ppb	# 84
18) 1,1-dichloroethene	7.08	96	111749	1.09	ppb	# 88
24) trans-1,2-dichloroethene	8.25	61	165630	1.14	ppb	99
29) cis-1,2-dichloroethene	9.45	61	177576	1.40	ppb	95
44) Trichloroethene	12.88	130	2005119	9.89	ppb	96

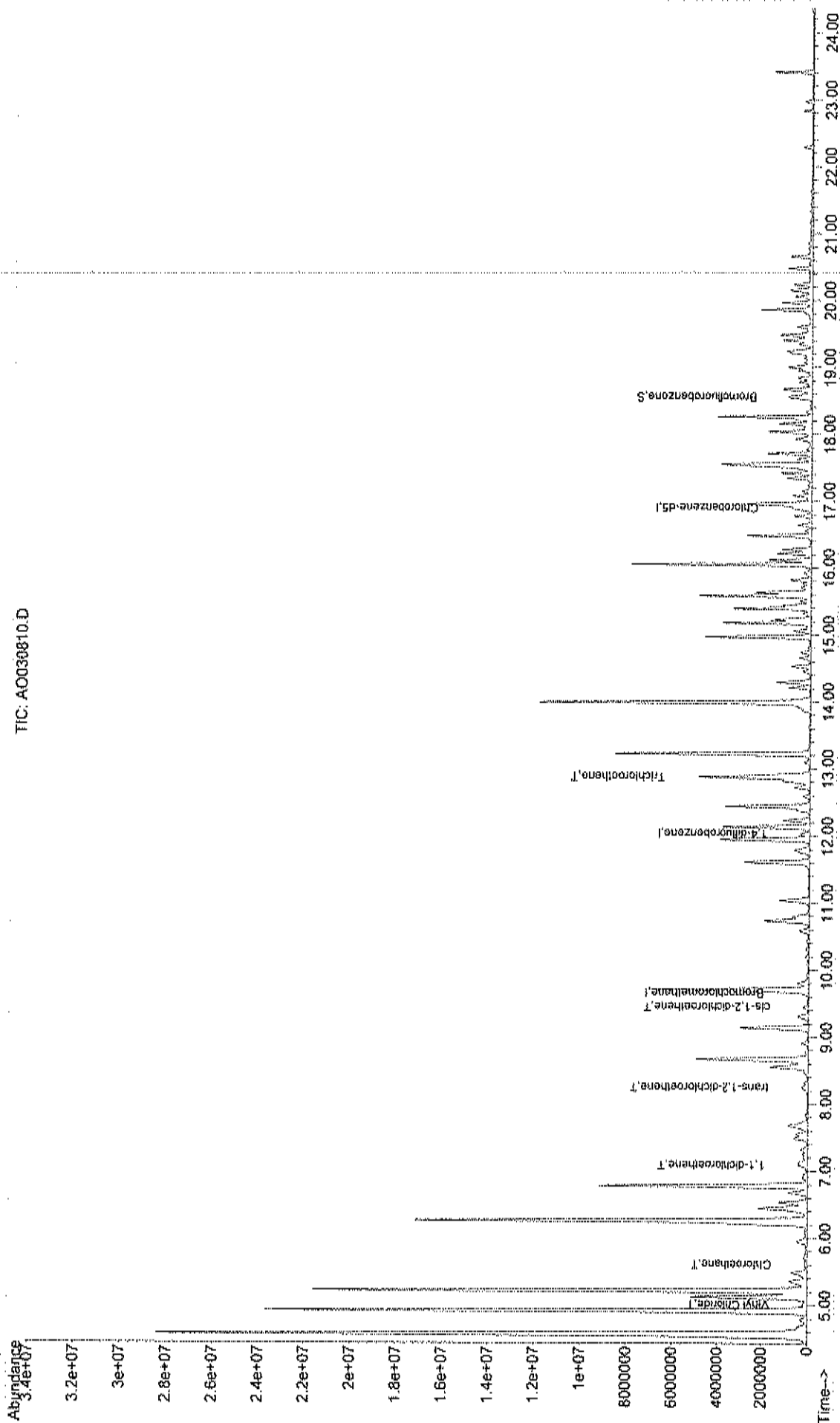
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 (#) = qualifier out of range (m) = manual integration (+) = signals summed  
 AO030810.D A227\_1UG.M Mon Mar 27 11:03:07 2017 MSD1

Data File : C:\HPCHEM\1\DATA\AO030810.D  
Acq On : 8 Mar 2017 3:32 pm  
Sample : C1703015-003A MS  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 10 12:31 2017

Vial: 25  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Mar 27 11:00:28 2017  
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\AO030811.D

Vial: 26

Acq On : 8 Mar 2017 4:18 pm

Operator: RJP

Sample : C1703015-003A MSD

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:21 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	9.66	128	79391	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.03	114	370639	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	337498	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	366795	1.73	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	173.00%#

## Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
6) Vinyl Chloride	5.03	62	137142	0.75	ppb	89
10) Chloroethane	5.63	64	69599	0.86	ppb	93
18) 1,1-dichloroethene	7.09	96	117735	1.16	ppb	92
24) trans-1,2-dichloroethene	8.25	61	169371	1.18	ppb	94
29) cis-1,2-dichloroethene	9.45	61	181684	1.44	ppb	97
44) Trichloroethene	12.87	130	2010156	9.55	ppb	96

Data File : C:\HPCHEM\1\DATA\A0030811.D  
Acq On : 8 Mar 2017 4:18 pm  
Sample : C1703015-003A MSD  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 10 12:36 2017

Vial: 26

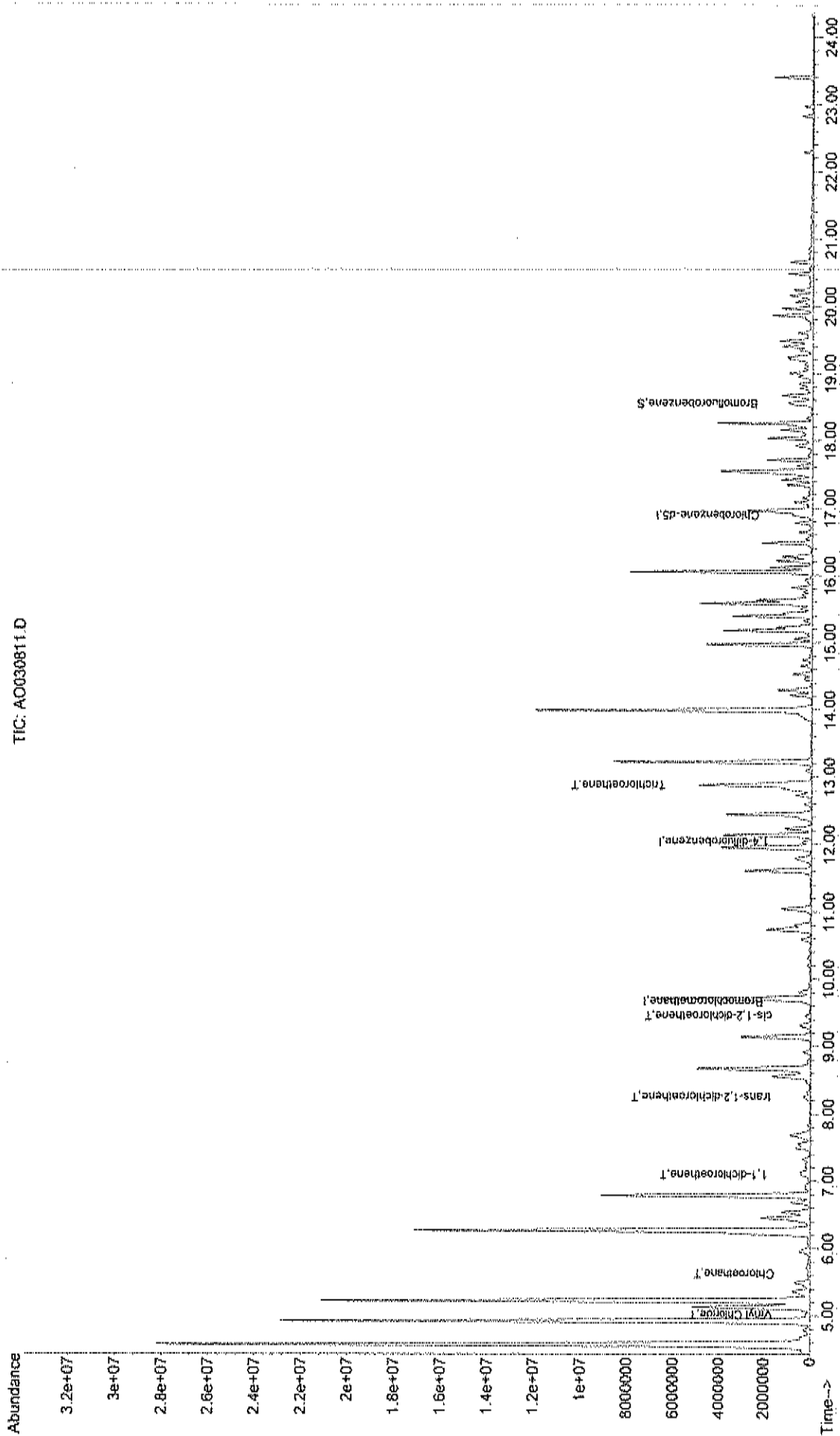
Operator: RJP

Inst : MSD #1

Multiplr: 1.00

Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Mar 27 11:00:28 2017  
Response via : Initial Calibration



Date: 27-Mar-17

## CENTEK LABORATORIES, LLC

## ANALYTICAL QC SUMMARY REPORT

CLIENT: LaBella Associates, P.C.  
 Work Order: C1703015  
 Project: 691 and 705 St Paul St

TestCode: 0.25CT-TCE-VC

Sample ID	ALCS1UG-030817	SampleType: LCS	TestCode: 0.25CT-TCE-	Units: ppbV	Prep Date:	RunNo: 12014					
Client ID: ZZZZZ	Batch ID: R12014	TestNo: TO-15	Analysis Date: 3/8/2017	SeqNo: 140535							
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	1.060	0.15	1	0	106	70	130				
Chloroethane	0.9000	0.15	1	0	90.0	70	130				
cis-1,2-Dichloroethene	1.100	0.15	1	0	110	70	130				
trans-1,2-Dichloroethene	1.070	0.15	1	0	107	70	130				
Trichloroethene	1.060	0.040	1	0	106	70	130				
Vinyl chloride	0.8700	0.040	1	0	87.0	70	130				

