

## ANALYTICAL REPORT

Job Number: 180-42975-1

Job Description: Harley Davidson

For:

Groundwater Sciences Corporation  
2601 Market Place Street, Suite 310  
Harrisburg, PA 17110-9307

Attention: Allan Miller



Approved for release.  
Carrie L Gamber  
Senior Project Manager  
4/17/2015 12:55 PM

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04/17/2015

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# Definitions/Glossary

Client: Groundwater Sciences Corporation  
Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## Qualifiers

### GC/MS VOA

Qualifier	Qualifier Description
U	Indicates the analyte was analyzed for but not detected.
*	LCS or LCSD exceeds the control limits
F1	MS and/or MSD Recovery exceeds the control limits
J	Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value.
E	Result exceeded calibration range.
B	Compound was found in the blank and sample.

## Glossary

Abbreviation	These commonly used abbreviations may or may not be present in this report.
D	Listed under the "D" column to designate that the result is reported on a dry weight basis
%R	Percent Recovery
CFL	Contains Free Liquid
CNF	Contains no Free Liquid
DER	Duplicate error ratio (normalized absolute difference)
Dil Fac	Dilution Factor
DL, RA, RE, IN	Indicates a Dilution, Re-analysis, Re-extraction, or additional Initial metals/anion analysis of the sample
DLC	Decision level concentration
MDA	Minimum detectable activity
EDL	Estimated Detection Limit
MDC	Minimum detectable concentration
MDL	Method Detection Limit
ML	Minimum Level (Dioxin)
NC	Not Calculated
ND	Not detected at the reporting limit (or MDL or EDL if shown)
PQL	Practical Quantitation Limit
QC	Quality Control
RER	Relative error ratio
RL	Reporting Limit or Requested Limit (Radiochemistry)
RPD	Relative Percent Difference, a measure of the relative difference between two points
TEF	Toxicity Equivalent Factor (Dioxin)
TEQ	Toxicity Equivalent Quotient (Dioxin)

## CASE NARRATIVE

**Client: Groundwater Sciences Corporation**

**Project: Harley Davidson**

**Report Number: 180-42975-1**

With the exceptions noted as flags or footnotes, standard analytical protocols were followed in the analysis of the samples and no problems were encountered or anomalies observed. In addition all laboratory quality control samples were within established control limits, with any exceptions noted below. Each sample was analyzed to achieve the lowest possible reporting limit within the constraints of the method. In some cases, due to interference or analytes present at high concentrations, samples were diluted. For diluted samples, the reporting limits are adjusted relative to the dilution required.

Calculations are performed before rounding to avoid round-off errors in calculated results.

All holding times were met and proper preservation noted for the methods performed on these samples, unless otherwise detailed in the individual sections below.

### **RECEIPT**

The samples were received on 04/11/2015; the samples arrived in good condition, properly preserved and on ice. The temperature of the coolers at receipt was 2.9 C.

### **VOLATILES**

The following sample was diluted to bring the concentration of target analytes within the calibration range: HD-MW-162-0/1-0 (180-42975-2). Elevated reporting limits (RLs) are provided.

Methylene Chloride was detected in method blank MB 180-138685/5 at a level that was above the method detection limit but below the reporting limit. The value should be considered an estimate, and has been flagged. If the associated sample reported a result above the MDL and/or RL, the result has been flagged.

The laboratory control sample (LCS) for batch 138583 recovered outside control limits for the following analytes: Carbon disulfide. A low-level LCS (LLCS), spiked at the reporting limit (RL), was prepared with this batch. The affected target analytes recovered within acceptance limits; therefore, the LLCS demonstrates the analytical system had sufficient sensitivity to detect the compounds had they been present. Since the affected target compounds were not detected in the samples, the data have been reported and qualified.

Carbon disulfide and Trichloroethene failed the recovery criteria low for the MS of sample HD-MW-165-0/1-0 (180-42975-1) in batch 180-138583. Carbon disulfide, Tetrachloroethene and Trichloroethene failed the recovery criteria low for the MSD of sample HD-MW-165-0/1-0 (180-42975-1) in batch 180-138583.

# Detection Summary

Client: Groundwater Sciences Corporation  
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## **Client Sample ID: HD-MW-165-0/1-0**

## **Lab Sample ID: 180-42975-1**

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Trichloroethene	16	F1	1.0	0.14	ug/L	1		8260C	Total/NA
Tetrachloroethene	7.4	F1	1.0	0.15	ug/L	1		8260C	Total/NA

## **Client Sample ID: HD-MW-162-0/1-0**

## **Lab Sample ID: 180-42975-2**

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
1,1-Dichloroethene	1.2	J	4.0	1.2	ug/L	4		8260C	Total/NA
Methylene Chloride	1.4	J	4.0	0.50	ug/L	4		8260C	Total/NA
Trichloroethene	190		4.0	0.57	ug/L	4		8260C	Total/NA
Tetrachloroethene	560	E	4.0	0.59	ug/L	4		8260C	Total/NA
Methylene Chloride - DL	22	J B	40	5.0	ug/L	40		8260C	Total/NA
Trichloroethene - DL	180		40	5.7	ug/L	40		8260C	Total/NA
Tetrachloroethene - DL	700		40	5.9	ug/L	40		8260C	Total/NA

## **Client Sample ID: HD-MW-169-0/1-0**

## **Lab Sample ID: 180-42975-3**

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Methylene Chloride	0.14	J B	1.0	0.13	ug/L	1		8260C	Total/NA

## **Client Sample ID: HD-QC2-0/1-2**

## **Lab Sample ID: 180-42975-4**

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Methylene Chloride	0.22	J	1.0	0.13	ug/L	1		8260C	Total/NA

This Detection Summary does not include radiochemical test results.

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# Client Sample Results

Client: Groundwater Sciences Corporation  
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## Method: 8260C - Volatile Organic Compounds (GC/MS)

**Client Sample ID: HD-MW-165-01-0**

**Date Collected: 04/10/15 08:40**

**Date Received: 04/11/15 09:30**

**Lab Sample ID: 180-42975-1**

**Matrix: Water**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Chloromethane	1.0	U	1.0	0.28	ug/L			04/15/15 15:09	1
Vinyl chloride	1.0	U	1.0	0.23	ug/L			04/15/15 15:09	1
Bromomethane	1.0	U	1.0	0.31	ug/L			04/15/15 15:09	1
Chloroethane	1.0	U	1.0	0.21	ug/L			04/15/15 15:09	1
1,1-Dichloroethene	1.0	U	1.0	0.30	ug/L			04/15/15 15:09	1
Acetone	5.0	U	5.0	2.5	ug/L			04/15/15 15:09	1
Carbon disulfide	1.0	U F1*	1.0	0.21	ug/L			04/15/15 15:09	1
Methylene Chloride	1.0	U	1.0	0.13	ug/L			04/15/15 15:09	1
trans-1,2-Dichloroethene	1.0	U	1.0	0.17	ug/L			04/15/15 15:09	1
Methyl tert-butyl ether	1.0	U	1.0	0.18	ug/L			04/15/15 15:09	1
1,1-Dichloroethane	1.0	U	1.0	0.12	ug/L			04/15/15 15:09	1
cis-1,2-Dichloroethene	1.0	U	1.0	0.24	ug/L			04/15/15 15:09	1
Bromochloromethane	1.0	U	1.0	0.18	ug/L			04/15/15 15:09	1
2-Butanone (MEK)	5.0	U	5.0	0.55	ug/L			04/15/15 15:09	1
Chloroform	1.0	U	1.0	0.17	ug/L			04/15/15 15:09	1
1,1,1-Trichloroethane	1.0	U	1.0	0.29	ug/L			04/15/15 15:09	1
Carbon tetrachloride	1.0	U	1.0	0.14	ug/L			04/15/15 15:09	1
Benzene	1.0	U	1.0	0.11	ug/L			04/15/15 15:09	1
1,2-Dichloroethane	1.0	U	1.0	0.21	ug/L			04/15/15 15:09	1
<b>Trichloroethene</b>	<b>16</b>	<b>F1</b>	1.0	0.14	ug/L			04/15/15 15:09	1
1,2-Dichloropropane	1.0	U	1.0	0.095	ug/L			04/15/15 15:09	1
Bromodichloromethane	1.0	U	1.0	0.13	ug/L			04/15/15 15:09	1
cis-1,3-Dichloropropene	1.0	U	1.0	0.19	ug/L			04/15/15 15:09	1
4-Methyl-2-pentanone (MIBK)	5.0	U	5.0	0.53	ug/L			04/15/15 15:09	1
Toluene	1.0	U	1.0	0.15	ug/L			04/15/15 15:09	1
trans-1,3-Dichloropropene	1.0	U	1.0	0.15	ug/L			04/15/15 15:09	1
1,1,2-Trichloroethane	1.0	U	1.0	0.20	ug/L			04/15/15 15:09	1
<b>Tetrachloroethene</b>	<b>7.4</b>	<b>F1</b>	1.0	0.15	ug/L			04/15/15 15:09	1
2-Hexanone	5.0	U	5.0	0.16	ug/L			04/15/15 15:09	1
Dibromochloromethane	1.0	U	1.0	0.14	ug/L			04/15/15 15:09	1
1,2-Dibromoethane (EDB)	1.0	U	1.0	0.18	ug/L			04/15/15 15:09	1
Chlorobenzene	1.0	U	1.0	0.14	ug/L			04/15/15 15:09	1
1,1,1,2-Tetrachloroethane	1.0	U	1.0	0.28	ug/L			04/15/15 15:09	1
Ethylbenzene	1.0	U	1.0	0.23	ug/L			04/15/15 15:09	1
Xylenes, Total	3.0	U	3.0	0.49	ug/L			04/15/15 15:09	1
Styrene	1.0	U	1.0	0.097	ug/L			04/15/15 15:09	1
Bromoform	1.0	U	1.0	0.19	ug/L			04/15/15 15:09	1
1,1,2,2-Tetrachloroethane	1.0	U	1.0	0.20	ug/L			04/15/15 15:09	1
Acrylonitrile	20	U	20	0.55	ug/L			04/15/15 15:09	1
1,4-Dioxane	200	U	200	34	ug/L			04/15/15 15:09	1
<b>Surrogate</b>	<b>%Recovery</b>	<b>Qualifier</b>	<b>Limits</b>			<b>Prepared</b>	<b>Analyzed</b>	<b>Dil Fac</b>	
1,2-Dichloroethane-d4 (Surr)	118		64 - 135						1
Toluene-d8 (Surr)	100		71 - 118						1
4-Bromofluorobenzene (Surr)	96		70 - 118						1
Dibromofluoromethane (Surr)	110		70 - 128						1

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# Client Sample Results

Client: Groundwater Sciences Corporation  
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## Method: 8260C - Volatile Organic Compounds (GC/MS)

**Client Sample ID: HD-MW-162-0/1-0**

**Date Collected: 04/10/15 09:50**

**Date Received: 04/11/15 09:30**

**Lab Sample ID: 180-42975-2**

**Matrix: Water**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Chloromethane	4.0	U	4.0	1.1	ug/L			04/15/15 20:47	4
Vinyl chloride	4.0	U	4.0	0.91	ug/L			04/15/15 20:47	4
Bromomethane	4.0	U	4.0	1.3	ug/L			04/15/15 20:47	4
Chloroethane	4.0	U	4.0	0.86	ug/L			04/15/15 20:47	4
<b>1,1-Dichloroethene</b>	<b>1.2</b>	<b>J</b>	4.0	1.2	ug/L			04/15/15 20:47	4
Acetone	20	U	20	10	ug/L			04/15/15 20:47	4
Carbon disulfide	4.0	U *	4.0	0.85	ug/L			04/15/15 20:47	4
<b>Methylene Chloride</b>	<b>1.4</b>	<b>J</b>	4.0	0.50	ug/L			04/15/15 20:47	4
trans-1,2-Dichloroethene	4.0	U	4.0	0.68	ug/L			04/15/15 20:47	4
Methyl tert-butyl ether	4.0	U	4.0	0.73	ug/L			04/15/15 20:47	4
1,1-Dichloroethane	4.0	U	4.0	0.47	ug/L			04/15/15 20:47	4
cis-1,2-Dichloroethene	4.0	U	4.0	0.95	ug/L			04/15/15 20:47	4
Bromochloromethane	4.0	U	4.0	0.72	ug/L			04/15/15 20:47	4
2-Butanone (MEK)	20	U	20	2.2	ug/L			04/15/15 20:47	4
Chloroform	4.0	U	4.0	0.68	ug/L			04/15/15 20:47	4
1,1,1-Trichloroethane	4.0	U	4.0	1.1	ug/L			04/15/15 20:47	4
Carbon tetrachloride	4.0	U	4.0	0.55	ug/L			04/15/15 20:47	4
Benzene	4.0	U	4.0	0.42	ug/L			04/15/15 20:47	4
1,2-Dichloroethane	4.0	U	4.0	0.85	ug/L			04/15/15 20:47	4
<b>Trichloroethene</b>	<b>190</b>		4.0	0.57	ug/L			04/15/15 20:47	4
1,2-Dichloropropane	4.0	U	4.0	0.38	ug/L			04/15/15 20:47	4
Bromodichloromethane	4.0	U	4.0	0.52	ug/L			04/15/15 20:47	4
cis-1,3-Dichloropropene	4.0	U	4.0	0.75	ug/L			04/15/15 20:47	4
4-Methyl-2-pentanone (MIBK)	20	U	20	2.1	ug/L			04/15/15 20:47	4
Toluene	4.0	U	4.0	0.60	ug/L			04/15/15 20:47	4
trans-1,3-Dichloropropene	4.0	U	4.0	0.59	ug/L			04/15/15 20:47	4
1,1,2-Trichloroethane	4.0	U	4.0	0.81	ug/L			04/15/15 20:47	4
<b>Tetrachloroethene</b>	<b>560</b>	<b>E</b>	4.0	0.59	ug/L			04/15/15 20:47	4
2-Hexanone	20	U	20	0.64	ug/L			04/15/15 20:47	4
Dibromochloromethane	4.0	U	4.0	0.55	ug/L			04/15/15 20:47	4
1,2-Dibromoethane (EDB)	4.0	U	4.0	0.72	ug/L			04/15/15 20:47	4
Chlorobenzene	4.0	U	4.0	0.54	ug/L			04/15/15 20:47	4
1,1,1,2-Tetrachloroethane	4.0	U	4.0	1.1	ug/L			04/15/15 20:47	4
Ethylbenzene	4.0	U	4.0	0.91	ug/L			04/15/15 20:47	4
Xylenes, Total	12	U	12	2.0	ug/L			04/15/15 20:47	4
Styrene	4.0	U	4.0	0.39	ug/L			04/15/15 20:47	4
Bromoform	4.0	U	4.0	0.77	ug/L			04/15/15 20:47	4
1,1,2,2-Tetrachloroethane	4.0	U	4.0	0.80	ug/L			04/15/15 20:47	4
Acrylonitrile	80	U	80	2.2	ug/L			04/15/15 20:47	4
1,4-Dioxane	800	U	800	140	ug/L			04/15/15 20:47	4
<b>Surrogate</b>	<b>%Recovery</b>	<b>Qualifier</b>	<b>Limits</b>			<b>Prepared</b>	<b>Analyzed</b>	<b>Dil Fac</b>	
1,2-Dichloroethane-d4 (Surr)	115		64 - 135						4
Toluene-d8 (Surr)	98		71 - 118						4
4-Bromofluorobenzene (Surr)	91		70 - 118						4
Dibromofluoromethane (Surr)	116		70 - 128						4