Sample ID	ALCS1UGD-030817	SampleType: LCSD	Batch ID: R12014	TestCode: 0.25CT-TCE-	Units: ppbv	Prep Date:	RunNo: 12014				
Client ID: ZZZZZ				TestNo: TO-15		Analysis Date: 3/9/2017	SeqNo: 140536				
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	1.100	0.15	1	0	110	70	130	1.06	3.70	30	
Chloroethane	1.000	0.15	1	0	100	70	130	0.9	10.5	30	
cis-1,2-Dichloroethene	1.130	0.15	1	0	113	70	130	1.1	2.69	30	
trans-1,2-Dichloroethene	1.100	0.15	1	0	110	70	130	1.07	2.76	30	
Trichloroethene	1.100	0.040	1	0	110	70	130	1.06	3.70	30	
Vinyl chloride	0.9100	0.040	1	0	91.0	70	130	0.87	4.49	30	

Qualifiers: J Results reported are not blank corrected  
 S Analyte detected below quantitation limit  
 R Spike Recovery outside accepted recovery limits  
 E Estimated Value above quantitation range  
 ND Not Detected at the limit of Detection  
 H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits

Page 1 of 1



Data File : C:\HPCHEM\1\DATA\AO030804.D

Vial: 4

Acq On : 8 Mar 2017 11:24 am

Operator: RJP

Sample : ALCS1UG-030817

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:14 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	9.65	128	54344	1.00	ppb	-0.01
35) 1,4-difluorobenzene	12.03	114	232438	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	196865	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.53	95	129241	1.05	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	= 105.00%	

## Target Compounds

						Qvalue
2) Propylene	4.55	41	45237	0.80	ppb	# 60
3) Freon 12	4.62	85	538483	1.07	ppb	98
4) Chloromethane	4.79	50	79156m	0.88	ppb	
5) Freon 114	4.89	85	420501	0.98	ppb	98
6) Vinyl Chloride	5.02	62	108615	0.87	ppb	91
7) Butane	5.22	43	121082	0.93	ppb	91
8) 1,3-butadiene	5.16	39	80846m	0.97	ppb	
9) Bromomethane	5.44	94	150085	0.98	ppb	91
10) Chloroethane	5.62	64	49422	0.90	ppb	# 85
11) Ethanol	5.72	45	33607	0.83	ppb	86
12) Acrolein	6.09	56	37853	0.91	ppb	85
13) Vinyl Bromide	5.97	106	156665	1.01	ppb	97
14) Freon 11	6.41	101	622060	1.03	ppb	99
15) Acetone	6.22	58	44784	0.93	ppb	# 56
16) Pentane	6.78	42	44290	0.81	ppb	# 19
17) Isopropyl alcohol	6.47	45	151920	0.87	ppb	# 100
18) 1,1-dichloroethene	7.09	96	74055	1.06	ppb	97
19) Freon 113	7.49	101	203450	1.07	ppb	98
20) t-Butyl alcohol	7.11	59	149440	1.09	ppb	# 95
21) Methylene chloride	7.20	84	66709	0.92	ppb	89
22) Allyl chloride	7.34	41	57833	0.98	ppb	88
23) Carbon disulfide	7.55	76	194545	0.90	ppb	78
24) trans-1,2-dichloroethene	8.24	61	104717	1.07	ppb	95
25) methyl tert-butyl ether	8.54	73	221033	1.28	ppb	98
26) 1,1-dichloroethane	8.46	63	129810	1.00	ppb	94
27) Vinyl acetate	8.62	43	135796m	1.06	ppb	
28) Methyl Ethyl Ketone	8.91	72	29584	0.96	ppb	# 83
29) cis-1,2-dichloroethene	9.45	61	94397	1.10	ppb	92
30) Hexane	9.70	57	72624	1.02	ppb	# 78
31) Ethyl acetate	9.70	43	156257	0.97	ppb	97
32) Chloroform	9.81	83	247209	1.12	ppb	98
33) Tetrahydrofuran	10.32	42	44151	0.94	ppb	# 57
34) 1,2-dichloroethane	10.71	62	189436	1.28	ppb	98
36) 1,1,1-trichloroethane	11.03	97	329169	1.25	ppb	98
37) Cyclohexane	11.95	56	79057	1.00	ppb	# 82
38) Carbon tetrachloride	11.79	117	415954	1.13	ppb	99
39) Benzene	11.61	78	209409	1.01	ppb	93
40) Methyl methacrylate	13.10	41	77526	1.07	ppb	# 83
41) 1,4-dioxane	12.89	58	30085	0.75	ppb	# 16
42) 2,2,4-trimethylpentane	12.91	57	250343	0.91	ppb	# 77
43) Heptane	13.22	43	77615	0.93	ppb	81
44) Trichloroethene	12.87	130	140034	1.06	ppb	94
45) 1,2-dichloropropane	12.59	63	58386	0.82	ppb	97

(#)=qualifier out of range (m)=manual integration

AO030804.D A227\_1UG.M

Mon Mar 27 11:03:00 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO030804.D

Vial: 4

Acq On : 8 Mar 2017 11:24 am

Operator: RJP

Sample : ALCS1UG-030817

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:14 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.82	83	282463	1.12	ppb	97
47) cis-1,3-dichloropropene	13.87	75	128833	1.06	ppb	93
48) trans-1,3-dichloropropene	14.46	75	142832	1.20	ppb	97
49) 1,1,2-trichloroethane	14.66	97	106228	0.94	ppb	98
51) Toluene	14.97	92	164977	1.10	ppb	87
52) Methyl Isobutyl Ketone	13.90	43	106895	0.80	ppb	93
53) Dibromochloromethane	15.44	129	307301	1.08	ppb	93
54) Methyl Butyl Ketone	15.25	43	88414m	0.71	ppb	
55) 1,2-dibromoethane	15.71	107	181960	0.99	ppb	95
56) Tetrachloroethylene	16.22	164	160622	1.18	ppb	98
57) Chlorobenzene	16.94	112	266095	1.09	ppb	96
58) Ethylbenzene	17.35	91	403144	1.17	ppb	97
59) m&p-xylene	17.53	91	772113	2.36	ppb	100
60) Nonane	18.25	43	123526	0.90	ppb	93
61) Styrene	17.92	104	226633	1.19	ppb	92
62) Bromoform	17.62	173	287231	1.12	ppb	99
63) o-xylene	18.04	91	367128	1.16	ppb	95
64) Cumene	18.67	105	509222	1.16	ppb	97
66) 1,1,2,2-tetrachloroethane	18.04	83	201677	0.82	ppb	90
67) Propylbenzene	19.24	120	144853	1.18	ppb	# 1
68) 2-Chlorotoluene	19.21	126	120955	1.16	ppb	# 1
69) 4-ethyltoluene	19.40	105	482912	1.19	ppb	98
70) 1,3,5-trimethylbenzene	19.49	105	451264	1.18	ppb	100
71) 1,2,4-trimethylbenzene	19.96	105	407582	1.26	ppb	95
72) 1,3-dichlorobenzene	20.16	146	297360	1.17	ppb	97
73) benzyl chloride	20.14	91	263130	1.19	ppb	87
74) 1,4-dichlorobenzene	20.24	146	263735	1.18	ppb	94
75) 1,2,3-trimethylbenzene	20.48	105	404680	1.20	ppb	95
76) 1,2-dichlorobenzene	20.66	146	284291	1.11	ppb	96
77) 1,2,4-trichlorobenzene	22.83	180	112850	1.22	ppb	95
78) Naphthalene	22.97	128	249180	1.01	ppb	95
79) Hexachloro-1,3-butadiene	23.41	225	273829	1.02	ppb	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed  
 AO030804.D A227\_1UG.M Mon Mar 27 11:03:01 2017 MSD1

Data File : C:\HPCHEM\1\DATA\AO030804.D  
Acq On : 8 Mar 2017 11:24 am  
Sample : ALCS1UG-030817  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 9 15:46 2017

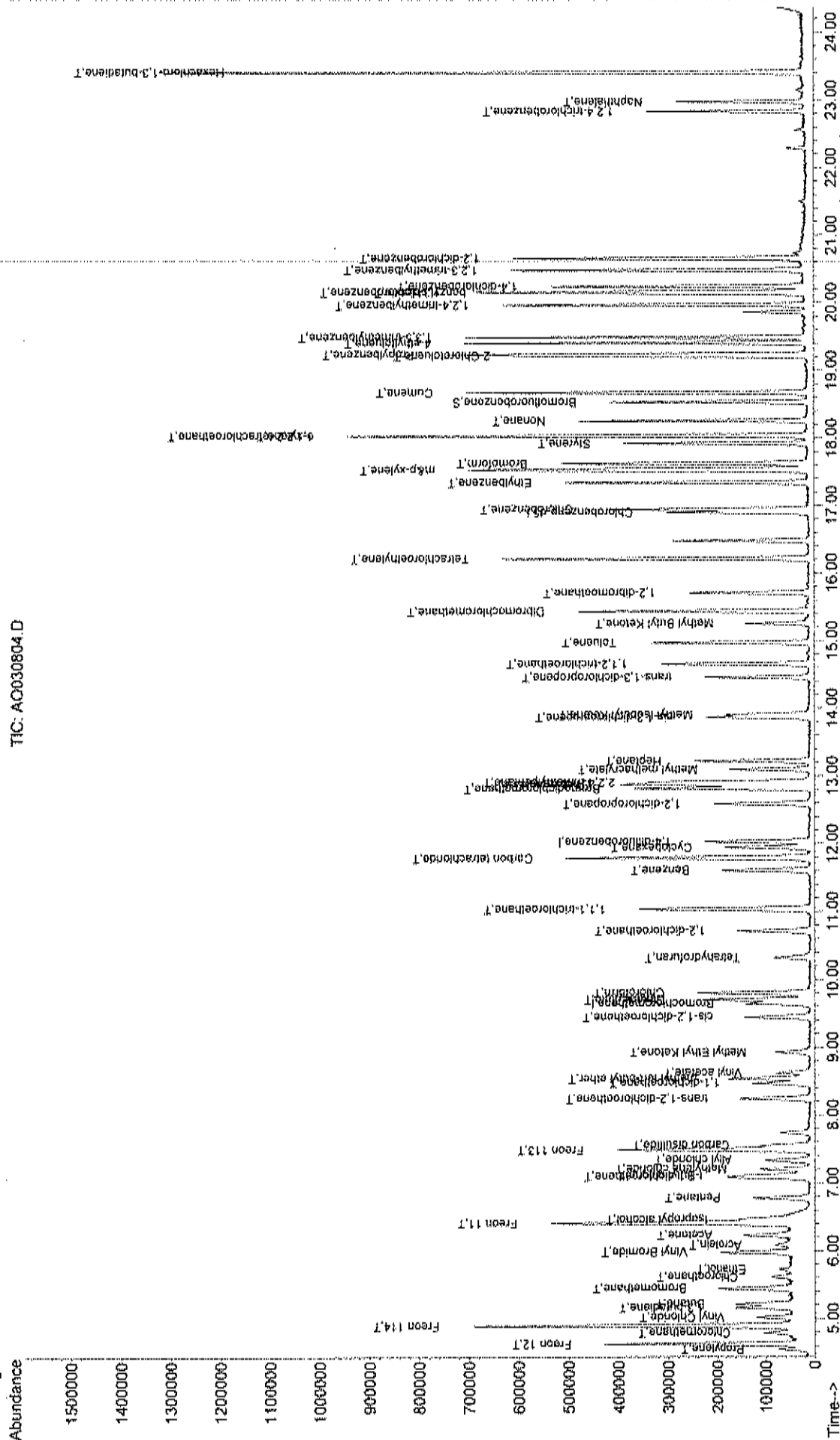
Vial: 4  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Mar 27 11:00:28 2017  
Response via : Initial Calibration

Abundance

TIC: AO030804.D



Data File : C:\HPCHEM\1\DATA\AO030828.D

Vial: 42

Acq On : 9 Mar 2017 3:17 am

Operator: RJP

Sample : ALCS1UGD-030817

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:38 2017

Quant Results File: A227\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.65	128	45106	1.00	ppb	-0.01
35) 1,4-difluorobenzene	12.03	114	199550	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	175594	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.53	95	114997	1.04	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	104.00%

## Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Propylene	4.55	41	39639	0.84	ppb	70
3) Freon 12	4.62	85	491685	1.18	ppb	98
4) Chloromethane	4.79	50	67019m $\beta$	0.90	ppb	
5) Freon 114	4.89	85	377457	1.06	ppb	98
6) Vinyl Chloride	5.02	62	94793	0.91	ppb	94
7) Butane	5.22	43	105897	0.98	ppb	90
8) 1,3-butadiene	5.16	39	74542m $\beta$	1.08	ppb	
9) Bromomethane	5.44	94	133551	1.05	ppb	88
10) Chloroethane	5.61	64	45895	1.00	ppb	91
11) Ethanol	5.71	45	28285	0.84	ppb	# 81
12) Acrolein	6.09	56	32467	0.94	ppb	85
13) Vinyl Bromide	5.97	106	131346	1.02	ppb	92
14) Freon 11	6.40	101	575654	1.15	ppb	98
15) Acetone	6.23	58	37110	0.93	ppb	# 38
16) Pentane	6.77	42	38902	0.85	ppb	# 22
17) Isopropyl alcohol	6.46	45	130448	0.90	ppb	# 100
18) 1,1-dichloroethene	7.08	96	63717	1.10	ppb	97
19) Freon 113	7.49	101	179775	1.13	ppb	98
20) t-Butyl alcohol	7.11	59	114628	1.01	ppb	# 93
21) Methylene chloride	7.21	84	58332	0.97	ppb	92
22) Allyl chloride	7.33	41	50490	1.03	ppb	88
23) Carbon disulfide	7.55	76	181827	1.01	ppb	82
24) trans-1,2-dichloroethene	8.25	61	89836	1.10	ppb	94
25) methyl tert-butyl ether	8.54	73	181878	1.27	ppb	97
26) 1,1-dichloroethane	8.47	63	115379	1.07	ppb	95
27) Vinyl acetate	8.63	43	110010m $\beta$	1.03	ppb	
28) Methyl Ethyl Ketone	8.93	72	25499	1.00	ppb	# 1
29) cis-1,2-dichloroethene	9.45	61	80954	1.13	ppb	90
30) Hexane	9.70	57	64532	1.10	ppb	# 81
31) Ethyl acetate	9.71	43	136266	1.02	ppb	97
32) Chloroform	9.80	83	221863	1.21	ppb	96
33) Tetrahydrofuran	10.32	42	38670	0.99	ppb	# 62
34) 1,2-dichloroethane	10.71	62	167113	1.36	ppb	95
36) 1,1,1-trichloroethane	11.04	97	292539	1.29	ppb	96
37) Cyclohexane	11.95	56	68437	1.01	ppb	# 79
38) Carbon tetrachloride	11.79	117	374325	1.19	ppb	99
39) Benzene	11.61	78	185808	1.05	ppb	91
40) Methyl methacrylate	13.10	41	69019m $\beta$	1.11	ppb	
41) 1,4-dioxane	12.90	58	27466	0.80	ppb	# 24
42) 2,2,4-trimethylpentane	12.92	57	222268	0.94	ppb	# 75
43) Heptane	13.22	43	70110	0.98	ppb	# 75
44) Trichloroethene	12.88	130	124500	1.10	ppb	94
45) 1,2-dichloropropane	12.59	63	54520	0.89	ppb	97

(#)=qualifier out of range (m)=manual integration

AO030828.D A227\_1UG.M

Mon Mar 27 11:03:13 2017

MSD1

Page 1

Data File : C:\HPCHEM\1\DATA\AO030828.D

Vial: 42

Acq On : 9 Mar 2017 3:17 am

Operator: RJP

Sample : ALCS1UGD-030817

Inst : MSD #1

Misc : A227\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Mar 09 11:10:38 2017

Quant Results File: A227\_1UG.REB

Quant Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Mar 06 15:15:36 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Bromodichloromethane	12.81	83	252724	1.16	ppb	97
47) cis-1,3-dichloropropene	13.87	75	111902	1.07	ppb	95
48) trans-1,3-dichloropropene	14.46	75	122008	1.19	ppb	98
49) 1,1,2-trichloroethane	14.66	97	97565	1.00	ppb	98
51) Toluene	14.97	92	146157	1.09	ppb	88
52) Methyl Isobutyl Ketone	13.90	43	80273	0.67	ppb	96
53) Dibromochloromethane	15.44	129	275186	1.09	ppb	92
54) Methyl Butyl Ketone	15.25	43	55926	0.51	ppb	98
55) 1,2-dibromoethane	15.71	107	161381	0.99	ppb	95
56) Tetrachloroethylene	16.22	164	144159	1.19	ppb	98
57) Chlorobenzene	16.95	112	236772	1.09	ppb	96
58) Ethylbenzene	17.34	91	354727	1.16	ppb	95
59) m&p-xylene	17.54	91	688068	2.35	ppb	98
60) Nonane	18.26	43	116714	0.96	ppb	96
61) Styrene	17.92	104	197514	1.16	ppb	90
62) Bromoform	17.62	173	250414	1.09	ppb	100
63) o-xylene	18.04	91	331452	1.17	ppb	96
64) Cumene	18.68	105	450851	1.15	ppb	97
66) 1,1,2,2-tetrachloroethane	18.03	83	187265	0.85	ppb	94
67) Propylbenzene	19.24	120	126078	1.15	ppb	# 1
68) 2-Chlorotoluene	19.21	126	106543	1.15	ppb	# 1
69) 4-ethyltoluene	19.40	105	422495	1.17	ppb	99
70) 1,3,5-trimethylbenzene	19.49	105	404941	1.19	ppb	99
71) 1,2,4-trimethylbenzene	19.96	105	382574	1.33	ppb	93
72) 1,3-dichlorobenzene	20.16	146	263375	1.16	ppb	99
73) benzyl chloride	20.13	91	224469	1.14	ppb	87
74) 1,4-dichlorobenzene	20.24	146	229005	1.15	ppb	96
75) 1,2,3-trimethylbenzene	20.48	105	349230	1.16	ppb	90
76) 1,2-dichlorobenzene	20.65	146	242933	1.07	ppb	95
77) 1,2,4-trichlorobenzene	22.82	180	84549	1.03	ppb	94
78) Naphthalene	22.97	128	191434	0.87	ppb	95
79) Hexachloro-1,3-butadiene	23.40	225	233885	0.98	ppb	98

Data File : C:\HPCHEM\1\DATA\A0030828.D  
Acq On : 9 Mar 2017 3:17 am  
Sample : ALCS1UGD-030817  
Misc : A227\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Mar 9 15:47 2017

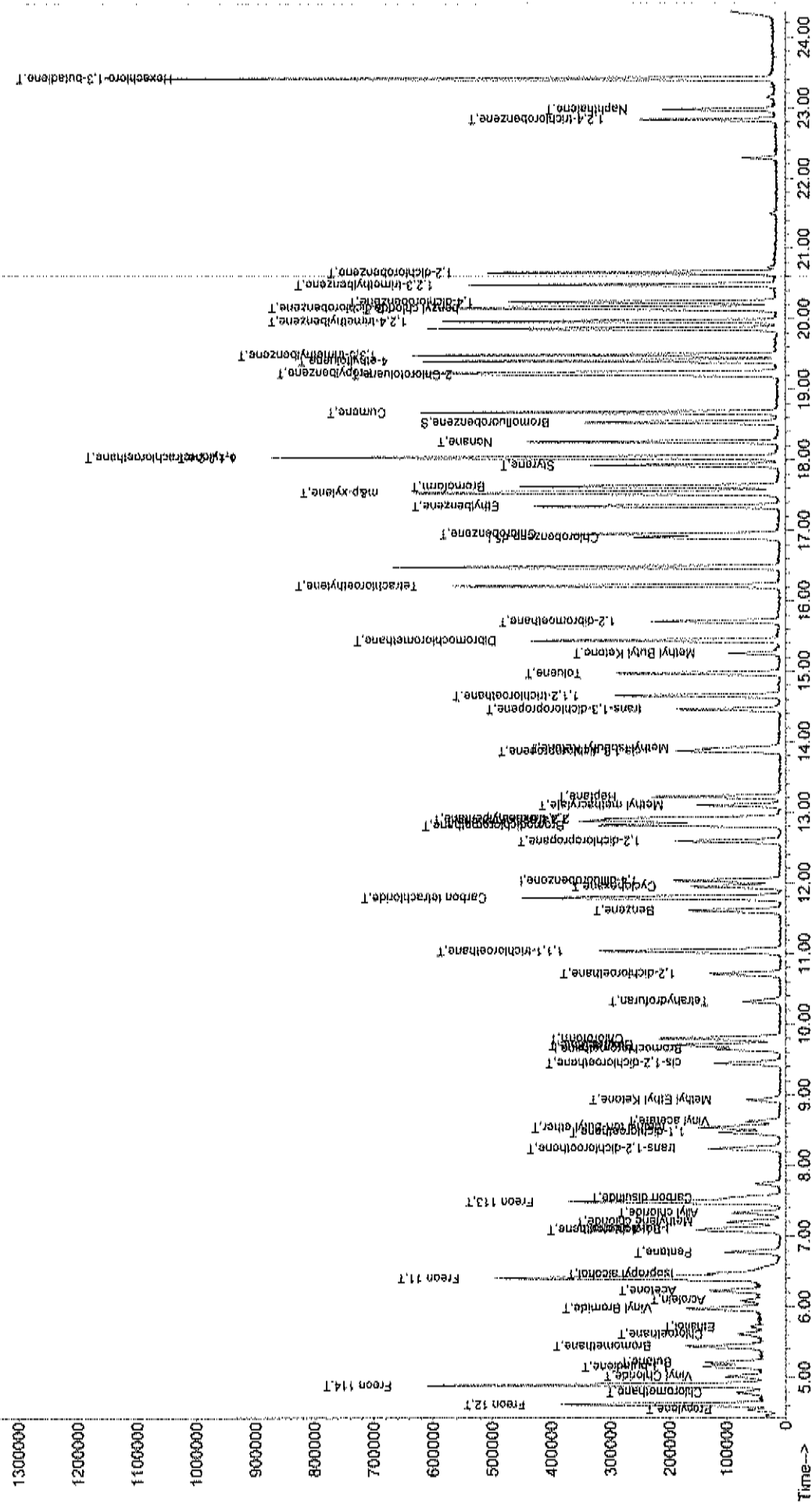
Vial: 42  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A227\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A227\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Mon Mar 27 11:00:28 2017  
Response via : Initial Calibration