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# Client Sample Results

Client: Groundwater Sciences Corporation  
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## Method: 8260C - Volatile Organic Compounds (GC/MS)

**Client Sample ID: HD-MW-169-0/1-0**

**Date Collected: 04/10/15 12:32**

**Date Received: 04/11/15 09:30**

**Lab Sample ID: 180-42975-3**

**Matrix: Water**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Chloromethane	1.0	U	1.0	0.28	ug/L			04/16/15 16:53	1
Vinyl chloride	1.0	U	1.0	0.23	ug/L			04/16/15 16:53	1
Bromomethane	1.0	U	1.0	0.31	ug/L			04/16/15 16:53	1
Chloroethane	1.0	U	1.0	0.21	ug/L			04/16/15 16:53	1
1,1-Dichloroethene	1.0	U	1.0	0.30	ug/L			04/16/15 16:53	1
Acetone	5.0	U	5.0	2.5	ug/L			04/16/15 16:53	1
Carbon disulfide	1.0	U	1.0	0.21	ug/L			04/16/15 16:53	1
<b>Methylene Chloride</b>	<b>0.14</b>	<b>J B</b>	1.0	0.13	ug/L			04/16/15 16:53	1
trans-1,2-Dichloroethene	1.0	U	1.0	0.17	ug/L			04/16/15 16:53	1
Methyl tert-butyl ether	1.0	U	1.0	0.18	ug/L			04/16/15 16:53	1
1,1-Dichloroethane	1.0	U	1.0	0.12	ug/L			04/16/15 16:53	1
cis-1,2-Dichloroethene	1.0	U	1.0	0.24	ug/L			04/16/15 16:53	1
Bromochloromethane	1.0	U	1.0	0.18	ug/L			04/16/15 16:53	1
2-Butanone (MEK)	5.0	U	5.0	0.55	ug/L			04/16/15 16:53	1
Chloroform	1.0	U	1.0	0.17	ug/L			04/16/15 16:53	1
1,1,1-Trichloroethane	1.0	U	1.0	0.29	ug/L			04/16/15 16:53	1
Carbon tetrachloride	1.0	U	1.0	0.14	ug/L			04/16/15 16:53	1
Benzene	1.0	U	1.0	0.11	ug/L			04/16/15 16:53	1
1,2-Dichloroethane	1.0	U	1.0	0.21	ug/L			04/16/15 16:53	1
Trichloroethene	1.0	U	1.0	0.14	ug/L			04/16/15 16:53	1
1,2-Dichloropropane	1.0	U	1.0	0.095	ug/L			04/16/15 16:53	1
Bromodichloromethane	1.0	U	1.0	0.13	ug/L			04/16/15 16:53	1
cis-1,3-Dichloropropene	1.0	U	1.0	0.19	ug/L			04/16/15 16:53	1
4-Methyl-2-pentanone (MIBK)	5.0	U	5.0	0.53	ug/L			04/16/15 16:53	1
Toluene	1.0	U	1.0	0.15	ug/L			04/16/15 16:53	1
trans-1,3-Dichloropropene	1.0	U	1.0	0.15	ug/L			04/16/15 16:53	1
1,1,2-Trichloroethane	1.0	U	1.0	0.20	ug/L			04/16/15 16:53	1
Tetrachloroethene	1.0	U	1.0	0.15	ug/L			04/16/15 16:53	1
2-Hexanone	5.0	U	5.0	0.16	ug/L			04/16/15 16:53	1
Dibromochloromethane	1.0	U	1.0	0.14	ug/L			04/16/15 16:53	1
1,2-Dibromoethane (EDB)	1.0	U	1.0	0.18	ug/L			04/16/15 16:53	1
Chlorobenzene	1.0	U	1.0	0.14	ug/L			04/16/15 16:53	1
1,1,1,2-Tetrachloroethane	1.0	U	1.0	0.28	ug/L			04/16/15 16:53	1
Ethylbenzene	1.0	U	1.0	0.23	ug/L			04/16/15 16:53	1
Xylenes, Total	3.0	U	3.0	0.49	ug/L			04/16/15 16:53	1
Styrene	1.0	U	1.0	0.097	ug/L			04/16/15 16:53	1
Bromoform	1.0	U	1.0	0.19	ug/L			04/16/15 16:53	1
1,1,2,2-Tetrachloroethane	1.0	U	1.0	0.20	ug/L			04/16/15 16:53	1
Acrylonitrile	20	U	20	0.55	ug/L			04/16/15 16:53	1
1,4-Dioxane	200	U	200	34	ug/L			04/16/15 16:53	1
<b>Surrogate</b>	<b>%Recovery</b>	<b>Qualifier</b>	<b>Limits</b>			<b>Prepared</b>	<b>Analyzed</b>	<b>Dil Fac</b>	
1,2-Dichloroethane-d4 (Surr)	115		64 - 135				04/16/15 16:53		1
Toluene-d8 (Surr)	98		71 - 118				04/16/15 16:53		1
4-Bromofluorobenzene (Surr)	92		70 - 118				04/16/15 16:53		1
Dibromofluoromethane (Surr)	109		70 - 128				04/16/15 16:53		1

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# Client Sample Results

Client: Groundwater Sciences Corporation  
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## Method: 8260C - Volatile Organic Compounds (GC/MS)

**Client Sample ID: HD-QC2-0/1-2**

**Date Collected: 04/10/15 12:00**

**Date Received: 04/11/15 09:30**

**Lab Sample ID: 180-42975-4**

**Matrix: Water**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Chloromethane	1.0	U	1.0	0.28	ug/L		04/15/15 15:33		1
Vinyl chloride	1.0	U	1.0	0.23	ug/L		04/15/15 15:33		1
Bromomethane	1.0	U	1.0	0.31	ug/L		04/15/15 15:33		1
Chloroethane	1.0	U	1.0	0.21	ug/L		04/15/15 15:33		1
1,1-Dichloroethene	1.0	U	1.0	0.30	ug/L		04/15/15 15:33		1
Acetone	5.0	U	5.0	2.5	ug/L		04/15/15 15:33		1
Carbon disulfide	1.0	U *	1.0	0.21	ug/L		04/15/15 15:33		1
<b>Methylene Chloride</b>	<b>0.22</b>	<b>J</b>	1.0	0.13	ug/L		04/15/15 15:33		1
trans-1,2-Dichloroethene	1.0	U	1.0	0.17	ug/L		04/15/15 15:33		1
Methyl tert-butyl ether	1.0	U	1.0	0.18	ug/L		04/15/15 15:33		1
1,1-Dichloroethane	1.0	U	1.0	0.12	ug/L		04/15/15 15:33		1
cis-1,2-Dichloroethene	1.0	U	1.0	0.24	ug/L		04/15/15 15:33		1
Bromochloromethane	1.0	U	1.0	0.18	ug/L		04/15/15 15:33		1
2-Butanone (MEK)	5.0	U	5.0	0.55	ug/L		04/15/15 15:33		1
Chloroform	1.0	U	1.0	0.17	ug/L		04/15/15 15:33		1
1,1,1-Trichloroethane	1.0	U	1.0	0.29	ug/L		04/15/15 15:33		1
Carbon tetrachloride	1.0	U	1.0	0.14	ug/L		04/15/15 15:33		1
Benzene	1.0	U	1.0	0.11	ug/L		04/15/15 15:33		1
1,2-Dichloroethane	1.0	U	1.0	0.21	ug/L		04/15/15 15:33		1
Trichloroethene	1.0	U	1.0	0.14	ug/L		04/15/15 15:33		1
1,2-Dichloropropane	1.0	U	1.0	0.095	ug/L		04/15/15 15:33		1
Bromodichloromethane	1.0	U	1.0	0.13	ug/L		04/15/15 15:33		1
cis-1,3-Dichloropropene	1.0	U	1.0	0.19	ug/L		04/15/15 15:33		1
4-Methyl-2-pentanone (MIBK)	5.0	U	5.0	0.53	ug/L		04/15/15 15:33		1
Toluene	1.0	U	1.0	0.15	ug/L		04/15/15 15:33		1
trans-1,3-Dichloropropene	1.0	U	1.0	0.15	ug/L		04/15/15 15:33		1
1,1,2-Trichloroethane	1.0	U	1.0	0.20	ug/L		04/15/15 15:33		1
Tetrachloroethene	1.0	U	1.0	0.15	ug/L		04/15/15 15:33		1
2-Hexanone	5.0	U	5.0	0.16	ug/L		04/15/15 15:33		1
Dibromochloromethane	1.0	U	1.0	0.14	ug/L		04/15/15 15:33		1
1,2-Dibromoethane (EDB)	1.0	U	1.0	0.18	ug/L		04/15/15 15:33		1
Chlorobenzene	1.0	U	1.0	0.14	ug/L		04/15/15 15:33		1
1,1,1,2-Tetrachloroethane	1.0	U	1.0	0.28	ug/L		04/15/15 15:33		1
Ethylbenzene	1.0	U	1.0	0.23	ug/L		04/15/15 15:33		1
Xylenes, Total	3.0	U	3.0	0.49	ug/L		04/15/15 15:33		1
Styrene	1.0	U	1.0	0.097	ug/L		04/15/15 15:33		1
Bromoform	1.0	U	1.0	0.19	ug/L		04/15/15 15:33		1
1,1,2,2-Tetrachloroethane	1.0	U	1.0	0.20	ug/L		04/15/15 15:33		1
Acrylonitrile	20	U	20	0.55	ug/L		04/15/15 15:33		1
1,4-Dioxane	200	U	200	34	ug/L		04/15/15 15:33		1
<b>Surrogate</b>	<b>%Recovery</b>	<b>Qualifier</b>	<b>Limits</b>			<b>Prepared</b>	<b>Analyzed</b>	<b>Dil Fac</b>	
1,2-Dichloroethane-d4 (Surr)	114		64 - 135				04/15/15 15:33		1
Toluene-d8 (Surr)	103		71 - 118				04/15/15 15:33		1
4-Bromofluorobenzene (Surr)	97		70 - 118				04/15/15 15:33		1
Dibromofluoromethane (Surr)	112		70 - 128				04/15/15 15:33		1

TestAmerica Pittsburgh

# Client Sample Results

Client: Groundwater Sciences Corporation  
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## Method: 8260C - Volatile Organic Compounds (GC/MS) - DL