Abundance

TIC: A0030828.D



**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**INJECTION LOG**

## Injection Log

Instrument # 1Internal Standard Stock # A1901Standard Stock # A1902LCS Stock # A1903

Directory: C:\HPCHEM\1\DATA

Misc Info: Ref: EPA TO-15 Injected: 1999

ne	Vial	FileName	Multiplier	SampleName		
	1	Ao030601.d	1.	BFB1UG	A227_1UG	6 Mar 2017 08:36
	2	Ao030602.d	1.	A1UG	A227_1UG	6 Mar 2017 09:38
	3	Ao030603.d	1.	A1UG	A227_1UG	6 Mar 2017 10:17
	4	Ao030604.d	1.	ALCS1UG-030617	A227_1UG	6 Mar 2017 10:58
	4	Ao030605.d	1.	AMB1UG-030617	A227_1UG	6 Mar 2017 11:33
	21	Ao030606.d	1.	WAC030617A	A227_1UG	6 Mar 2017 13:43
	22	Ao030607.d	1.	WAC030617B	A227_1UG	6 Mar 2017 14:20
	23	Ao030608.d	1.	WAC030617C	A227_1UG	6 Mar 2017 14:58
	24	Ao030609.d	1.	WAC030617D	A227_1UG	6 Mar 2017 15:36
)	25	Ao030610.d	1.	WAC030617E	A227_1UG	6 Mar 2017 16:13
1	26	Ao030611.d	1.	WAC030617F N	A227_1UG	6 Mar 2017 16:51
2	27	Ao030612.d	1.	WAC030617G	A227_1UG	6 Mar 2017 17:28
3	28	Ao030613.d	1.	WAC030617H	A227_1UG	6 Mar 2017 18:06
4	29	Ao030614.d	1.	WAC030617I	A227_1UG	6 Mar 2017 18:43
5	1	Ao030615.d	1.	C1703007-001A	A227_1UG	6 Mar 2017 19:23
6	2	Ao030616.d	1.	C1703006-001A	A227_1UG	6 Mar 2017 20:03
7	3	Ao030617.d	1.	C1703006-002A	A227_1UG	6 Mar 2017 20:43
8	4	Ao030618.d	1.	C1703006-003A	A227_1UG	6 Mar 2017 21:23
9	5	Ao030619.d	1.	C1703006-004A	A227_1UG	6 Mar 2017 22:03
0	6	Ao030620.d	1.	C1703006-005A	A227_1UG	6 Mar 2017 22:43
1	7	Ao030621.d	1.	C1703005-001A	A227_1UG	6 Mar 2017 23:24
2	8	Ao030622.d	1.	C1703005-002A	A227_1UG	7 Mar 2017 00:06
3	9	Ao030623.d	1.	ALKCS1UGD-030617	A227_1UG	7 Mar 2017 00:45
4	10	Ao030624.d	1.	ALKCS1UGD	A227_1UG	7 Mar 2017 01:25
5	10	Ao030625.d	1.	C1703007-001A 10X	A227_1UG	7 Mar 2017 02:01
6	11	Ao030626.d	1.	C1703006-001A 10X	A227_1UG	7 Mar 2017 02:38
7	12	Ao030627.d	1.	C1703006-002A 10X	A227_1UG	7 Mar 2017 03:15
8	13	Ao030628.d	1.	C1703006-003A 10X	A227_1UG	7 Mar 2017 03:52
9	14	Ao030629.d	1.	C1703006-004A 10X	A227_1UG	7 Mar 2017 04:29
0	15	Ao030630.d	1.	C1703006-005A 10X	A227_1UG	7 Mar 2017 05:06
1	16	Ao030631.d	1.	C1703005-001A 10X	A227_1UG	7 Mar 2017 05:43
2	17	Ao030632.d	1.	C1703005	A227_1UG -001A 40X	7 Mar 2017 06:19
3	18	Ao030633.d	1.	C1703005-002A 10X	A227_1UG	7 Mar 2017 06:56
4	19	Ao030634.d	1.	C1703006-003A 20X	A227_1UG	7 Mar 2017 07:33
5	20	Ao030635.d	1.	C1703006-004A 20X	A227_1UG	7 Mar 2017 08:10
6	30	Ao030636.d	1.	C1703006-004A 270X	A227_1UG	7 Mar 2017 09:15
7	31	Ao030637.d	1.	C1703006-005A 20X	A227_1UG	7 Mar 2017 09:52
8		Ao030638.d	1.	No MS or GC data present		
9	1	Ao030801.d	1.	BFB1UG	A227_1UG	8 Mar 2017 09:00
0	2	Ao030802.d	1.	A1UG	A227_1UG	8 Mar 2017 09:54
1	3	Ao030803.d	1.	A1UG_1.0	A227_1UG	8 Mar 2017 10:33
2	4	Ao030804.d	1.	ALCS1UG-030817	A227_1UG	8 Mar 2017 11:24
3	5	Ao030805.d	1.	AMB1UG-030817	A227_1UG	8 Mar 2017 12:00
4	21	Ao030806.d	1.	C1703015-006A	A227_1UG	8 Mar 2017 12:48
5	22	Ao030807.d	1.	C1703015-001A	A227_1UG	8 Mar 2017 13:28
6	23	Ao030808.d	1.	C1703015-002A	A227_1UG	8 Mar 2017 14:09
7	24	Ao030809.d	1.	C1703015-003A	A227_1UG	8 Mar 2017 14:48
8	25	Ao030810.d	1.	C1703015-003A MS	A227_1UG	8 Mar 2017 15:32
9	26	Ao030811.d	1.	C1703015-003A MSD	A227_1UG	8 Mar 2017 16:18
0	27	Ao030812.d	1.	C1703015-004A	A227_1UG	8 Mar 2017 16:59
1	28	Ao030813.d	1.	C1703015-005A	A227_1UG	8 Mar 2017 17:39
2	29	Ao030814.d	1.	C1703014-001A	A227_1UG	8 Mar 2017 18:20
3	30	Ao030815.d	1.	C1703014-002A	A227_1UG	8 Mar 2017 19:00
4	31	Ao030816.d	1.	C1703014-003A	A227_1UG	8 Mar 2017 19:48
5	32	Ao030817.d	1.	C1703014-001A RE	A227_1UG	8 Mar 2017 20:31



## Injection Log

Directory: C:\HPCHEM\1\DATA

Instrument # 1  
 Internal Standard Stock # A01901  
 Standard Stock # A1902  
 LCS Stock # A1903  
 Injected  
 Method Ref: EPA 8210-10 / Jan. 1999

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
3	32	Ao030818.d	1.	C1703015	A227_1UG -001A 10X	8 Mar 2017 21:07
7	33	Ao030819.d	1.	C1703015	A227_1UG -001A 20X	8 Mar 2017 21:44
8	34	Ao030820.d	1.	C1703015	A227_1UG -002A 10X	8 Mar 2017 22:20
9	35	Ao030821.d	1.	C1703015	A227_1UG -002A 20X	8 Mar 2017 22:57
0	36	Ao030822.d	1.	C1703015-003A 10X	A227_1UG	8 Mar 2017 23:34
1	37	Ao030823.d	1.	C1703013	A227_1UG -003A 10X	9 Mar 2017 00:10
2	38	Ao030824.d	1.	C1703015-004A 10X	A227_1UG	9 Mar 2017 00:47
3	39	Ao030825.d	1.	C1703013	A227_1UG -004A 10X	9 Mar 2017 01:24
4	40	Ao030826.d	1.	C1703015-005A 10X	A227_1UG	9 Mar 2017 02:01
5	41	Ao030827.d	1.	C1703013	A227_1UG -005A 10X	9 Mar 2017 02:37
6	42	Ao030828.d	1.	ALCS1UGD-030817	A227_1UG	9 Mar 2017 03:17
7	43	Ao030829.d	1.	C1703014-001A 10X	A227_1UG	9 Mar 2017 03:54
8	44	Ao030830.d	1.	C1703014-002A 10X	A227_1UG	9 Mar 2017 04:30
9	45	Ao030831.d	1.	C1703014-003A 10X	A227_1UG	9 Mar 2017 05:07
0	46	Ao030832.d	1.	C1703014-002A 90X	A227_1UG	9 Mar 2017 08:55
1	47	Ao030833.d	1.	C1703014-003A 90X	A227_1UG	9 Mar 2017 09:31
2		Ao030834.d	1.	No MS or GC data present		
3	1	Ao030901.d	1.	BFB1UG	A227_1UG	9 Mar 2017 13:43
4	2	Ao030902.d	1.	A1UG_1.0	A227_1UG	9 Mar 2017 14:37
5	3	Ao030903.d	1.	ALCS1UG-030917	A227_1UG	9 Mar 2017 15:35
6	4	Ao030904.d	1.	AMB1UG-030917	A227_1UG	9 Mar 2017 16:11
7	1	Ao030905.d	1.	WAC030917A	A227_1UG	9 Mar 2017 16:48
8	2	Ao030906.d	1.	WAC030917B	A227_1UG	9 Mar 2017 17:25
9	3	Ao030907.d	1.	WAC030917C	A227_1UG	9 Mar 2017 18:03
0	4	Ao030908.d	1.	WAC030917D	A227_1UG	9 Mar 2017 18:40
1	5	Ao030909.d	1.	WAC030917E	A227_1UG	9 Mar 2017 19:17
2	1	Ao030910.d	1.	C1703026-004A	A227_1UG	9 Mar 2017 19:58
3	2	Ao030911.d	1.	C1703026-008A	A227_1UG	9 Mar 2017 20:40
4	3	Ao030912.d	1.	C1703026-001A	A227_1UG	9 Mar 2017 21:20
5	4	Ao030913.d	1.	C1703026-002A	A227_1UG	9 Mar 2017 22:01
6	5	Ao030914.d	1.	C1703026-003A	A227_1UG	9 Mar 2017 22:43
7	6	Ao030915.d	1.	C1703026-005A	A227_1UG	9 Mar 2017 23:25
8	7	Ao030916.d	1.	C1703026-006A	A227_1UG	10 Mar 2017 00:07
9	8	Ao030917.d	1.	C1703026-007A	A227_1UG	10 Mar 2017 00:48
0	9	Ao030918.d	1.	C1703026-004A 10x	A227_1UG	10 Mar 2017 01:25
1	10	Ao030919.d	1.	C1703026	A227_1UG -008A 10x	10 Mar 2017 02:02
2	11	Ao030920.d	1.	C1703026-001A 10x	A227_1UG	10 Mar 2017 02:38
3	12	Ao030921.d	1.	C1703026-002A 10x	A227_1UG	10 Mar 2017 03:15
4	13	Ao030922.d	1.	C1703026-003A 10x	A227_1UG	10 Mar 2017 03:53
5	14	Ao030923.d	1.	ALCS1UGD-030917	A227_1UG	10 Mar 2017 04:33
6	15	Ao030924.d	1.	C1703028-001A	A227_1UG	10 Mar 2017 08:03
7	16	Ao030925.d	1.	C1703028-002A	A227_1UG	10 Mar 2017 08:43
8	17	Ao030926.d	1.	C1703028-003A	A227_1UG	10 Mar 2017 09:23
9	18	Ao030927.d	1.	C1703028-004A	A227_1UG	10 Mar 2017 10:03
0	19	Ao030928.d	1.	C1703026-005A 10x	A227_1UG	10 Mar 2017 10:41
1	20	Ao030929.d	1.	C1703026-006A 10x	A227_1UG	10 Mar 2017 11:18
2	21	Ao030930.d	1.	C1703026-007A 10x	A227_1UG	10 Mar 2017 11:55
3	22	Ao030931.d	1.	C1703028-001A 9x	A227_1UG	10 Mar 2017 12:34
4	23	Ao030932.d	1.	C1703028-002A 20x	A227_1UG	10 Mar 2017 13:11
5	24	Ao030933.d	1.	C1703028-003A 10x	A227_1UG	10 Mar 2017 13:48
6	25	Ao030934.d	1.	C1703028-004A 20x	A227_1UG	10 Mar 2017 14:24
7	26	Ao030935.d	1.	C1703028-001A 90x	A227_1UG	10 Mar 2017 15:01
8		Ao030936.d	1.	No MS or GC data present		
9	1	Ao031201.d	1.	BFB1UG	A311_1UG	12 Mar 2017 12:01
0	2	Ao031202.d	1.	A1UG	A312_1UG	12 Mar 2017 12:52

**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**STANDARDS LOG**

Std #	Date Prep	Date Exp	Description	Stock #	Stock Conc	Initial Vol (psig)	Final Vol (psia)	Final Conc (ppb)	Prep by	Chkd
A-1788	12/22/16	12/29/16	TO15 STLX	<del>A1088</del> A1089	500ppb	3.0	30	50	ZZ	
A-1789			↓	A0270	1ppm	1.5	30	50		
A-1790			↓	A0269	10ppm	1.5	30	500		
A-1791			TO15 ING IS	A1782	50ppb	0.9	45	1		
A-1792			↓	STD	↓	↓	↓	↓		
A-1793			↓	LCS	↓	↓	↓	↓		
A-1794	12/29/16	1/5/17	TO15 IS	A1289	1ppm	1.5	30	50	WD	
<del>A-1795</del>			STD	A1203	↓	↓	↓	↓		
A-1796			LCS	A1204	↓	↓	↓	↓		
A-1797			4PCB	9519	↓	↓	↓	↓		
A-1798			4PCB	A1797	50ppb	3.0	30	5		
A-1799			FORM	A0974	11.5ppm	0.20	45	50		
A-1800			SILOX	<del>A1088</del> A1089	500ppb	3.0	30	50		
A-1801			SULF	A0270	1ppm	1.5	30	50		
A-1802			H2S	A0269	10ppm	1.5	30	500		
A-1803			TO15 ING IS	A1794	50ppb	0.9	45	1		
A-1804			↓	A1795	↓	↓	↓	↓		
A-1805			↓	LCS	↓	↓	↓	↓		
A-1806	1/5/17	1/5/18	TO15 IS	FF-47206	LINDE	LINDE	2000psig	1ppm	WD	
A-1807	1/5/17	1/5/18	STOCK TO15 STD	FF-45347	LINDE	LINDE	2200psig	1ppm	WD	
A-1808	1/6/17	1/6/18	TO15 <del>STD</del> LCS	<del>A1796</del>	1ppm	<del>1.5</del>	<del>30</del>	<del>50ppb</del>	M	

## Centek Laboratories, LLC

## GC/MS Calibration Standards Logbook

Std #	Date Prep	Date Exp	Description	Stock #	Stock Conc	Initial Vol (psig)	Final Vol (psia)	Final Conc (ppb)	Prep by	Chkd by
A-1892	3/3/11	3/10/11	TO15 IS	A1806	1ppm	1.5	30	50	PO	
A-1893			STD	A1807						
A-1894			LCS	A1808						
A-1895			HACH	9519						
A-1896			4PCHS	A1893	50ppm	3.0		5		
A-1897			FORM	A0974	11.5ppm	0.20	45	50		
A-1898			SILOX	<del>A1808</del>	500ppm	3.0	30			
A-1899			SUL	A0270	1ppm	1.5				
A-1900			HAS	A0269	10ppm					
A-1901			TO15/10/6 IS	A1892	50ppm	0.5	45	1ppm		
A-1902			STD	A1893						
A-1903			LCS	A1894						
A-										
A-										
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**GC/MS VOLATILES-WHOLE AIR**

**METHOD TO-15**

**CANISTER CLEANING LOG**

Instrument: Entech 3100

Canister Number	Canister Size	QC Can Number	# of Cycles	Int & Date Cleaned	QC Batch Number	Detection Limits	Leak Test 24hr Int & Date
1202	14L	218	30	11/02/16 RSP	WAC110316A	1ug/m <sup>3</sup> + 0.25	+ 30 11/02/16 RSP
1321							+ 30
1206							+ 30
209							+ 30
218							+ 30
1320		1207			WAC110316B		+ 30
1208							+ 30
1200							+ 30
484							+ 30
1207							+ 30
216		215			WAC110316C		+ 30
1323							+ 30
210							+ 30
487							+ 30
215							+ 30
1319		1203			WAC110316D		+ 30
485							+ 30
483							+ 30
1201							+ 30
1203							+ 30
171	LL	419			WAC110316E		+ 30
193							+ 30
542							+ 30
243							+ 30
419							+ 30

## QC Canister Cleaning Logbook

Centek Laboratories, LLC

Instrument: Entech 3100

Canister Number	Canister Size	QC Can Number	# of Cycles	Int & Date Cleaned	QC Batch Number	Detection Limits	Leak Test 24hr Int & Date
143	1L	569	20	2/16/17	WPC0207A	1ppm ± 0.25	+ 30 2/19/17
1191							+ 30
93							+ 30
563							+ 30
569							+ 30
225		1318					+ 30
131							+ 30
542							+ 30
237							+ 30
1318							+ 30
556		<del>1318</del> 94					+ 30
1184							+ 30
489							+ 30
290							+ 30
<del>1318</del> 94							+ 30
136		362					+ 30
359							+ 30
222							+ 30
1195							+ 30
362							+ 30
233		163					+ 30
595							+ 30
137							+ 30
96							+ 30
163							+ 30

Data File : C:\HPCHEM\1\DATA2\AO021706.D

Vial: 1

Acq On : 17 Feb 2017 6:47 pm

Operator: RJP

Sample : WAC021717A

Inst : MSD #1

Misc : A120\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 18 08:10:52 2017

Quant Results File: A120\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A120\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Jan 30 16:04:21 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	9.64	128	52270	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.02	114	198116	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	175690	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.54	95	99310	0.89	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	89.00%

Target Compounds

Qvalue



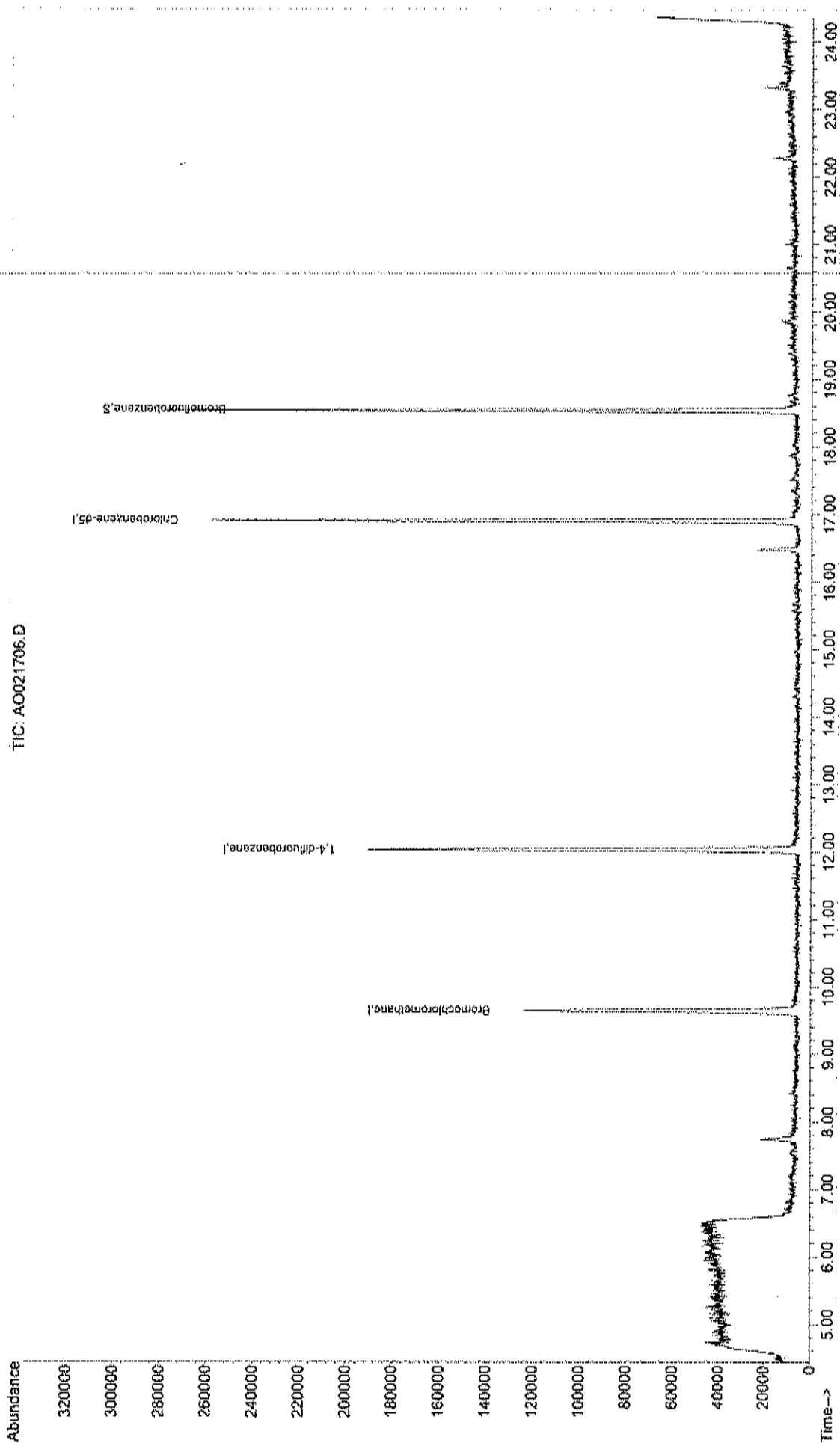
## Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA2\AO021706.D  
Acq On : 17 Feb 2017 6:47 pm  
Sample : WAC021717A  
Misc : A120\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Feb 18 9:10 2017