**Client Sample ID: HD-MW-162-01-0**

**Date Collected: 04/10/15 09:50**

**Date Received: 04/11/15 09:30**

**Lab Sample ID: 180-42975-2**

**Matrix: Water**

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Chloromethane	40	U	40	11	ug/L			04/16/15 16:29	40
Vinyl chloride	40	U	40	9.1	ug/L			04/16/15 16:29	40
Bromomethane	40	U	40	13	ug/L			04/16/15 16:29	40
Chloroethane	40	U	40	8.6	ug/L			04/16/15 16:29	40
1,1-Dichloroethene	40	U	40	12	ug/L			04/16/15 16:29	40
Acetone	200	U	200	100	ug/L			04/16/15 16:29	40
Carbon disulfide	40	U	40	8.5	ug/L			04/16/15 16:29	40
<b>Methylene Chloride</b>	<b>22 J B</b>		40	5.0	ug/L			04/16/15 16:29	40
trans-1,2-Dichloroethene	40	U	40	6.8	ug/L			04/16/15 16:29	40
Methyl tert-butyl ether	40	U	40	7.3	ug/L			04/16/15 16:29	40
1,1-Dichloroethane	40	U	40	4.7	ug/L			04/16/15 16:29	40
cis-1,2-Dichloroethene	40	U	40	9.5	ug/L			04/16/15 16:29	40
Bromochloromethane	40	U	40	7.2	ug/L			04/16/15 16:29	40
2-Butanone (MEK)	200	U	200	22	ug/L			04/16/15 16:29	40
Chloroform	40	U	40	6.8	ug/L			04/16/15 16:29	40
1,1,1-Trichloroethane	40	U	40	11	ug/L			04/16/15 16:29	40
Carbon tetrachloride	40	U	40	5.5	ug/L			04/16/15 16:29	40
Benzene	40	U	40	4.2	ug/L			04/16/15 16:29	40
1,2-Dichloroethane	40	U	40	8.5	ug/L			04/16/15 16:29	40
<b>Trichloroethene</b>	<b>180</b>		40	5.7	ug/L			04/16/15 16:29	40
1,2-Dichloropropane	40	U	40	3.8	ug/L			04/16/15 16:29	40
Bromodichloromethane	40	U	40	5.2	ug/L			04/16/15 16:29	40
cis-1,3-Dichloropropene	40	U	40	7.5	ug/L			04/16/15 16:29	40
4-Methyl-2-pentanone (MIBK)	200	U	200	21	ug/L			04/16/15 16:29	40
Toluene	40	U	40	6.0	ug/L			04/16/15 16:29	40
trans-1,3-Dichloropropene	40	U	40	5.9	ug/L			04/16/15 16:29	40
1,1,2-Trichloroethane	40	U	40	8.1	ug/L			04/16/15 16:29	40
<b>Tetrachloroethene</b>	<b>700</b>		40	5.9	ug/L			04/16/15 16:29	40
2-Hexanone	200	U	200	6.4	ug/L			04/16/15 16:29	40
Dibromochloromethane	40	U	40	5.5	ug/L			04/16/15 16:29	40
1,2-Dibromoethane (EDB)	40	U	40	7.2	ug/L			04/16/15 16:29	40
Chlorobenzene	40	U	40	5.4	ug/L			04/16/15 16:29	40
1,1,1,2-Tetrachloroethane	40	U	40	11	ug/L			04/16/15 16:29	40
Ethylbenzene	40	U	40	9.1	ug/L			04/16/15 16:29	40
Xylenes, Total	120	U	120	20	ug/L			04/16/15 16:29	40
Styrene	40	U	40	3.9	ug/L			04/16/15 16:29	40
Bromoform	40	U	40	7.7	ug/L			04/16/15 16:29	40
1,1,2,2-Tetrachloroethane	40	U	40	8.0	ug/L			04/16/15 16:29	40
Acrylonitrile	800	U	800	22	ug/L			04/16/15 16:29	40
1,4-Dioxane	8000	U	8000	1400	ug/L			04/16/15 16:29	40
<b>Surrogate</b>	<b>%Recovery</b>	<b>Qualifier</b>		<b>Limits</b>		<b>Prepared</b>	<b>Analyzed</b>	<b>Dil Fac</b>	
1,2-Dichloroethane-d4 (Surr)	107			64 - 135					40
Toluene-d8 (Surr)	102			71 - 118					40
4-Bromofluorobenzene (Surr)	98			70 - 118					40
Dibromofluoromethane (Surr)	110			70 - 128					40

TestAmerica Pittsburgh

# Default Detection Limits

Client: Groundwater Sciences Corporation  
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## **Method: 8260C - Volatile Organic Compounds (GC/MS)**

Analyte	RL	MDL	Units	Method
1,1,1,2-Tetrachloroethane	1.0	0.28	ug/L	8260C
1,1,1-Trichloroethane	1.0	0.29	ug/L	8260C
1,1,2,2-Tetrachloroethane	1.0	0.20	ug/L	8260C
1,1,2-Trichloroethane	1.0	0.20	ug/L	8260C
1,1-Dichloroethane	1.0	0.12	ug/L	8260C
1,1-Dichloroethene	1.0	0.30	ug/L	8260C
1,2-Dibromoethane (EDB)	1.0	0.18	ug/L	8260C
1,2-Dichloroethane	1.0	0.21	ug/L	8260C
1,2-Dichloropropane	1.0	0.095	ug/L	8260C
1,4-Dioxane	200	34	ug/L	8260C
2-Butanone (MEK)	5.0	0.55	ug/L	8260C
2-Hexanone	5.0	0.16	ug/L	8260C
4-Methyl-2-pentanone (MIBK)	5.0	0.53	ug/L	8260C
Acetone	5.0	2.5	ug/L	8260C
Acrylonitrile	20	0.55	ug/L	8260C
Benzene	1.0	0.11	ug/L	8260C
Bromochloromethane	1.0	0.18	ug/L	8260C
Bromodichloromethane	1.0	0.13	ug/L	8260C
Bromoform	1.0	0.19	ug/L	8260C
Bromomethane	1.0	0.31	ug/L	8260C
Carbon disulfide	1.0	0.21	ug/L	8260C
Carbon tetrachloride	1.0	0.14	ug/L	8260C
Chlorobenzene	1.0	0.14	ug/L	8260C
Chloroethane	1.0	0.21	ug/L	8260C
Chloroform	1.0	0.17	ug/L	8260C
Chloromethane	1.0	0.28	ug/L	8260C
cis-1,2-Dichloroethene	1.0	0.24	ug/L	8260C
cis-1,3-Dichloropropene	1.0	0.19	ug/L	8260C
Dibromochloromethane	1.0	0.14	ug/L	8260C
Ethylbenzene	1.0	0.23	ug/L	8260C
Methyl tert-butyl ether	1.0	0.18	ug/L	8260C
Methylene Chloride	1.0	0.13	ug/L	8260C
Styrene	1.0	0.097	ug/L	8260C
Tetrachloroethene	1.0	0.15	ug/L	8260C
Toluene	1.0	0.15	ug/L	8260C
trans-1,2-Dichloroethene	1.0	0.17	ug/L	8260C
trans-1,3-Dichloropropene	1.0	0.15	ug/L	8260C
Trichloroethene	1.0	0.14	ug/L	8260C
Vinyl chloride	1.0	0.23	ug/L	8260C
Xylenes, Total	3.0	0.49	ug/L	8260C

# Surrogate Summary

Client: Groundwater Sciences Corporation  
Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## Method: 8260C - Volatile Organic Compounds (GC/MS)

Matrix: Water

Prep Type: Total/NA

Lab Sample ID	Client Sample ID	Percent Surrogate Recovery (Acceptance Limits)			
		12DCE (64-135)	TOL (71-118)	BFB (70-118)	DBFM (70-128)
180-42975-1	HD-MW-165-0/1-0	118	100	96	110
180-42975-1 MS	HD-MW-165-0/1-0	90	99	89	92
180-42975-1 MSD	HD-MW-165-0/1-0	96	96	93	92
180-42975-2	HD-MW-162-0/1-0	115	98	91	116
180-42975-2 - DL	HD-MW-162-0/1-0	107	102	98	110
180-42975-3	HD-MW-169-0/1-0	115	98	92	109
180-42975-4	HD-QC2-0/1-2	114	103	97	112
LCS 180-138583/8	Lab Control Sample	95	99	94	91
LCS 180-138685/8	Lab Control Sample	93	101	94	93
LCSD 180-138685/10	Lab Control Sample Dup	92	97	92	89
MB 180-138583/5	Method Blank	113	101	93	107
MB 180-138685/5	Method Blank	116	100	97	111

### Surrogate Legend

12DCE = 1,2-Dichloroethane-d4 (Surr)

TOL = Toluene-d8 (Surr)

BFB = 4-Bromofluorobenzene (Surr)

DBFM = Dibromofluoromethane (Surr)

# QC Sample Results

Client: Groundwater Sciences Corporation  
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## Method: 8260C - Volatile Organic Compounds (GC/MS)

Lab Sample ID: MB 180-138583/5

Matrix: Water

Analysis Batch: 138583

Client Sample ID: Method Blank

Prep Type: Total/NA

Analyte	MB	MB	D	Prepared	Analyzed	Dil Fac
	Result	Qualifier				
Chloromethane	1.0	U	1.0	0.28	ug/L	1
Vinyl chloride	1.0	U	1.0	0.23	ug/L	1
Bromomethane	1.0	U	1.0	0.31	ug/L	1
Chloroethane	1.0	U	1.0	0.21	ug/L	1
1,1-Dichloroethene	1.0	U	1.0	0.30	ug/L	1
Acetone	5.0	U	5.0	2.5	ug/L	1
Carbon disulfide	1.0	U	1.0	0.21	ug/L	1
Methylene Chloride	1.0	U	1.0	0.13	ug/L	1
trans-1,2-Dichloroethene	1.0	U	1.0	0.17	ug/L	1
Methyl tert-butyl ether	1.0	U	1.0	0.18	ug/L	1
1,1-Dichloroethane	1.0	U	1.0	0.12	ug/L	1
cis-1,2-Dichloroethene	1.0	U	1.0	0.24	ug/L	1
Bromochloromethane	1.0	U	1.0	0.18	ug/L	1
2-Butanone (MEK)	5.0	U	5.0	0.55	ug/L	1
Chloroform	1.0	U	1.0	0.17	ug/L	1
1,1,1-Trichloroethane	1.0	U	1.0	0.29	ug/L	1
Carbon tetrachloride	1.0	U	1.0	0.14	ug/L	1
Benzene	1.0	U	1.0	0.11	ug/L	1
1,2-Dichloroethane	1.0	U	1.0	0.21	ug/L	1
Trichloroethene	1.0	U	1.0	0.14	ug/L	1
1,2-Dichloropropane	1.0	U	1.0	0.095	ug/L	1
Bromodichloromethane	1.0	U	1.0	0.13	ug/L	1
cis-1,3-Dichloropropene	1.0	U	1.0	0.19	ug/L	1
4-Methyl-2-pentanone (MIBK)	5.0	U	5.0	0.53	ug/L	1
Toluene	1.0	U	1.0	0.15	ug/L	1
trans-1,3-Dichloropropene	1.0	U	1.0	0.15	ug/L	1
1,1,2-Trichloroethane	1.0	U	1.0	0.20	ug/L	1
Tetrachloroethene	1.0	U	1.0	0.15	ug/L	1
2-Hexanone	5.0	U	5.0	0.16	ug/L	1
Dibromochloromethane	1.0	U	1.0	0.14	ug/L	1
1,2-Dibromoethane (EDB)	1.0	U	1.0	0.18	ug/L	1
Chlorobenzene	1.0	U	1.0	0.14	ug/L	1
1,1,1,2-Tetrachloroethane	1.0	U	1.0	0.28	ug/L	1
Ethylbenzene	1.0	U	1.0	0.23	ug/L	1
Xylenes, Total	3.0	U	3.0	0.49	ug/L	1
Styrene	1.0	U	1.0	0.097	ug/L	1
Bromoform	1.0	U	1.0	0.19	ug/L	1
1,1,2,2-Tetrachloroethane	1.0	U	1.0	0.20	ug/L	1
Acrylonitrile	20	U	20	0.55	ug/L	1
1,4-Dioxane	200	U	200	34	ug/L	1

Surrogate	MB	MB	Limits	Prepared	Analyzed	Dil Fac
	%Recovery	Qualifier				
1,2-Dichloroethane-d4 (Surr)	113		64 - 135		04/15/15 14:33	1
Toluene-d8 (Surr)	101		71 - 118		04/15/15 14:33	1
4-Bromofluorobenzene (Surr)	93		70 - 118		04/15/15 14:33	1
Dibromofluoromethane (Surr)	107		70 - 128		04/15/15 14:33	1

TestAmerica Pittsburgh

# QC Sample Results

Client: Groundwater Sciences Corporation  
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## Method: 8260C - Volatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: LCS 180-138583/8**

**Matrix: Water**

**Analysis Batch: 138583**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**

Analyte	Spike	LCS	LCS	Unit	D	%Rec	%Rec.
	Added	Result	Qualifier				Limits
Chloromethane	10.0	10.1		ug/L		101	50 - 139
Vinyl chloride	10.0	10.0		ug/L		100	53 - 138
Bromomethane	10.0	8.98		ug/L		90	33 - 150
Chloroethane	10.0	10.5		ug/L		105	36 - 142
1,1-Dichloroethene	10.0	7.04		ug/L		70	65 - 136
Acetone	20.0	19.0		ug/L		95	22 - 150
Carbon disulfide	10.0	3.87	*	ug/L		39	54 - 132
Methylene Chloride	10.0	7.62		ug/L		76	63 - 129
trans-1,2-Dichloroethene	10.0	7.74		ug/L		77	73 - 126
Methyl tert-butyl ether	10.0	7.94		ug/L		79	64 - 123
1,1-Dichloroethane	10.0	8.31		ug/L		83	73 - 126
cis-1,2-Dichloroethene	10.0	7.88		ug/L		79	70 - 120
Bromochloromethane	10.0	8.32		ug/L		83	70 - 127
2-Butanone (MEK)	20.0	16.7		ug/L		83	39 - 138
Chloroform	10.0	9.11		ug/L		91	72 - 127
1,1,1-Trichloroethane	10.0	8.87		ug/L		89	63 - 133
Carbon tetrachloride	10.0	9.33		ug/L		93	55 - 150
Benzene	10.0	8.86		ug/L		89	80 - 120
1,2-Dichloroethane	10.0	8.84		ug/L		88	68 - 132
Trichloroethene	10.0	8.40		ug/L		84	73 - 120
1,2-Dichloropropane	10.0	9.27		ug/L		93	76 - 124
Bromodichloromethane	10.0	8.85		ug/L		88	66 - 130
cis-1,3-Dichloropropene	10.0	9.60		ug/L		96	66 - 120
4-Methyl-2-pentanone (MIBK)	20.0	16.8		ug/L		84	45 - 145
Toluene	10.0	9.81		ug/L		98	80 - 123
trans-1,3-Dichloropropene	10.0	11.6		ug/L		116	65 - 125
1,1,2-Trichloroethane	10.0	11.2		ug/L		112	77 - 127
Tetrachloroethene	10.0	9.33		ug/L		93	70 - 135
2-Hexanone	20.0	15.3		ug/L		76	25 - 132
Dibromochloromethane	10.0	10.1		ug/L		101	60 - 140
1,2-Dibromoethane (EDB)	10.0	9.99		ug/L		100	74 - 123
Chlorobenzene	10.0	10.2		ug/L		102	80 - 120
1,1,1,2-Tetrachloroethane	10.0	11.8		ug/L		118	63 - 140
Ethylbenzene	10.0	9.72		ug/L		97	72 - 126
Xylenes, Total	20.0	19.0		ug/L		95	76 - 128
Styrene	10.0	9.70		ug/L		97	71 - 127
Bromoform	10.0	9.35		ug/L		93	46 - 150
1,1,2,2-Tetrachloroethane	10.0	10.9		ug/L		109	62 - 125
1,4-Dioxane	200	179	J	ug/L		90	10 - 160
<b>Surrogate</b>		<b>LCS</b>	<b>LCS</b>				
<b>Surrogate</b>		<b>%Recovery</b>	<b>Qualifier</b>	<b>Limits</b>			
1,2-Dichloroethane-d4 (Surr)		95		64 - 135			
Toluene-d8 (Surr)		99		71 - 118			
4-Bromofluorobenzene (Surr)		94		70 - 118			
Dibromofluoromethane (Surr)		91		70 - 128			