Vial: 1  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A120\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Wed Mar 15 10:58:20 2017  
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA2\AO021707.D

Vial: 2

Acq On : 17 Feb 2017 7:25 pm

Operator: RJP

Sample : WAC021717B

Inst : MSD #1

Misc : A120\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 18 08:10:53 2017

Quant Results File: A120\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A120\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Jan 30 16:04:21 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.63	128	49106	1.00	ppb	-0.01
35) 1,4-difluorobenzene	12.03	114	191028	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	168917	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.53	95	93615	0.87	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	87.00%

Target Compounds

Qvalue

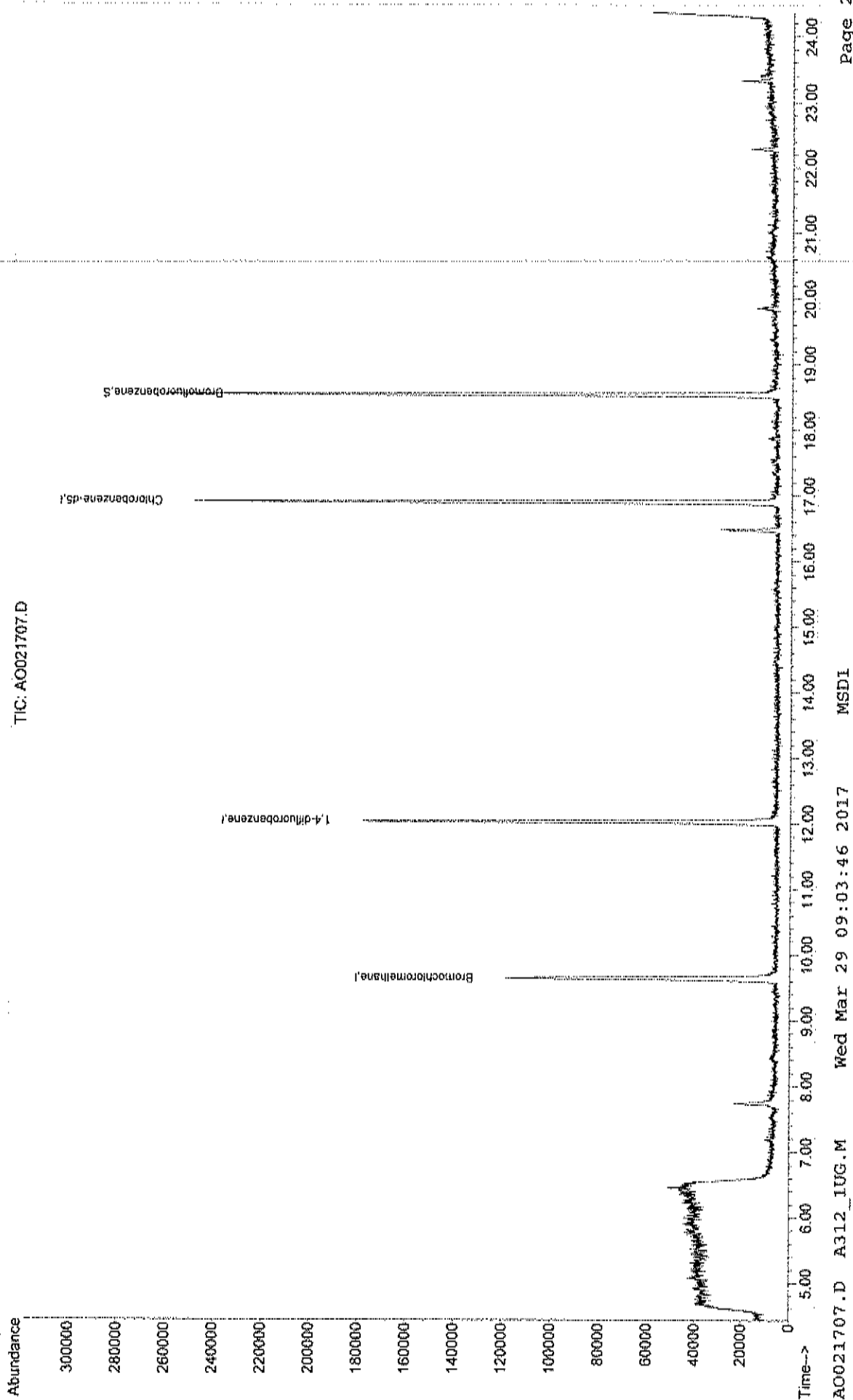
## Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA2\AO021707.D  
Acq On : 17 Feb 2017 7:25 pm  
Sample : WAC021717B  
Misc : A120\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Feb 18 9:10 2017

Vial: 2  
Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A120\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A112\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Wed Mar 15 10:58:20 2017  
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA2\AO021708.D

Vial: 3

Acq On : 17 Feb 2017 8:03 pm

Operator: RJP

Sample : WAC021717C

Inst : MSD #1

Misc : A120\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 18 08:10:54 2017

Quant Results File: A120\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A120\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Jan 30 16:04:21 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	9.64	128	48029	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.02	114	188739	1.00	ppb	0.00
50) Chlorobenzene-d5	16.89	117	164706	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.53	95	93087	0.89	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	89.00%

Target Compounds

Qvalue

## Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA2\AO021708.D  
Acq On : 17 Feb 2017 8:03 pm  
Sample : WAC021717C  
Misc : A120\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Feb 20 16:34 2017

Vial: 3

Operator: RJP

Inst : MSD #1

Multiplr: 1.00

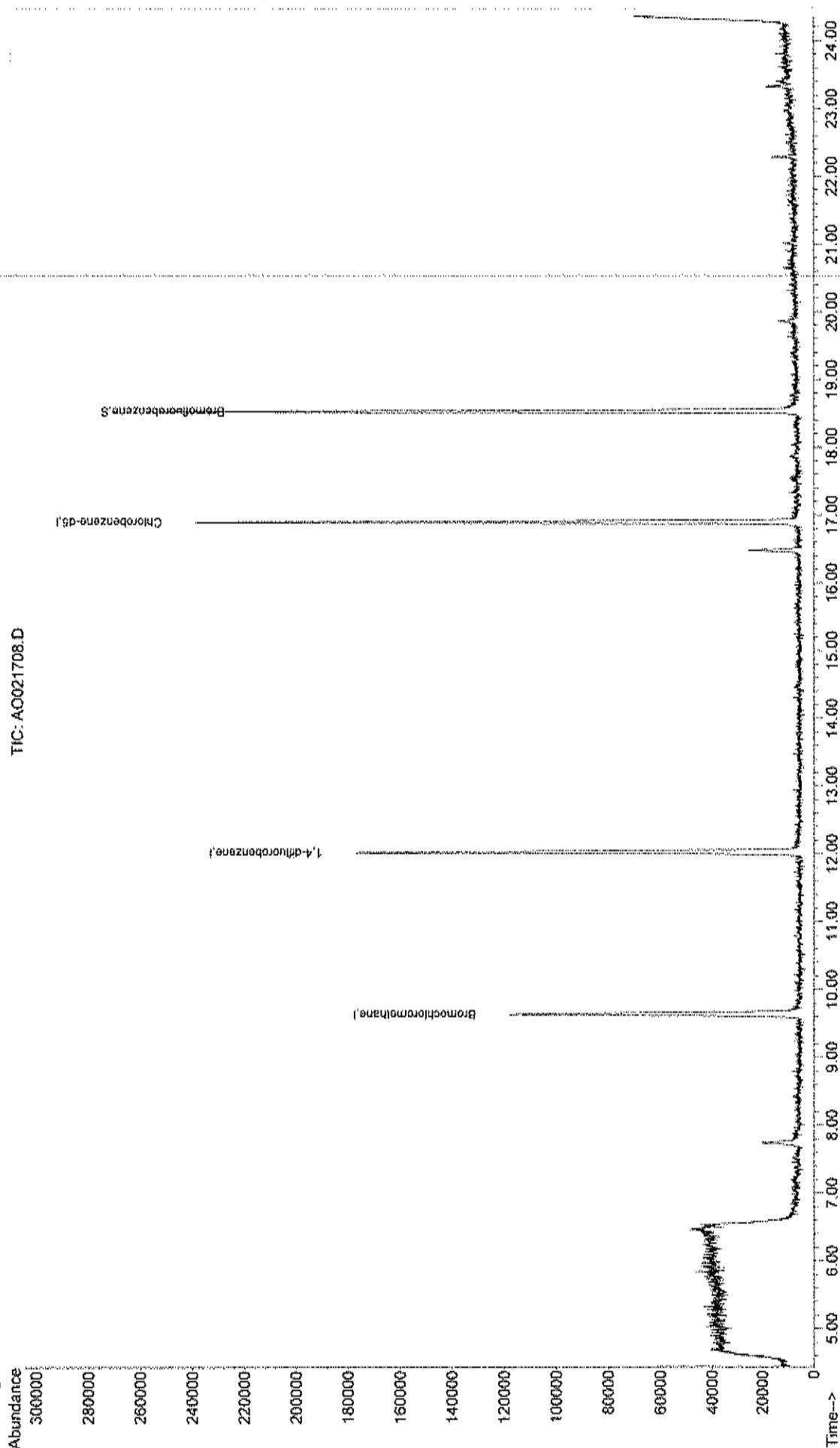
Quant Results File: A120\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Wed Mar 15 10:58:20 2017

Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA2\AO021709.D

Vial: 4

Acq On : 17 Feb 2017 8:41 pm

Operator: RJP

Sample : WAC021717D

Inst : MSD #1

Misc : A120\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 18 08:10:55 2017

Quant Results File: A120\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A120\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Jan 30 16:04:21 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.64	128	47964	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.03	114	178904	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	160775	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.53	95	89341	0.87	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	87.00%

Target Compounds

Qvalue

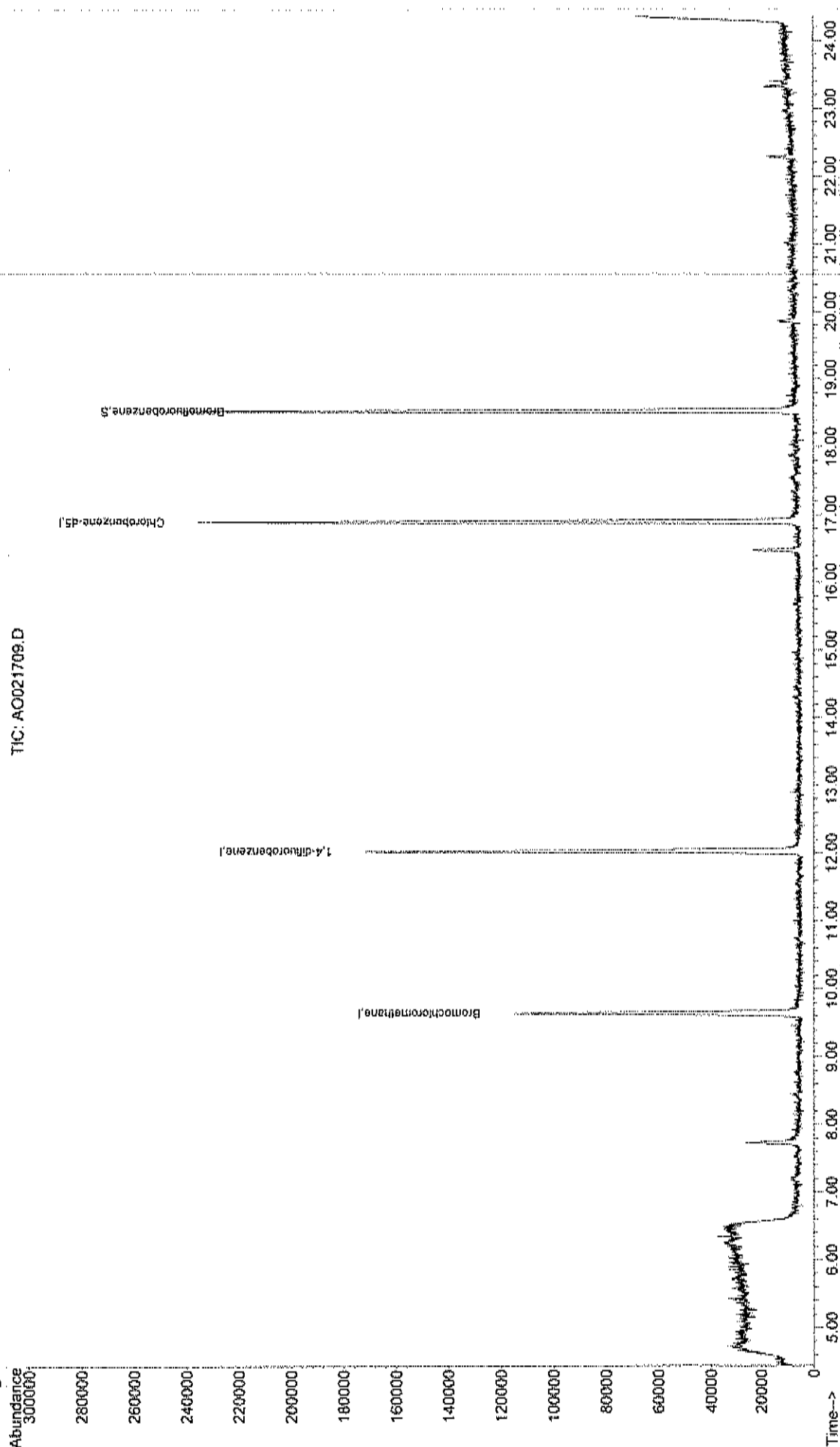
Data File : C:\HPCHEM\1\DATA2\AO021709.D  
Acq On : 17 Feb 2017 8:41 pm  
Sample : WAC021717D  
Misc : A120\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Feb 18 9:10 2017

Vial: 4

Operator: RJP  
Inst : MSD #1  
Multiplr: 1.00

Quant Results File: A120\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Wed Mar 15 10:58:20 2017  
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA2\AO021710.D

Vial: 5

Acq On : 17 Feb 2017 9:19 pm

Operator: RJP

Sample : WAC021717E

Inst : MSD #1

Misc : A120\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 18 08:10:56 2017

Quant Results File: A120\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A120\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Jan 30 16:04:21 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	9.64	128	48138	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.03	114	175951	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	158218	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.53	95	88691	0.88	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	88.00%

Target Compounds

Qvalue



## Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA2\AO021710.D  
Acq On : 17 Feb 2017 9:19 pm  
Sample : WAC021717E  
Misc : A120\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Feb 18 9:10 2017

Vial: 5

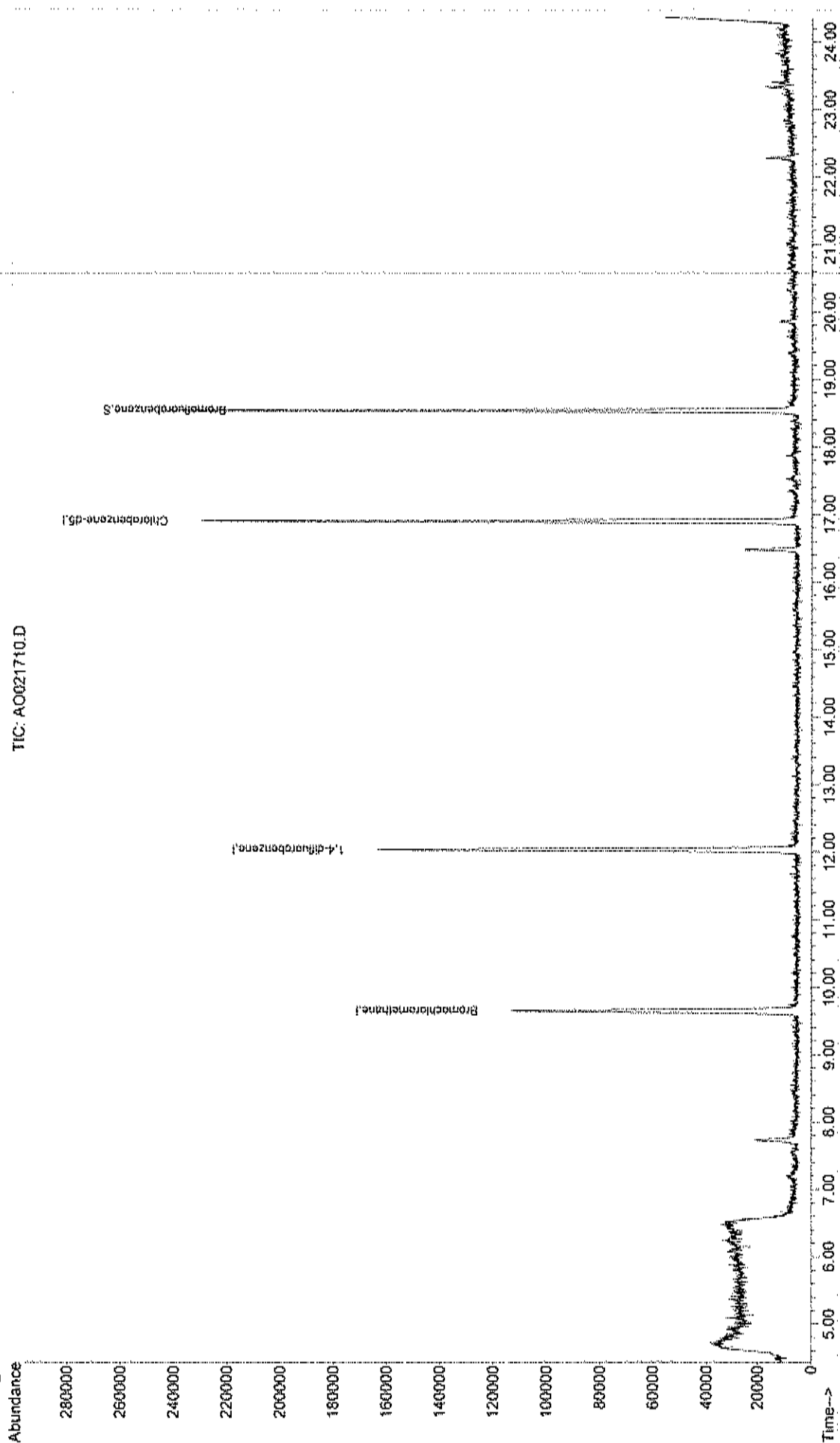
Operator: RJP

Inst : MSD #1

Multiplr: 1.00

Quant Results File: A120\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Wed Mar 15 10:58:20 2017  
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA2\AO021711.D

Vial: 6

Acq On : 17 Feb 2017 9:57 pm

Operator: RJP

Sample : WAC021717F

Inst : MSD #1

Misc : A120\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Feb 18 08:10:57 2017

Quant Results File: A120\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\A120\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Mon Jan 30 16:04:21 2017

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
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1) Bromochloromethane	9.65	128	46520	1.00	ppb	0.00
35) 1,4-difluorobenzene	12.03	114	179438	1.00	ppb	0.00
50) Chlorobenzene-d5	16.90	117	160329	1.00	ppb	0.00

## System Monitoring Compounds

65) Bromofluorobenzene	18.53	95	89777	0.88	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	88.00%

Target Compounds

Qvalue

Data File : C:\HPCHEM\1\DATA2\AO021711.D  
Acq On : 17 Feb 2017 9:57 pm  
Sample : WAC021717F  
Misc : A120\_1UG  
MS Integration Params: RTEINT.P  
Quant Time: Feb 20 16:36 2017