TestAmerica Pittsburgh

# QC Sample Results

Client: Groundwater Sciences Corporation  
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## Method: 8260C - Volatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: 180-42975-1 MS**

**Client Sample ID: HD-MW-165-0/1-0**

**Matrix: Water**

**Prep Type: Total/NA**

**Analysis Batch: 138583**

Analyte	Sample	Sample	Spike	MS	MS	Unit	D	%Rec	%Rec.
	Result	Qualifier	Added	Result	Qualifier				
Chloromethane	1.0	U	10.0	10.5		ug/L		105	50 - 139
Vinyl chloride	1.0	U	10.0	9.56		ug/L		96	53 - 138
Bromomethane	1.0	U	10.0	9.20		ug/L		92	33 - 150
Chloroethane	1.0	U	10.0	10.3		ug/L		103	36 - 142
1,1-Dichloroethene	1.0	U	10.0	6.75		ug/L		67	65 - 136
Acetone	5.0	U	20.0	20.2		ug/L		101	22 - 150
Carbon disulfide	1.0	U F1 *	10.0	3.81	F1	ug/L		38	54 - 132
Methylene Chloride	1.0	U	10.0	7.73		ug/L		77	63 - 129
trans-1,2-Dichloroethene	1.0	U	10.0	7.54		ug/L		75	73 - 126
Methyl tert-butyl ether	1.0	U	10.0	8.13		ug/L		81	64 - 123
1,1-Dichloroethane	1.0	U	10.0	8.29		ug/L		83	73 - 126
cis-1,2-Dichloroethene	1.0	U	10.0	8.26		ug/L		83	70 - 120
Bromochloromethane	1.0	U	10.0	7.70		ug/L		77	70 - 127
2-Butanone (MEK)	5.0	U	20.0	16.6		ug/L		83	39 - 138
Chloroform	1.0	U	10.0	9.04		ug/L		90	72 - 127
1,1,1-Trichloroethane	1.0	U	10.0	8.69		ug/L		87	63 - 133
Carbon tetrachloride	1.0	U	10.0	8.96		ug/L		90	55 - 150
Benzene	1.0	U	10.0	8.62		ug/L		86	80 - 120
1,2-Dichloroethane	1.0	U	10.0	8.54		ug/L		85	68 - 132
Trichloroethene	16	F1	10.0	19.5	F1	ug/L		37	73 - 120
1,2-Dichloropropane	1.0	U	10.0	8.78		ug/L		88	76 - 124
Bromodichloromethane	1.0	U	10.0	8.29		ug/L		83	66 - 130
cis-1,3-Dichloropropene	1.0	U	10.0	8.80		ug/L		88	66 - 120
4-Methyl-2-pentanone (MIBK)	5.0	U	20.0	16.8		ug/L		84	45 - 145
Toluene	1.0	U	10.0	9.59		ug/L		96	80 - 123
trans-1,3-Dichloropropene	1.0	U	10.0	11.0		ug/L		110	65 - 125
1,1,2-Trichloroethane	1.0	U	10.0	10.7		ug/L		107	77 - 127
Tetrachloroethene	7.4	F1	10.0	15.2		ug/L		78	70 - 135
2-Hexanone	5.0	U	20.0	15.4		ug/L		77	25 - 132
Dibromochloromethane	1.0	U	10.0	9.09		ug/L		91	60 - 140
1,2-Dibromoethane (EDB)	1.0	U	10.0	9.76		ug/L		98	74 - 123
Chlorobenzene	1.0	U	10.0	9.83		ug/L		98	80 - 120
1,1,1,2-Tetrachloroethane	1.0	U	10.0	10.7		ug/L		107	63 - 140
Ethylbenzene	1.0	U	10.0	9.18		ug/L		92	72 - 126
Xylenes, Total	3.0	U	20.0	18.5		ug/L		93	76 - 128
Styrene	1.0	U	10.0	9.29		ug/L		93	71 - 127
Bromoform	1.0	U	10.0	8.75		ug/L		87	46 - 150
1,1,2,2-Tetrachloroethane	1.0	U	10.0	10.1		ug/L		101	62 - 125
1,4-Dioxane	200	U	200	175	J	ug/L		87	10 - 160
<hr/>									
Surrogate	MS		MS		Limits				
	%Recovery	Qualifier							
1,2-Dichloroethane-d4 (Surr)	90				64 - 135				
Toluene-d8 (Surr)	99				71 - 118				
4-Bromofluorobenzene (Surr)	89				70 - 118				
Dibromofluoromethane (Surr)	92				70 - 128				

TestAmerica Pittsburgh

# QC Sample Results

Client: Groundwater Sciences Corporation  
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## Method: 8260C - Volatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: 180-42975-1 MSD**

**Matrix: Water**

**Analysis Batch: 138583**

**Client Sample ID: HD-MW-165-0/1-0**

**Prep Type: Total/NA**

Analyte	Sample	Sample	Spike	MSD	MSD	Unit	D	%Rec	%Rec.	RPD	RPD Limit
	Result	Qualifier	Added	Result	Qualifier				Limits		
Chloromethane	1.0	U	10.0	10.8		ug/L		108	50 - 139	3	35
Vinyl chloride	1.0	U	10.0	9.37		ug/L		94	53 - 138	2	35
Bromomethane	1.0	U	10.0	10.1		ug/L		101	33 - 150	9	35
Chloroethane	1.0	U	10.0	10.4		ug/L		104	36 - 142	0	35
1,1-Dichloroethene	1.0	U	10.0	6.67		ug/L		67	65 - 136	1	35
Acetone	5.0	U	20.0	19.1		ug/L		96	22 - 150	5	35
Carbon disulfide	1.0	U F1 *	10.0	3.88	F1	ug/L		39	54 - 132	2	35
Methylene Chloride	1.0	U	10.0	8.19		ug/L		82	63 - 129	6	35
trans-1,2-Dichloroethene	1.0	U	10.0	7.67		ug/L		77	73 - 126	2	35
Methyl tert-butyl ether	1.0	U	10.0	8.73		ug/L		87	64 - 123	7	35
1,1-Dichloroethane	1.0	U	10.0	8.39		ug/L		84	73 - 126	1	35
cis-1,2-Dichloroethene	1.0	U	10.0	8.61		ug/L		86	70 - 120	4	35
Bromochloromethane	1.0	U	10.0	9.01		ug/L		90	70 - 127	16	35
2-Butanone (MEK)	5.0	U	20.0	18.7		ug/L		94	39 - 138	12	35
Chloroform	1.0	U	10.0	9.15		ug/L		92	72 - 127	1	35
1,1,1-Trichloroethane	1.0	U	10.0	8.67		ug/L		87	63 - 133	0	35
Carbon tetrachloride	1.0	U	10.0	8.65		ug/L		86	55 - 150	4	35
Benzene	1.0	U	10.0	8.79		ug/L		88	80 - 120	2	32
1,2-Dichloroethane	1.0	U	10.0	9.38		ug/L		94	68 - 132	9	32
Trichloroethene	16	F1	10.0	19.7	F1	ug/L		38	73 - 120	1	35
1,2-Dichloropropane	1.0	U	10.0	9.80		ug/L		98	76 - 124	11	34
Bromodichloromethane	1.0	U	10.0	9.01		ug/L		90	66 - 130	8	35
cis-1,3-Dichloropropene	1.0	U	10.0	9.60		ug/L		96	66 - 120	9	35
4-Methyl-2-pentanone (MIBK)	5.0	U	20.0	18.0		ug/L		90	45 - 145	7	35
Toluene	1.0	U	10.0	9.38		ug/L		94	80 - 123	2	35
trans-1,3-Dichloropropene	1.0	U	10.0	11.3		ug/L		113	65 - 125	3	35
1,1,2-Trichloroethane	1.0	U	10.0	10.4		ug/L		104	77 - 127	3	35
Tetrachloroethene	7.4	F1	10.0	14.2	F1	ug/L		68	70 - 135	7	35
2-Hexanone	5.0	U	20.0	16.2		ug/L		81	25 - 132	5	35
Dibromochloromethane	1.0	U	10.0	9.15		ug/L		91	60 - 140	1	35
1,2-Dibromoethane (EDB)	1.0	U	10.0	10.2		ug/L		102	74 - 123	4	35
Chlorobenzene	1.0	U	10.0	9.90		ug/L		99	80 - 120	1	29
1,1,1,2-Tetrachloroethane	1.0	U	10.0	10.8		ug/L		108	63 - 140	0	34
Ethylbenzene	1.0	U	10.0	9.33		ug/L		93	72 - 126	2	33
Xylenes, Total	3.0	U	20.0	18.3		ug/L		91	76 - 128	1	32
Styrene	1.0	U	10.0	9.54		ug/L		95	71 - 127	3	34
Bromoform	1.0	U	10.0	8.52		ug/L		85	46 - 150	3	35
1,1,2,2-Tetrachloroethane	1.0	U	10.0	10.5		ug/L		105	62 - 125	4	35
1,4-Dioxane	200	U	200	191	J	ug/L		95	10 - 160	9	35
<b>MSD MSD</b>											
Surrogate	%Recovery	Qualifier		<b>Limits</b>							
1,2-Dichloroethane-d4 (Surr)	96			64 - 135							
Toluene-d8 (Surr)	96			71 - 118							
4-Bromofluorobenzene (Surr)	93			70 - 118							
Dibromofluoromethane (Surr)	92			70 - 128							

TestAmerica Pittsburgh

# QC Sample Results

Client: Groundwater Sciences Corporation  
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## Method: 8260C - Volatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: MB 180-138685/5**

**Matrix: Water**

**Analysis Batch: 138685**

**Client Sample ID: Method Blank**

**Prep Type: Total/NA**

Analyte	MB	MB	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
	Result	Qualifier									
Chloromethane	1.0	U			1.0	0.28	ug/L			04/16/15 11:32	1
Vinyl chloride	1.0	U			1.0	0.23	ug/L			04/16/15 11:32	1
Bromomethane	1.0	U			1.0	0.31	ug/L			04/16/15 11:32	1
Chloroethane	1.0	U			1.0	0.21	ug/L			04/16/15 11:32	1
1,1-Dichloroethene	1.0	U			1.0	0.30	ug/L			04/16/15 11:32	1
Acetone	5.0	U			5.0	2.5	ug/L			04/16/15 11:32	1
Carbon disulfide	1.0	U			1.0	0.21	ug/L			04/16/15 11:32	1
Methylene Chloride	0.162	J			1.0	0.13	ug/L			04/16/15 11:32	1
trans-1,2-Dichloroethene	1.0	U			1.0	0.17	ug/L			04/16/15 11:32	1
Methyl tert-butyl ether	1.0	U			1.0	0.18	ug/L			04/16/15 11:32	1
1,1-Dichloroethane	1.0	U			1.0	0.12	ug/L			04/16/15 11:32	1
cis-1,2-Dichloroethene	1.0	U			1.0	0.24	ug/L			04/16/15 11:32	1
Bromochloromethane	1.0	U			1.0	0.18	ug/L			04/16/15 11:32	1
2-Butanone (MEK)	5.0	U			5.0	0.55	ug/L			04/16/15 11:32	1
Chloroform	1.0	U			1.0	0.17	ug/L			04/16/15 11:32	1
1,1,1-Trichloroethane	1.0	U			1.0	0.29	ug/L			04/16/15 11:32	1
Carbon tetrachloride	1.0	U			1.0	0.14	ug/L			04/16/15 11:32	1
Benzene	1.0	U			1.0	0.11	ug/L			04/16/15 11:32	1
1,2-Dichloroethane	1.0	U			1.0	0.21	ug/L			04/16/15 11:32	1
Trichloroethene	1.0	U			1.0	0.14	ug/L			04/16/15 11:32	1
1,2-Dichloropropane	1.0	U			1.0	0.095	ug/L			04/16/15 11:32	1
Bromodichloromethane	1.0	U			1.0	0.13	ug/L			04/16/15 11:32	1
cis-1,3-Dichloropropene	1.0	U			1.0	0.19	ug/L			04/16/15 11:32	1
4-Methyl-2-pentanone (MIBK)	5.0	U			5.0	0.53	ug/L			04/16/15 11:32	1
Toluene	1.0	U			1.0	0.15	ug/L			04/16/15 11:32	1
trans-1,3-Dichloropropene	1.0	U			1.0	0.15	ug/L			04/16/15 11:32	1
1,1,2-Trichloroethane	1.0	U			1.0	0.20	ug/L			04/16/15 11:32	1
Tetrachloroethene	1.0	U			1.0	0.15	ug/L			04/16/15 11:32	1
2-Hexanone	5.0	U			5.0	0.16	ug/L			04/16/15 11:32	1
Dibromochloromethane	1.0	U			1.0	0.14	ug/L			04/16/15 11:32	1
1,2-Dibromoethane (EDB)	1.0	U			1.0	0.18	ug/L			04/16/15 11:32	1
Chlorobenzene	1.0	U			1.0	0.14	ug/L			04/16/15 11:32	1
1,1,1,2-Tetrachloroethane	1.0	U			1.0	0.28	ug/L			04/16/15 11:32	1
Ethylbenzene	1.0	U			1.0	0.23	ug/L			04/16/15 11:32	1
Xylenes, Total	3.0	U			3.0	0.49	ug/L			04/16/15 11:32	1
Styrene	1.0	U			1.0	0.097	ug/L			04/16/15 11:32	1
Bromoform	1.0	U			1.0	0.19	ug/L			04/16/15 11:32	1
1,1,2,2-Tetrachloroethane	1.0	U			1.0	0.20	ug/L			04/16/15 11:32	1
Acrylonitrile	20	U			20	0.55	ug/L			04/16/15 11:32	1
1,4-Dioxane	200	U			200	34	ug/L			04/16/15 11:32	1

Surrogate	MB	MB	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
	Result	Qualifier						
1,2-Dichloroethane-d4 (Surr)	116		64 - 135					1
Toluene-d8 (Surr)	100		71 - 118					1
4-Bromofluorobenzene (Surr)	97		70 - 118					1
Dibromofluoromethane (Surr)	111		70 - 128					1

# QC Sample Results

Client: Groundwater Sciences Corporation  
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## Method: 8260C - Volatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: LCS 180-138685/8**

**Matrix: Water**

**Analysis Batch: 138685**

**Client Sample ID: Lab Control Sample**

**Prep Type: Total/NA**

Analyte	Spike	LCS	LCS	Unit	D	%Rec	%Rec.
	Added	Result	Qualifier				Limits
Chloromethane	10.0	10.0		ug/L		100	50 - 139
Vinyl chloride	10.0	10.5		ug/L		105	53 - 138
Bromomethane	10.0	9.71		ug/L		97	33 - 150
Chloroethane	10.0	11.0		ug/L		110	36 - 142
1,1-Dichloroethene	10.0	9.09		ug/L		91	65 - 136
Acetone	20.0	21.0		ug/L		105	22 - 150
Carbon disulfide	10.0	6.75		ug/L		68	54 - 132
Methylene Chloride	10.0	9.92		ug/L		99	63 - 129
trans-1,2-Dichloroethene	10.0	9.42		ug/L		94	73 - 126
Methyl tert-butyl ether	10.0	9.22		ug/L		92	64 - 123
1,1-Dichloroethane	10.0	9.66		ug/L		97	73 - 126
cis-1,2-Dichloroethene	10.0	9.13		ug/L		91	70 - 120
Bromochloromethane	10.0	9.67		ug/L		97	70 - 127
2-Butanone (MEK)	20.0	17.2		ug/L		86	39 - 138
Chloroform	10.0	10.1		ug/L		101	72 - 127
1,1,1-Trichloroethane	10.0	10.6		ug/L		106	63 - 133
Carbon tetrachloride	10.0	11.0		ug/L		110	55 - 150
Benzene	10.0	10.2		ug/L		102	80 - 120
1,2-Dichloroethane	10.0	10.0		ug/L		100	68 - 132
Trichloroethene	10.0	9.34		ug/L		93	73 - 120
1,2-Dichloropropane	10.0	9.73		ug/L		97	76 - 124
Bromodichloromethane	10.0	9.17		ug/L		92	66 - 130
cis-1,3-Dichloropropene	10.0	9.80		ug/L		98	66 - 120
4-Methyl-2-pentanone (MIBK)	20.0	16.2		ug/L		81	45 - 145
Toluene	10.0	10.5		ug/L		105	80 - 123
trans-1,3-Dichloropropene	10.0	10.8		ug/L		108	65 - 125
1,1,2-Trichloroethane	10.0	10.7		ug/L		107	77 - 127
Tetrachloroethene	10.0	10.2		ug/L		102	70 - 135
2-Hexanone	20.0	16.8		ug/L		84	25 - 132
Dibromochloromethane	10.0	9.51		ug/L		95	60 - 140
1,2-Dibromoethane (EDB)	10.0	10.2		ug/L		102	74 - 123
Chlorobenzene	10.0	10.3		ug/L		103	80 - 120
1,1,1,2-Tetrachloroethane	10.0	11.1		ug/L		111	63 - 140
Ethylbenzene	10.0	9.73		ug/L		97	72 - 126
Xylenes, Total	20.0	19.8		ug/L		99	76 - 128
Styrene	10.0	9.71		ug/L		97	71 - 127
Bromoform	10.0	9.19		ug/L		92	46 - 150
1,1,2,2-Tetrachloroethane	10.0	10.0		ug/L		100	62 - 125
1,4-Dioxane	200	167 J		ug/L		84	10 - 160

Surrogate	LCS	LCS	Limits
	%Recovery	Qualifier	
1,2-Dichloroethane-d4 (Surr)	93		64 - 135
Toluene-d8 (Surr)	101		71 - 118
4-Bromofluorobenzene (Surr)	94		70 - 118
Dibromofluoromethane (Surr)	93		70 - 128

TestAmerica Pittsburgh

# QC Sample Results

Client: Groundwater Sciences Corporation  
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## Method: 8260C - Volatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: LCSD 180-138685/10**

**Client Sample ID: Lab Control Sample Dup**

**Prep Type: Total/NA**

**Matrix: Water**

**Analysis Batch: 138685**

Analyte	Spike	LCSD	LCSD	Unit	D	%Rec	%Rec.		RPD	RPD	RPD
	Added	Result	Qualifier				Limits	Limit			
Chloromethane	10.0	10.8		ug/L		108	50 - 139		8		35
Vinyl chloride	10.0	10.3		ug/L		103	53 - 138		1		35
Bromomethane	10.0	8.47		ug/L		85	33 - 150		14		35
Chloroethane	10.0	10.7		ug/L		107	36 - 142		3		35
1,1-Dichloroethene	10.0	9.04		ug/L		90	65 - 136		1		35
Acetone	20.0	20.2		ug/L		101	22 - 150		4		35
Carbon disulfide	10.0	6.12		ug/L		61	54 - 132		10		35
Methylene Chloride	10.0	10.2		ug/L		102	63 - 129		3		35
trans-1,2-Dichloroethene	10.0	9.27		ug/L		93	73 - 126		2		35
Methyl tert-butyl ether	10.0	9.56		ug/L		96	64 - 123		4		35
1,1-Dichloroethane	10.0	9.47		ug/L		95	73 - 126		2		35
cis-1,2-Dichloroethene	10.0	9.45		ug/L		95	70 - 120		3		35
Bromochloromethane	10.0	9.04		ug/L		90	70 - 127		7		35
2-Butanone (MEK)	20.0	17.7		ug/L		89	39 - 138		3		35
Chloroform	10.0	9.80		ug/L		98	72 - 127		3		35
1,1,1-Trichloroethane	10.0	10.2		ug/L		102	63 - 133		3		35
Carbon tetrachloride	10.0	10.3		ug/L		103	55 - 150		7		35
Benzene	10.0	10.1		ug/L		101	80 - 120		1		32
1,2-Dichloroethane	10.0	9.65		ug/L		96	68 - 132		4		32
Trichloroethene	10.0	9.24		ug/L		92	73 - 120		1		35
1,2-Dichloropropane	10.0	9.60		ug/L		96	76 - 124		1		34
Bromodichloromethane	10.0	8.86		ug/L		89	66 - 130		3		35
cis-1,3-Dichloropropene	10.0	9.88		ug/L		99	66 - 120		1		35
4-Methyl-2-pentanone (MIBK)	20.0	16.9		ug/L		84	45 - 145		4		35
Toluene	10.0	10.4		ug/L		104	80 - 123		2		35
trans-1,3-Dichloropropene	10.0	10.3		ug/L		103	65 - 125		5		35
1,1,2-Trichloroethane	10.0	10.4		ug/L		104	77 - 127		2		35
Tetrachloroethene	10.0	10.0		ug/L		100	70 - 135		2		35
2-Hexanone	20.0	17.3		ug/L		86	25 - 132		3		35
Dibromochloromethane	10.0	9.55		ug/L		95	60 - 140		0		35
1,2-Dibromoethane (EDB)	10.0	10.1		ug/L		101	74 - 123		1		35
Chlorobenzene	10.0	10.2		ug/L		102	80 - 120		1		29
1,1,1,2-Tetrachloroethane	10.0	10.7		ug/L		107	63 - 140		4		34
Ethylbenzene	10.0	9.71		ug/L		97	72 - 126		0		33
Xylenes, Total	20.0	19.2		ug/L		96	76 - 128		3		32
Styrene	10.0	9.79		ug/L		98	71 - 127		1		34
Bromoform	10.0	8.77		ug/L		88	46 - 150		5		35
1,1,2,2-Tetrachloroethane	10.0	9.87		ug/L		99	62 - 125		2		35
1,4-Dioxane	200	193 J		ug/L		96	10 - 160		14		35

Surrogate	LCSD	LCSD	Limits
	%Recovery	Qualifier	
1,2-Dichloroethane-d4 (Surr)	92		64 - 135
Toluene-d8 (Surr)	97		71 - 118
4-Bromofluorobenzene (Surr)	92		70 - 118
Dibromofluoromethane (Surr)	89		70 - 128

TestAmerica Pittsburgh

# QC Association Summary

Client: Groundwater Sciences Corporation  
Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

## GC/MS VOA

### Analysis Batch: 138583

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
180-42975-1	HD-MW-165-0/1-0	Total/NA	Water	8260C	
180-42975-1 MS	HD-MW-165-0/1-0	Total/NA	Water	8260C	
180-42975-1 MSD	HD-MW-165-0/1-0	Total/NA	Water	8260C	
180-42975-2	HD-MW-162-0/1-0	Total/NA	Water	8260C	
180-42975-4	HD-QC2-0/1-2	Total/NA	Water	8260C	
LCS 180-138583/8	Lab Control Sample	Total/NA	Water	8260C	
MB 180-138583/5	Method Blank	Total/NA	Water	8260C	

### Analysis Batch: 138685

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
180-42975-2 - DL	HD-MW-162-0/1-0	Total/NA	Water	8260C	
180-42975-3	HD-MW-169-0/1-0	Total/NA	Water	8260C	
LCS 180-138685/8	Lab Control Sample	Total/NA	Water	8260C	
LCSD 180-138685/10	Lab Control Sample Dup	Total/NA	Water	8260C	
MB 180-138685/5	Method Blank	Total/NA	Water	8260C	

# Lab Chronicle

Client: Groundwater Sciences Corporation  
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

**Client Sample ID: HD-MW-165-0/1-0**

**Lab Sample ID: 180-42975-1**

**Matrix: Water**

Date Collected: 04/10/15 08:40  
 Date Received: 04/11/15 09:30

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	8260C		1	5 mL	5 mL	138583	04/15/15 15:09	DLF	TAL PIT

Instrument ID: CHHP5

**Client Sample ID: HD-MW-162-0/1-0**

**Lab Sample ID: 180-42975-2**

**Matrix: Water**

Date Collected: 04/10/15 09:50  
 Date Received: 04/11/15 09:30

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	8260C		4	5 mL	5 mL	138583	04/15/15 20:47	DLF	TAL PIT
		Instrument ID: CHHP5								
Total/NA	Analysis	8260C	DL	40	5 mL	5 mL	138685	04/16/15 16:29	DLF	TAL PIT
		Instrument ID: CHHP5								

**Client Sample ID: HD-MW-169-0/1-0**

**Lab Sample ID: 180-42975-3**

**Matrix: Water**

Date Collected: 04/10/15 12:32  
 Date Received: 04/11/15 09:30

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	8260C		1	5 mL	5 mL	138685	04/16/15 16:53	DLF	TAL PIT

Instrument ID: CHHP5

**Client Sample ID: HD-QC2-0/1-2**

**Lab Sample ID: 180-42975-4**

**Matrix: Water**

Date Collected: 04/10/15 12:00  
 Date Received: 04/11/15 09:30

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	8260C		1	5 mL	5 mL	138583	04/15/15 15:33	DLF	TAL PIT

Instrument ID: CHHP5

**Laboratory References:**

TAL PIT = TestAmerica Pittsburgh, 301 Alpha Drive, RIDC Park, Pittsburgh, PA 15238, TEL (412)963-7058

**Analyst References:**

Lab: TAL PIT

Batch Type: Analysis

DLF = Donald Ferguson

TestAmerica Pittsburgh

## Certification Summary

Client: Groundwater Sciences Corporation  
Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

### Laboratory: TestAmerica Pittsburgh

The certifications listed below are applicable to this report.