Vial: 6

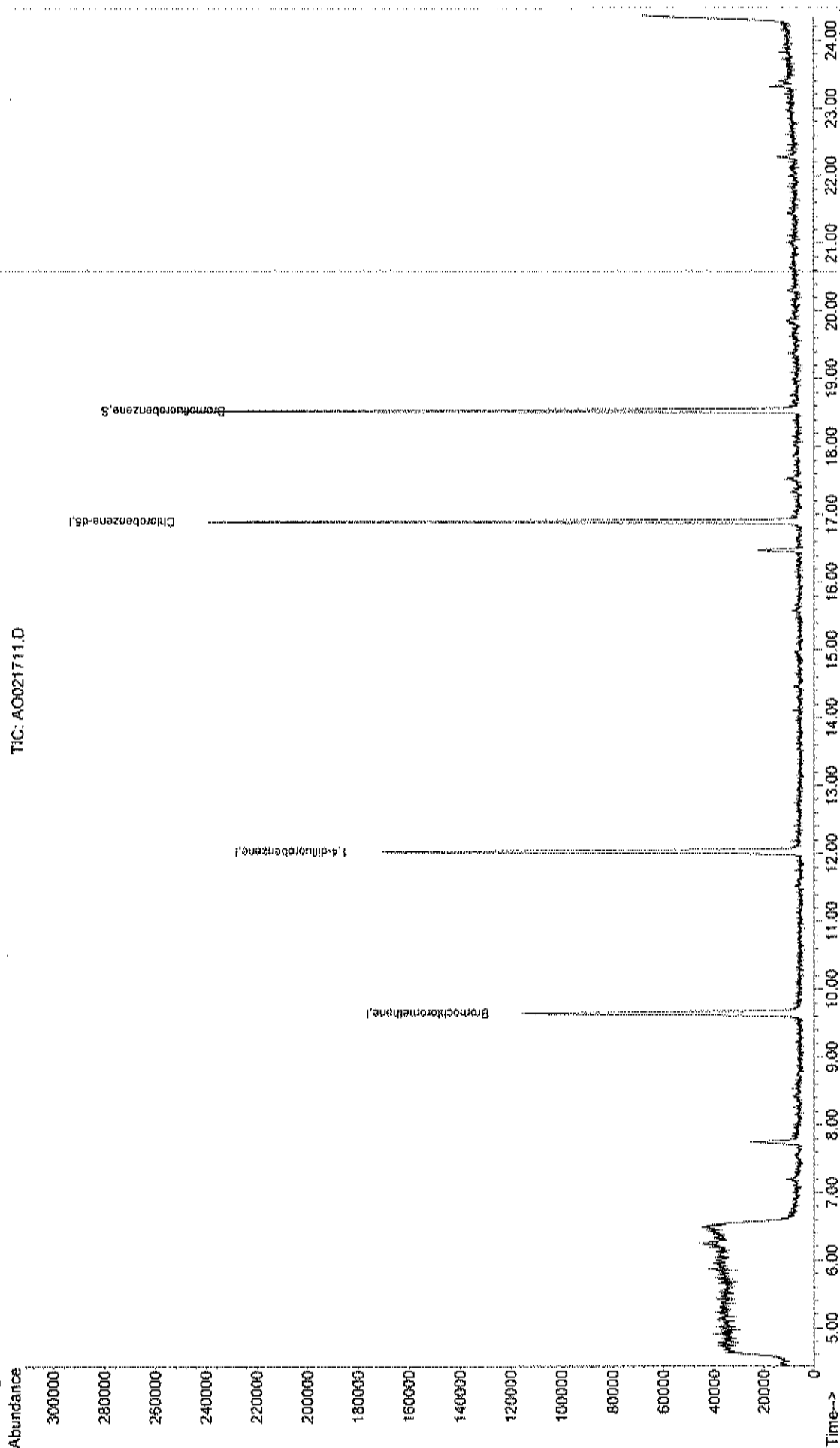
Operator: RJP

Inst : MSD #1

Multiplr: 1.00

Quant Results File: A120\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Wed Mar 15 10:58:20 2017  
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA2\2016NOV\AN110306.D

Vial: 6

Acq On : 3 Nov 2016 12:34 pm

Operator: RJP

Sample : WAC110316A

Inst : MSD #1

Misc : AO26\_1UG

Multiplr: 1.00

MS Integration Params: RTEINT.P

Quant Time: Nov 08 15:11:32 2016

Quant Results File: AO26\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\AO26\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Thu Oct 27 07:19:53 2016

Response via : Initial Calibration

DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	9.55	128	51628	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.85	114	215735	1.00	ppb	0.01
50) Chlorobenzene-d5	16.39	117	192732	1.00	ppb	0.00

## System Monitoring Compounds

66) Bromofluorobenzene	17.96	95	109128	0.84	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	84.00%

Target Compounds

Qvalue

Data File : C:\HPCHEM\1\DATA2\2016NOV\AN110306.D

Acq On : 3 Nov 2016 12:34 pm

Sample : WAC110316A

Misc : AO26\_1UG

MS Integration Params: RTEINT.P

Quant Time: Nov 8 16:11 2016

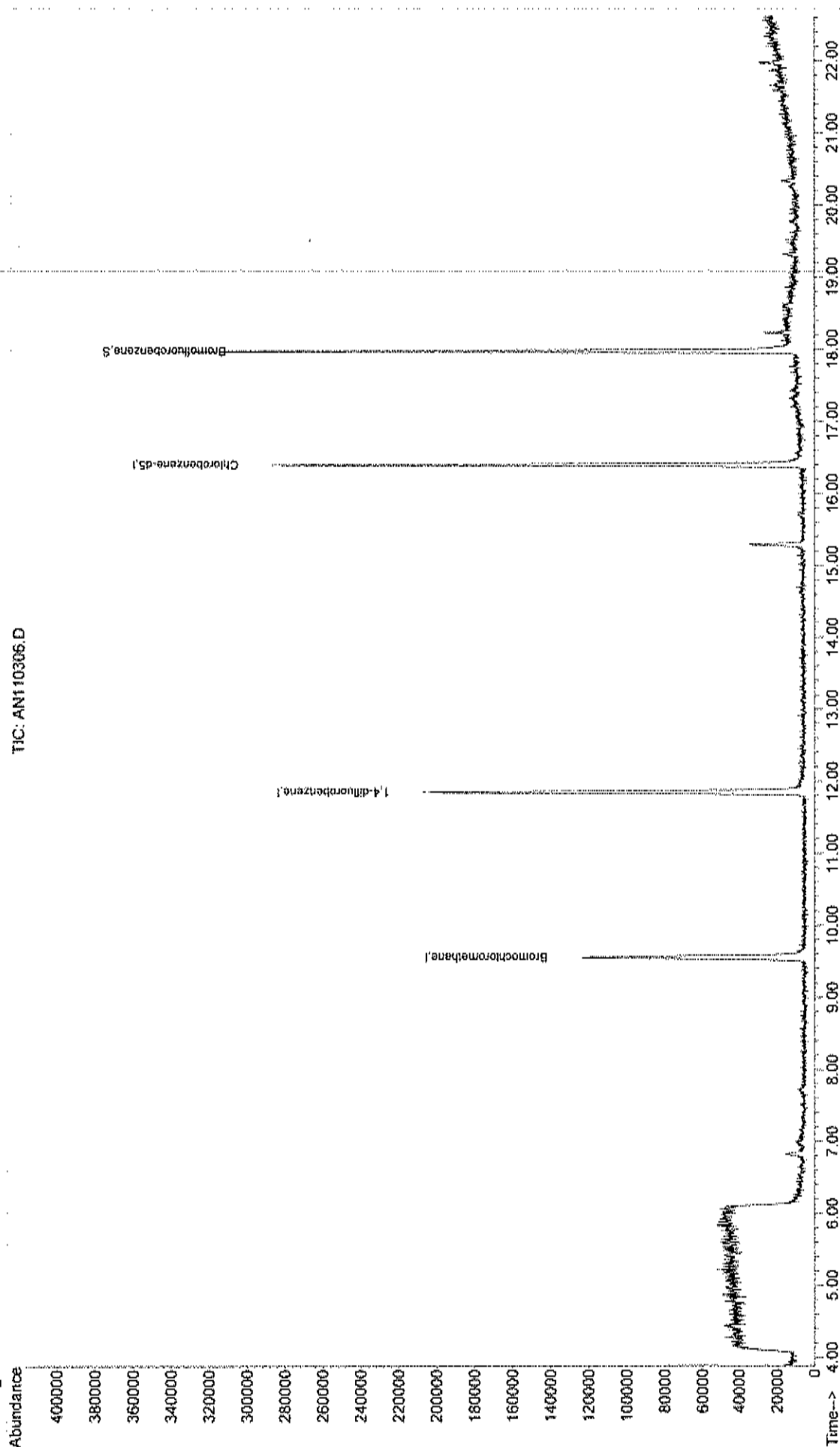
Quant Results File: AO26\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)

Title : TO-15 VOA Standards for 5 point calibration

Last Update : Wed Mar 15 10:58:20 2017

Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA2\2016NOV\AN110307.D Vial: 7  
Acq On : 3 Nov 2016 1:10 pm Operator: RJP  
Sample : WAC110316B Inst : MSD #1  
Misc : AO26\_1UG Multiplr: 1.00  
MS Integration Params: RTEINT.P  
Quant Time: Nov 08 15:11:33 2016 Quant Results File: AO26\_1UG.RES

Quant Method : C:\HPCHEM\1\METHODS\AO26\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Thu Oct 27 07:19:53 2016  
Response via : Initial Calibration  
DataAcq Meth : 1UG\_RUN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	9.56	128	46894	1.00	ppb	0.00
35) 1,4-difluorobenzene	11.84	114	198486	1.00	ppb	0.00
50) Chlorobenzene-d5	16.39	117	177747	1.00	ppb	0.00

## System Monitoring Compounds

66) Bromofluorobenzene	17.96	95	95638	0.80	ppb	0.00
Spiked Amount	1.000	Range	70 - 130	Recovery	=	80.00%

## Target Compounds

Qvalue

Data File : C:\HPCHEM\1\DATA2\2016NOV\AN110307.D Vial: 7  
Acq On : 3 Nov 2016 1:10 pm Operator: RJP  
Sample : WAC110316B Inst : MSD #1  
Misc : AO26\_1UG Multiplr: 1.00  
MS Integration Params: RTEINT.P  
Quant Time: Nov 8 16:11 2016 Quant Results File: AO26\_1UG.RES

Method : C:\HPCHEM\1\METHODS\A312\_1UG.M (RTE Integrator)  
Title : TO-15 VOA Standards for 5 point calibration  
Last Update : Wed Mar 15 10:58:20 2017  
Response via : Initial Calibration