Authority	Program	EPA Region	Certification ID	Expiration Date
Pennsylvania	NELAP	3	02-00416	04-30-15 *

\* Certification renewal pending - certification considered valid.

## Method Summary

Client: Groundwater Sciences Corporation  
Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

Method	Method Description	Protocol	Laboratory
8260C	Volatile Organic Compounds (GC/MS)	SW846	TAL PIT

**Protocol References:**

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

**Laboratory References:**

TAL PIT = TestAmerica Pittsburgh, 301 Alpha Drive, RIDC Park, Pittsburgh, PA 15238, TEL (412)963-7058

## Sample Summary

Client: Groundwater Sciences Corporation  
Project/Site: Harley Davidson

TestAmerica Job ID: 180-42975-1

Lab Sample ID	Client Sample ID	Matrix	Collected	Received
180-42975-1	HD-MW-165-0/1-0	Water	04/10/15 08:40	04/11/15 09:30
180-42975-2	HD-MW-162-0/1-0	Water	04/10/15 09:50	04/11/15 09:30
180-42975-3	HD-MW-169-0/1-0	Water	04/10/15 12:32	04/11/15 09:30
180-42975-4	HD-QC2-0/1-2	Water	04/10/15 12:00	04/11/15 09:30

## GC/MS VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.:

Instrument ID: CHHP5 Analysis Batch Number: 135593

Lab Sample ID: IC 180-135593/4 Client Sample ID:

Date Analyzed: 03/16/15 12:41 Lab File ID: 50316004.D GC Column: DB-624 ID: 0.18 (mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Bromomethane	2.25	Baseline	fergusond	03/17/15 09:42

Lab Sample ID: ICIS 180-135593/5 Client Sample ID:

Date Analyzed: 03/16/15 13:05 Lab File ID: 50316005.D GC Column: DB-624 ID: 0.18 (mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
1,4-Dioxane	8.06	Peak Tail	fergusond	03/17/15 09:27

Lab Sample ID: IC 180-135593/6 Client Sample ID:

Date Analyzed: 03/16/15 13:29 Lab File ID: 50316006.D GC Column: DB-624 ID: 0.18 (mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Isobutyl alcohol	6.95	Peak Tail	fergusond	03/17/15 09:45

Lab Sample ID: IC 180-135593/7 Client Sample ID:

Date Analyzed: 03/16/15 13:53 Lab File ID: 50316007.D GC Column: DB-624 ID: 0.18 (mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Isobutyl alcohol	6.94	Peak Tail	fergusond	03/17/15 09:48

## GC/MS VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.:

Instrument ID: CHHP5

Analysis Batch Number: 135593

Lab Sample ID: IC 180-135593/13

Client Sample ID:

Date Analyzed: 03/16/15 16:17

Lab File ID: 50316013.D

GC Column: DB-624

ID: 0.18 (mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Chloromethane	1.77	Poor chromatography	fergusond	03/17/15 10:01
Acrolein	3.25	Poor chromatography	fergusond	03/17/15 10:01
2-Hexanone	9.66	Poor chromatography	fergusond	03/17/15 10:01
trans-1,4-Dichloro-2-butene	11.74	Poor chromatography	fergusond	03/17/15 10:01

## GC/MS VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.:

Instrument ID: CHHP5 Analysis Batch Number: 138583

Lab Sample ID: CCVIS 180-138583/2 Client Sample ID:

Date Analyzed: 04/15/15 13:21 Lab File ID: 50415002.D GC Column: DB-624 ID: 0.18 (mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
1,4-Dioxane	8.05	Peak Tail	fergusond	04/15/15 13:44

Lab Sample ID: 180-42975-1 Client Sample ID: HD-MW-165-0/1-0

Date Analyzed: 04/15/15 15:09 Lab File ID: 50415006.D GC Column: DB-624 ID: 0.18 (mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Methylene Chloride	4.15	Split Peak	fergusond	04/16/15 07:33

Lab Sample ID: 180-42975-1 MS Client Sample ID: HD-MW-165-0/1-0 MS

Date Analyzed: 04/15/15 16:21 Lab File ID: 50415009.D GC Column: DB-624 ID: 0.18 (mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
1,4-Dioxane	8.06	Peak Tail	fergusond	04/16/15 07:36

Lab Sample ID: 180-42975-1 MSD Client Sample ID: HD-MW-165-0/1-0 MSD

Date Analyzed: 04/15/15 16:46 Lab File ID: 50415010.D GC Column: DB-624 ID: 0.18 (mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
1,4-Dioxane	8.05	Peak Tail	fergusond	04/16/15 07:37

Lab Sample ID: 180-42975-2 Client Sample ID: HD-MW-162-0/1-0

Date Analyzed: 04/15/15 20:47 Lab File ID: 50415020.D GC Column: DB-624 ID: 0.18 (mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
1,1-Dichloroethene	3.40	Poor chromatography	fergusond	04/16/15 08:55
Methylene Chloride	4.15	Poor chromatography	fergusond	04/16/15 08:55
1,1,1-Trichloroethane	6.53	Poor chromatography	fergusond	04/16/15 08:55

## GC/MS VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.:

Instrument ID: CHHP5 Analysis Batch Number: 138685

Lab Sample ID: CCVIS 180-138685/2 Client Sample ID:

Date Analyzed: 04/16/15 10:19 Lab File ID: 50416002.D GC Column: DB-624 ID: 0.18 (mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Iodomethane	3.62	Poor chromatography	fergusond	04/16/15 11:02

Lab Sample ID: MB 180-138685/5 Client Sample ID:

Date Analyzed: 04/16/15 11:32 Lab File ID: 50416005.D GC Column: DB-624 ID: 0.18 (mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Methylene Chloride	4.18	Split Peak	fergusond	04/16/15 11:51

Lab Sample ID: LCS 180-138685/8 Client Sample ID:

Date Analyzed: 04/16/15 13:05 Lab File ID: 50416008.D GC Column: DB-624 ID: 0.18 (mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
1,4-Dioxane	8.06	Peak Tail	fergusond	04/16/15 13:37

Lab Sample ID: LCSD 180-138685/10 Client Sample ID:

Date Analyzed: 04/16/15 13:53 Lab File ID: 50416010.D GC Column: DB-624 ID: 0.18 (mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
1,4-Dioxane	8.06	Peak Tail	fergusond	04/16/15 14:11

Lab Sample ID: 180-42975-2 DL Client Sample ID: HD-MW-162-0/1-0 DL

Date Analyzed: 04/16/15 16:29 Lab File ID: 50416016.D GC Column: DB-624 ID: 0.18 (mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Methylene Chloride	4.13	Poor chromatography	fergusond	04/16/15 16:58

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
VOA8260INT_00030	04/10/15	03/10/15	Methanol, Lot 85233	10 mL	VOA8260INTRES_00091	1 mL	1,4-Dichlorobenzene-d4	25 ug/mL
							Chlorobenzene-d5	25 ug/mL
							Fluorobenzene (IS)	25 ug/mL
							TBA-d9 (IS)	500 ug/mL
.VOA8260INTRES_00091	07/31/19		Restek, Lot A0104742		(Purchased Reagent)		1,4-Dichlorobenzene-d4	250 ug/mL
							Chlorobenzene-d5	250 ug/mL
							Fluorobenzene (IS)	250 ug/mL
							TBA-d9 (IS)	5000 ug/mL
VOA8260SURR_00032	04/10/15	03/10/15	Methanol, Lot 85233	100 mL	VOA8260SURRES_00063	1 mL	1,2-Dichloroethane-d4 (Surr)	25 ug/mL
							4-Bromofluorobenzene (Surr)	25 ug/mL
							Dibromofluoromethane (Surr)	25 ug/mL
							Toluene-d8 (Surr)	25 ug/mL
.VOA8260SURRES_00063	01/31/19		Restek, Lot A0100424		(Purchased Reagent)		1,2-Dichloroethane-d4 (Surr)	2500 ug/mL
							4-Bromofluorobenzene (Surr)	2500 ug/mL
							Dibromofluoromethane (Surr)	2500 ug/mL
							Toluene-d8 (Surr)	2500 ug/mL
VOA8260SURR_00033	05/03/15	04/03/15	Methanol, Lot 85233	100 mL	VOA8260SURRES_00087	1 mL	1,2-Dichloroethane-d4 (Surr)	25 ug/mL
							4-Bromofluorobenzene (Surr)	25 ug/mL
							Dibromofluoromethane (Surr)	25 ug/mL
							Toluene-d8 (Surr)	25 ug/mL
.VOA8260SURRES_00087	04/30/19		Restek, Lot A0102817		(Purchased Reagent)		1,2-Dichloroethane-d4 (Surr)	2500 ug/mL
							4-Bromofluorobenzene (Surr)	2500 ug/mL
							Dibromofluoromethane (Surr)	2500 ug/mL
							Toluene-d8 (Surr)	2500 ug/mL
VOA8260VOA2ND_00111	04/22/16	04/15/15	Methanol, Lot 85233	10 mL	VOA8260GAS2ND_00093	0.1 mL	Bromomethane	25 ug/mL
							Chloroethane	25 ug/mL
							Chloromethane	25 ug/mL
							Vinyl chloride	25 ug/mL
					VOA8260VOA2ND_00107	1.25 mL	1,1,1,2-Tetrachloroethane	25 ug/mL
							1,1,1-Trichloroethane	25 ug/mL
							1,1,2,2-Tetrachloroethane	25 ug/mL
							1,1,2-Trichloroethane	25 ug/mL
							1,1-Dichloroethane	25 ug/mL
							1,1-Dichloroethene	25 ug/mL
							1,2-Dibromoethane (EDB)	25 ug/mL
							1,2-Dichloroethane	25 ug/mL
							1,2-Dichloropropane	25 ug/mL
							1,4-Dioxane	500 ug/mL
							Acrylonitrile	250 ug/mL
							Benzene	25 ug/mL
							Bromochloromethane	25 ug/mL
							Bromodichloromethane	25 ug/mL
							Bromoform	25 ug/mL
							Carbon disulfide	25 ug/mL
							Carbon tetrachloride	25 ug/mL
							Chlorobenzene	25 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Chloroform	25 ug/mL
							cis-1,2-Dichloroethene	25 ug/mL
							cis-1,3-Dichloropropene	25 ug/mL
							Dibromochloromethane	25 ug/mL
							Ethylbenzene	25 ug/mL
							Methyl tert-butyl ether	25 ug/mL
							Methylene Chloride	25 ug/mL
							Styrene	25 ug/mL
							Tetrachloroethene	25 ug/mL
							Toluene	25 ug/mL
							trans-1,2-Dichloroethene	25 ug/mL
							trans-1,3-Dichloropropene	25 ug/mL
							Trichloroethene	25 ug/mL
							Xylenes, Total	50 ug/mL
.VOA8260GAS2ND_00093	01/31/18	Restek, Lot A0108226			(Purchased Reagent)		Bromomethane	2500 ug/mL
							Chloroethane	2500 ug/mL
							Chloromethane	2500 ug/mL
							Vinyl chloride	2500 ug/mL
.VOA8260VOA2ND_00107	04/19/15	03/19/15	Methanol, Lot 85233	10 mL	VOA8260MEGA2_00011	1 mL	1,1,1,2-Tetrachloroethane	200 ug/mL
							1,1,1-Trichloroethane	200 ug/mL
							1,1,2,2-Tetrachloroethane	200 ug/mL
							1,1,2-Trichloroethane	200 ug/mL
							1,1-Dichloroethane	200 ug/mL
							1,1-Dichloroethene	200 ug/mL
							1,2-Dibromoethane (EDB)	200 ug/mL
							1,2-Dichloroethane	200 ug/mL
							1,2-Dichloropropane	200 ug/mL
							1,4-Dioxane	4000 ug/mL
							Acrylonitrile	2000 ug/mL
							Benzene	200 ug/mL
							Bromochloromethane	200 ug/mL
							Bromodichloromethane	200 ug/mL
							Bromoform	200 ug/mL
							Carbon disulfide	200 ug/mL
							Carbon tetrachloride	200 ug/mL
							Chlorobenzene	200 ug/mL
							Chloroform	200 ug/mL
							cis-1,2-Dichloroethene	200 ug/mL
							cis-1,3-Dichloropropene	200 ug/mL
							Dibromochloromethane	200 ug/mL
							Ethylbenzene	200 ug/mL
							Methyl tert-butyl ether	200 ug/mL
							Methylene Chloride	200 ug/mL
							Styrene	200 ug/mL
							Tetrachloroethene	200 ug/mL
							Toluene	200 ug/mL
							trans-1,2-Dichloroethene	200 ug/mL
							trans-1,3-Dichloropropene	200 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..VOA8260MEGA2_00011	02/01/16		Restek, Lot A093733		(Purchased Reagent)		Trichloroethene	200 ug/mL
							Xylenes, Total	400 ug/mL
							1,1,1,2-Tetrachloroethane	2000 ug/mL
							1,1,1-Trichloroethane	2000 ug/mL
							1,1,2,2-Tetrachloroethane	2000 ug/mL
							1,1,2-Trichloroethane	2000 ug/mL
							1,1-Dichloroethane	2000 ug/mL
							1,1-Dichloroethene	2000 ug/mL
							1,2-Dibromoethane (EDB)	2000 ug/mL
							1,2-Dichloroethane	2000 ug/mL
							1,2-Dichloropropane	2000 ug/mL
							1,4-Dioxane	40000 ug/mL
							Acrylonitrile	20000 ug/mL
							Benzene	2000 ug/mL
							Bromochloromethane	2000 ug/mL
							Bromodichloromethane	2000 ug/mL
							Bromoform	2000 ug/mL
							Carbon disulfide	2000 ug/mL
							Carbon tetrachloride	2000 ug/mL
							Chlorobenzene	2000 ug/mL
							Chloroform	2000 ug/mL
							cis-1,2-Dichloroethene	2000 ug/mL
							cis-1,3-Dichloropropene	2000 ug/mL
							Dibromochloromethane	2000 ug/mL
							Ethylbenzene	2000 ug/mL
							Methyl tert-butyl ether	2000 ug/mL
							Methylene Chloride	2000 ug/mL
							Styrene	2000 ug/mL
							Tetrachloroethene	2000 ug/mL
							Toluene	2000 ug/mL
							trans-1,2-Dichloroethene	2000 ug/mL
							trans-1,3-Dichloropropene	2000 ug/mL
							Trichloroethene	2000 ug/mL
							Xylenes, Total	4000 ug/mL
VOA8260VOAPRI_00105	03/20/15	03/13/15	Methanol, Lot 85233	8 mL	VOA8260GAS1ST_00091	0.08 mL	Bromomethane	25 ug/mL
							Butadiene	25 ug/mL
							Chloroethane	25 ug/mL
							Chloromethane	25 ug/mL
							Dichlorodifluoromethane	25 ug/mL
							Dichlorofluoromethane	25 ug/mL
							Trichlorofluoromethane	25 ug/mL
							Vinyl chloride	25 ug/mL
					VOA8260VOAPRI_00101	1 mL	2-Butanone (MEK)	25 ug/mL
							2-Hexanone	25 ug/mL
							4-Methyl-2-pentanone (MIBK)	25 ug/mL
							Acetone	25 ug/mL
							1,1,1,2-Tetrachloroethane	25 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							1,1,1-Trichloroethane	25 ug/mL
							1,1,2,2-Tetrachloroethane	25 ug/mL
							1,1,2-Trichloro-1,2,2-trifluor oethane	25 ug/mL
							1,1,2-Trichloroethane	25 ug/mL
							1,1-Dichloroethane	25 ug/mL
							1,1-Dichloroethene	25 ug/mL
							1,1-Dichloropropene	25 ug/mL
							1,2,3-Trichlorobenzene	25 ug/mL
							1,2,3-Trichloropropane	25 ug/mL
							1,2,4-Trichlorobenzene	25 ug/mL
							1,2,4-Trimethylbenzene	25 ug/mL
							1,2-Dibromo-3-Chloropropane	25 ug/mL
							1,2-Dibromoethane (EDB)	25 ug/mL
							1,2-Dichlorobenzene	25 ug/mL
							1,2-Dichloroethane	25 ug/mL
							1,2-Dichloropropane	25 ug/mL
							1,3,5-Trimethylbenzene	25 ug/mL
							1,3-Dichlorobenzene	25 ug/mL
							1,3-Dichloropropane	25 ug/mL
							1,4-Dichlorobenzene	25 ug/mL
							1,4-Dioxane	500 ug/mL
							2,2-Dichloropropane	25 ug/mL
							2-Chlorotoluene	25 ug/mL
							2-Methyl-2-propanol	250 ug/mL
							3-Chloro-1-propene	25 ug/mL
							4-Chlorotoluene	25 ug/mL
							4-Isopropyltoluene	25 ug/mL
							Acrylonitrile	250 ug/mL
							Benzene	25 ug/mL
							Bromobenzene	25 ug/mL
							Bromochloromethane	25 ug/mL
							Bromodichloromethane	25 ug/mL
							Bromoform	25 ug/mL
							Carbon disulfide	25 ug/mL
							Carbon tetrachloride	25 ug/mL
							Chlorobenzene	25 ug/mL
							Chloroform	25 ug/mL
							cis-1,2-Dichloroethene	25 ug/mL
							cis-1,3-Dichloropropene	25 ug/mL
							Cyclohexane	25 ug/mL
							Dibromochloromethane	25 ug/mL
							Dibromomethane	25 ug/mL
							Ethyl ether	25 ug/mL
							Ethyl methacrylate	25 ug/mL
							Ethylbenzene	25 ug/mL
							Hexachlorobutadiene	25 ug/mL
							Hexane	25 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Iodomethane	25 ug/mL
							Isobutyl alcohol	625 ug/mL
							Isopropylbenzene	25 ug/mL
							m-Xylene & p-Xylene	25 ug/mL
							Methyl acetate	125 ug/mL
							Methyl tert-butyl ether	25 ug/mL
							Methylcyclohexane	25 ug/mL
							Methylene Chloride	25 ug/mL
							n-Butylbenzene	25 ug/mL
							n-Heptane	25 ug/mL
							N-Propylbenzene	25 ug/mL
							Naphthalene	25 ug/mL
							o-Xylene	25 ug/mL
							sec-Butylbenzene	25 ug/mL
							Styrene	25 ug/mL
							tert-Butylbenzene	25 ug/mL
							Tetrachloroethene	25 ug/mL
							Tetrahydrofuran	50 ug/mL
							Toluene	25 ug/mL
							trans-1,2-Dichloroethene	25 ug/mL
							trans-1,3-Dichloropropene	25 ug/mL
							trans-1,4-Dichloro-2-butene	25 ug/mL
							Trichloroethene	25 ug/mL
.VOA8260GAS1ST_00091	09/30/16	Restek, Lot A0108198			(Purchased Reagent)		Bromomethane	2500 ug/mL
							Butadiene	2500 ug/mL
							Chloroethane	2500 ug/mL
							Chloromethane	2500 ug/mL
							Dichlorodifluoromethane	2500 ug/mL
							Dichlorofluoromethane	2500 ug/mL
							Trichlorofluoromethane	2500 ug/mL
							Vinyl chloride	2500 ug/mL
.VOA8260VOAPRI_00101	03/24/15	02/24/15	Methanol, Lot 85233	10 mL	VOA8260KET1ST_00036	0.2 mL	2-Butanone (MEK)	200 ug/mL
							2-Hexanone	200 ug/mL
							4-Methyl-2-pentanone (MIBK)	200 ug/mL
							Acetone	200 ug/mL
					VOA8260MEGA1_00027	1 mL	1,1,1,2-Tetrachloroethane	200 ug/mL
							1,1,1-Trichloroethane	200 ug/mL
							1,1,2,2-Tetrachloroethane	200 ug/mL
							1,1,2-Trichloro-1,2,2-trifluoroethane	200 ug/mL
							1,1,2-Trichloroethane	200 ug/mL
							1,1-Dichloroethane	200 ug/mL
							1,1-Dichloroethene	200 ug/mL
							1,1-Dichloropropene	200 ug/mL
							1,2,3-Trichlorobenzene	200 ug/mL
							1,2,3-Trichloropropane	200 ug/mL
							1,2,4-Trichlorobenzene	200 ug/mL
							1,2,4-Trimethylbenzene	200 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							1,2-Dibromo-3-Chloropropane	200 ug/mL
							1,2-Dibromoethane (EDB)	200 ug/mL
							1,2-Dichlorobenzene	200 ug/mL
							1,2-Dichloroethane	200 ug/mL
							1,2-Dichloropropane	200 ug/mL
							1,3,5-Trimethylbenzene	200 ug/mL
							1,3-Dichlorobenzene	200 ug/mL
							1,3-Dichloropropane	200 ug/mL
							1,4-Dichlorobenzene	200 ug/mL
							1,4-Dioxane	4000 ug/mL
							2,2-Dichloropropane	200 ug/mL
							2-Chlorotoluene	200 ug/mL
							2-Methyl-2-propanol	2000 ug/mL
							3-Chloro-1-propene	200 ug/mL
							4-Chlorotoluene	200 ug/mL
							4-Isopropyltoluene	200 ug/mL
							Acrylonitrile	2000 ug/mL
							Benzene	200 ug/mL
							Bromobenzene	200 ug/mL
							Bromochloromethane	200 ug/mL
							Bromodichloromethane	200 ug/mL
							Bromoform	200 ug/mL
							Carbon disulfide	200 ug/mL
							Carbon tetrachloride	200 ug/mL
							Chlorobenzene	200 ug/mL
							Chloroform	200 ug/mL
							cis-1,2-Dichloroethene	200 ug/mL
							cis-1,3-Dichloropropene	200 ug/mL
							Cyclohexane	200 ug/mL
							Dibromochloromethane	200 ug/mL
							Dibromomethane	200 ug/mL
							Ethyl ether	200 ug/mL
							Ethyl methacrylate	200 ug/mL
							Ethylbenzene	200 ug/mL
							Hexachlorobutadiene	200 ug/mL
							Hexane	200 ug/mL
							Iodomethane	200 ug/mL
							Isobutyl alcohol	5000 ug/mL
							Isopropylbenzene	200 ug/mL
							m-Xylene & p-Xylene	200 ug/mL
							Methyl acetate	1000 ug/mL
							Methyl tert-butyl ether	200 ug/mL
							Methylcyclohexane	200 ug/mL
							Methylene Chloride	200 ug/mL
							n-Butylbenzene	200 ug/mL
							n-Heptane	200 ug/mL
							N-Propylbenzene	200 ug/mL
							Naphthalene	200 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							o-Xylene	200 ug/mL
							sec-Butylbenzene	200 ug/mL
							Styrene	200 ug/mL
							tert-Butylbenzene	200 ug/mL
							Tetrachloroethene	200 ug/mL
							Tetrahydrofuran	400 ug/mL
							Toluene	200 ug/mL
							trans-1,2-Dichloroethene	200 ug/mL
							trans-1,3-Dichloropropene	200 ug/mL
							trans-1,4-Dichloro-2-butene	200 ug/mL
							Trichloroethene	200 ug/mL
..VOA8260KET1ST_00036	02/28/16	Restek, Lot A093365			(Purchased Reagent)		2-Butanone (MEK)	10000 ug/mL
							2-Hexanone	10000 ug/mL
							4-Methyl-2-pentanone (MIBK)	10000 ug/mL
							Acetone	10000 ug/mL
..VOA8260MEGA1_00027	02/28/16	Restek, Lot A093581			(Purchased Reagent)		1,1,1,2-Tetrachloroethane	2000 ug/mL
							1,1,1-Trichloroethane	2000 ug/mL
							1,1,2,2-Tetrachloroethane	2000 ug/mL
							1,1,2-Trichloro-1,2,2-trifluoroethane	2000 ug/mL
							1,1,2-Trichloroethane	2000 ug/mL
							1,1-Dichloroethane	2000 ug/mL
							1,1-Dichloroethene	2000 ug/mL
							1,1-Dichloropropene	2000 ug/mL
							1,2,3-Trichlorobenzene	2000 ug/mL
							1,2,3-Trichloropropane	2000 ug/mL
							1,2,4-Trichlorobenzene	2000 ug/mL
							1,2,4-Trimethylbenzene	2000 ug/mL
							1,2-Dibromo-3-Chloropropane	2000 ug/mL
							1,2-Dibromoethane (EDB)	2000 ug/mL
							1,2-Dichlorobenzene	2000 ug/mL
							1,2-Dichloroethane	2000 ug/mL
							1,2-Dichloropropane	2000 ug/mL
							1,3,5-Trimethylbenzene	2000 ug/mL
							1,3-Dichlorobenzene	2000 ug/mL
							1,3-Dichloropropane	2000 ug/mL
							1,4-Dichlorobenzene	2000 ug/mL
							1,4-Dioxane	40000 ug/mL
							2,2-Dichloropropane	2000 ug/mL
							2-Chlorotoluene	2000 ug/mL
							2-Methyl-2-propanol	20000 ug/mL
							3-Chloro-1-propene	2000 ug/mL
							4-Chlorotoluene	2000 ug/mL
							4-Isopropyltoluene	2000 ug/mL
							Acrylonitrile	20000 ug/mL
							Benzene	2000 ug/mL
							Bromobenzene	2000 ug/mL
							Bromochloromethane	2000 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					Bromodichloromethane	2000 ug/mL		
					Bromoform	2000 ug/mL		
					Carbon disulfide	2000 ug/mL		
					Carbon tetrachloride	2000 ug/mL		
					Chlorobenzene	2000 ug/mL		
					Chloroform	2000 ug/mL		
					cis-1,2-Dichloroethene	2000 ug/mL		
					cis-1,3-Dichloropropene	2000 ug/mL		
					Cyclohexane	2000 ug/mL		
					Dibromochloromethane	2000 ug/mL		
					Dibromomethane	2000 ug/mL		
					Ethyl ether	2000 ug/mL		
					Ethyl methacrylate	2000 ug/mL		
					Ethylbenzene	2000 ug/mL		
					Hexachlorobutadiene	2000 ug/mL		
					Hexane	2000 ug/mL		
					Iodomethane	2000 ug/mL		
					Isobutyl alcohol	50000 ug/mL		
					Isopropylbenzene	2000 ug/mL		
					m-Xylene & p-Xylene	2000 ug/mL		
					Methyl acetate	10000 ug/mL		
					Methyl tert-butyl ether	2000 ug/mL		
					Methylcyclohexane	2000 ug/mL		
					Methylene Chloride	2000 ug/mL		
					n-Butylbenzene	2000 ug/mL		
					n-Heptane	2000 ug/mL		
					N-Propylbenzene	2000 ug/mL		
					Naphthalene	2000 ug/mL		
					o-Xylene	2000 ug/mL		
					sec-Butylbenzene	2000 ug/mL		
					Styrene	2000 ug/mL		
					tert-Butylbenzene	2000 ug/mL		
					Tetrachloroethene	2000 ug/mL		
					Tetrahydrofuran	4000 ug/mL		
					Toluene	2000 ug/mL		
					trans-1,2-Dichloroethene	2000 ug/mL		
					trans-1,3-Dichloropropene	2000 ug/mL		
					trans-1,4-Dichloro-2-butene	2000 ug/mL		
					Trichloroethene	2000 ug/mL		
VOA8260VOAPRI_00110	04/22/15	04/15/15	Methanol, Lot 85233	10 mL	VOA8260GAS1ST_00095	0.1 mL	Bromomethane	25 ug/mL
							Chloroethane	25 ug/mL
							Chloromethane	25 ug/mL
							Vinyl chloride	25 ug/mL
							VOA8260VOAPRI_00106	1.25 mL
							1,1,1,2-Tetrachloroethane	25 ug/mL
							1,1,1-Trichloroethane	25 ug/mL
							1,1,2,2-Tetrachloroethane	25 ug/mL
							1,1,2-Trichloroethane	25 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							1,1-Dichloroethane	25 ug/mL
							1,1-Dichloroethene	25 ug/mL
							1,2-Dibromoethane (EDB)	25 ug/mL
							1,2-Dichloroethane	25 ug/mL
							1,2-Dichloropropane	25 ug/mL
							1,4-Dioxane	500 ug/mL
							Acrylonitrile	250 ug/mL
							Benzene	25 ug/mL
							Bromochloromethane	25 ug/mL
							Bromodichloromethane	25 ug/mL
							Bromoform	25 ug/mL
							Carbon disulfide	25 ug/mL
							Carbon tetrachloride	25 ug/mL
							Chlorobenzene	25 ug/mL
							Chloroform	25 ug/mL
							cis-1,2-Dichloroethene	25 ug/mL
							cis-1,3-Dichloropropene	25 ug/mL
							Dibromochloromethane	25 ug/mL
							Ethylbenzene	25 ug/mL
							Methyl tert-butyl ether	25 ug/mL
							Methylene Chloride	25 ug/mL
							Styrene	25 ug/mL
							Tetrachloroethene	25 ug/mL
							Toluene	25 ug/mL
							trans-1,2-Dichloroethene	25 ug/mL
							trans-1,3-Dichloropropene	25 ug/mL
							Trichloroethene	25 ug/mL
							Xylenes, Total	50 ug/mL
.VOA8260GAS1ST_00095	01/31/18	Restek, Lot A0108198			(Purchased Reagent)		Bromomethane	2500 ug/mL
.VOA8260VOAPRI_00106	04/19/15	03/19/15	Methanol, Lot 85233	10 mL	VOA8260MEGA1_00014	1 mL	Chloroethane	2500 ug/mL
							Chloromethane	2500 ug/mL
							Vinyl chloride	2500 ug/mL
							1,1,1,2-Tetrachloroethane	200 ug/mL
							1,1,1-Trichloroethane	200 ug/mL
							1,1,2,2-Tetrachloroethane	200 ug/mL
							1,1,2-Trichloroethane	200 ug/mL
							1,1-Dichloroethane	200 ug/mL
							1,1-Dichloroethene	200 ug/mL
							1,2-Dibromoethane (EDB)	200 ug/mL
							1,2-Dichloroethane	200 ug/mL
							1,2-Dichloropropane	200 ug/mL
							1,4-Dioxane	4000 ug/mL
							Acrylonitrile	2000 ug/mL
							Benzene	200 ug/mL
							Bromochloromethane	200 ug/mL
							Bromodichloromethane	200 ug/mL
							Bromoform	200 ug/mL
							Carbon disulfide	200 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Carbon tetrachloride	200 ug/mL
							Chlorobenzene	200 ug/mL
							Chloroform	200 ug/mL
							cis-1,2-Dichloroethene	200 ug/mL
							cis-1,3-Dichloropropene	200 ug/mL
							Dibromochloromethane	200 ug/mL
							Ethylbenzene	200 ug/mL
							Methyl tert-butyl ether	200 ug/mL
							Methylene Chloride	200 ug/mL
							Styrene	200 ug/mL
							Tetrachloroethene	200 ug/mL
							Toluene	200 ug/mL
							trans-1,2-Dichloroethene	200 ug/mL
							trans-1,3-Dichloropropene	200 ug/mL
							Trichloroethene	200 ug/mL
							Xylenes, Total	400 ug/mL
..VOA8260MEGA1_00014	02/28/16	Restek, Lot A093581		(Purchased Reagent)			1,1,1,2-Tetrachloroethane	2000 ug/mL
							1,1,1-Trichloroethane	2000 ug/mL
							1,1,2,2-Tetrachloroethane	2000 ug/mL
							1,1,2-Trichloroethane	2000 ug/mL
							1,1-Dichloroethane	2000 ug/mL
							1,1-Dichloroethene	2000 ug/mL
							1,2-Dibromoethane (EDB)	2000 ug/mL
							1,2-Dichloroethane	2000 ug/mL
							1,2-Dichloropropane	2000 ug/mL
							1,4-Dioxane	40000 ug/mL
							Acrylonitrile	20000 ug/mL
							Benzene	2000 ug/mL
							Bromochloromethane	2000 ug/mL
							Bromodichloromethane	2000 ug/mL
							Bromoform	2000 ug/mL
							Carbon disulfide	2000 ug/mL
							Carbon tetrachloride	2000 ug/mL
							Chlorobenzene	2000 ug/mL
							Chloroform	2000 ug/mL
							cis-1,2-Dichloroethene	2000 ug/mL
							cis-1,3-Dichloropropene	2000 ug/mL
							Dibromochloromethane	2000 ug/mL
							Ethylbenzene	2000 ug/mL
							Methyl tert-butyl ether	2000 ug/mL
							Methylene Chloride	2000 ug/mL
							Styrene	2000 ug/mL
							Tetrachloroethene	2000 ug/mL
							Toluene	2000 ug/mL
							trans-1,2-Dichloroethene	2000 ug/mL
							trans-1,3-Dichloropropene	2000 ug/mL
							Trichloroethene	2000 ug/mL
							Xylenes, Total	4000 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
VOAACRPRI_00003	03/31/15	03/03/15	Methanol, Lot 85233	100 mL	VOAACRORES_00064	0.125 mL	Acrolein	25 ug/mL
.VOAACRORES_00064	03/31/15		Restek, Lot A0107338		(Purchased Reagent)		Acrolein	20000 ug/mL
VOAVAPRI_00005	04/13/15	03/13/15	Methanol, Lot 85233	50 mL	VOA8260VARES_00050	0.25 mL	Vinyl acetate	25 ug/mL
.VOA8260VARES_00050	07/31/15		Restek, Lot A0108225		(Purchased Reagent)		Vinyl acetate	5000 ug/mL
voaWEpri Res_00003	03/30/15	03/02/15	Methanol, Lot 85233	25 mL	VOARESEE1ST_00008	0.125 mL	1,2-dichloro-4-(trifluoromethyl)benzene 2,3,6-Trichlorotoluene 2,4,5-Trichlorotoluene 2,4-Dichloro-1-(trifluoromethyl)-benzene 2,5-Dichlorobenzotrifluoride 2-Chlorobenzotrifluoride 3-Chlorobenzotrifluoride 3-Chlorotoluene 4-Chlorobenzotrifluoride	25 ug/mL 25 ug/mL 25 ug/mL 25 ug/mL 25 ug/mL 25 ug/mL 25 ug/mL 25 ug/mL 25 ug/mL
.VOARESEE1ST_00008	02/28/15		Restek, Lot A097285		(Purchased Reagent)		1,2-dichloro-4-(trifluoromethyl)benzene 2,3,6-Trichlorotoluene 2,4,5-Trichlorotoluene 2,4-Dichloro-1-(trifluoromethyl)-benzene 2,5-Dichlorobenzotrifluoride 2-Chlorobenzotrifluoride 3-Chlorobenzotrifluoride 3-Chlorotoluene 4-Chlorobenzotrifluoride	5000 ug/mL 5000 ug/mL 5000 ug/mL 5000 ug/mL 5000 ug/mL 5000 ug/mL 5000 ug/mL 5000 ug/mL 5000 ug/mL
voaWKet2 Rest_00002	04/16/15	03/16/15	Methanol, Lot 85233	50 mL	VOA8260KET2ND_00042	0.1 mL	2-Butanone (MEK) 2-Hexanone 4-Methyl-2-pentanone (MIBK) Acetone	25 ug/mL 25 ug/mL 25 ug/mL 25 ug/mL
.VOA8260KET2ND_00042	01/31/18		Restek, Lot A0108157		(Purchased Reagent)		2-Butanone (MEK) 2-Hexanone 4-Methyl-2-pentanone (MIBK) Acetone	12500 ug/mL 12500 ug/mL 12500 ug/mL 12500 ug/mL
voaWKetpri Re_00003	03/26/15	02/24/15	Methanol, Lot 85233	50 mL	VOA8260KET1ST_00037	0.125 mL	2-Butanone (MEK) 2-Hexanone 4-Methyl-2-pentanone (MIBK) Acetone	25 ug/mL 25 ug/mL 25 ug/mL 25 ug/mL
.VOA8260KET1ST_00037	02/28/16		Restek, Lot A093365		(Purchased Reagent)		2-Butanone (MEK) 2-Hexanone 4-Methyl-2-pentanone (MIBK) Acetone	10000 ug/mL 10000 ug/mL 10000 ug/mL 10000 ug/mL
voaWKetpri Re_00004	04/30/15	03/30/15	Methanol, Lot 85233	50 mL	VOA8260KET1ST_00039	0.1 mL	2-Butanone (MEK) 2-Hexanone 4-Methyl-2-pentanone (MIBK) Acetone	25 ug/mL 25 ug/mL 25 ug/mL 25 ug/mL
.VOA8260KET1ST_00039	01/31/18		Restek, Lot A0108151		(Purchased Reagent)		2-Butanone (MEK)	12500 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2-Hexanone	12500 ug/mL
							4-Methyl-2-pentanone (MIBK)	12500 ug/mL
							Acetone	12500 ug/mL

Reagent

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**VOA8260GAS1ST\_00091**



# CERTIFIED REFERENCE MATERIAL

110 Benner Circle  
Bellefonte, PA 16823-8812  
Tel: (800)356-1688  
Fax: (814)353-1309

[www.restek.com](http://www.restek.com)

## Certificate of Analysis



### FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

*This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.*

**Catalog No. :** 569722

**Lot No.:** A0108198

**Description :** 8260 List 1 / Std #3 Gases (2015)

8260 List 1 / Std #3 Gases (2015) 2,000 ug/ml, P&T Methanol, 1 ml/ampul

**Container Size :** 2 mL

**Pkg Amt:** > 1 mL

**Expiration Date :** January 31, 2018

**Storage:** 0°C or colder

### C E R T I F I E D   V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Dichlorodifluoromethane (CFC-12)	2,504.8 µg/mL	+/- 21.9788	µg/mL	Gravimetric
	CAS # 75-71-8	(Lot Q167-08)	+/- 32.6918	µg/mL	Unstressed
	Purity 99%		+/- 36.4326	µg/mL	Stressed
2	Chloromethane (methyl chloride)	2,509.8 µg/mL	+/- 19.6377	µg/mL	Gravimetric
	CAS # 74-87-3	(Lot SHBC8470V)	+/- 31.2039	µg/mL	Unstressed
	Purity 99%		+/- 35.1185	µg/mL	Stressed
3	Vinyl chloride	2,515.3 µg/mL	+/- 22.1368	µg/mL	Gravimetric
	CAS # 75-01-4	(Lot 17542)	+/- 32.8734	µg/mL	Unstressed
	Purity 99%		+/- 36.6254	µg/mL	Stressed
4	1,3-Butadiene	2,498.0 µg/mL	+/- 23.6713	µg/mL	Gravimetric
	CAS # 106-99-0	(Lot SHBD5808V)	+/- 33.8065	µg/mL	Unstressed
	Purity 99%		+/- 37.4176	µg/mL	Stressed
5	Bromomethane (methyl bromide)	2,503.7 µg/mL	+/- 30.8470	µg/mL	Gravimetric
	CAS # 74-83-9	(Lot 101604)	+/- 39.2011	µg/mL	Unstressed
	Purity 99%		+/- 42.3685	µg/mL	Stressed
6	Chloroethane (ethyl chloride)	2,507.7 µg/mL	+/- 21.9404	µg/mL	Gravimetric
	CAS # 75-00-3	(Lot SHBD1717V)	+/- 32.6873	µg/mL	Unstressed
	Purity 99%		+/- 36.4370	µg/mL	Stressed
7	Dichlorodifluoromethane (CFC-21)	2,500.7 µg/mL	+/- 26.0039	µg/mL	Gravimetric
	CAS # 75-43-4	(Lot Q9B-58)	+/- 35.4965	µg/mL	Unstressed
	Purity 99%		+/- 38.9583	µg/mL	Stressed

8	Trichlorofluoromethane (CFC-11)	2,501.9 $\mu\text{g/mL}$	+/- 21.5914 $\mu\text{g/mL}$	Gravimetric
CAS #	75-69-4	(Lot SHBD5121V)	+/- 32.4119 $\mu\text{g/mL}$	Unstressed
Purity	99%		+/- 36.1734 $\mu\text{g/mL}$	Stressed

**Solvent:** P&T Methanol  
**CAS #** 67-56-1  
**Purity** 99%

**Column:**  
60m x 0.25mm x 1.4 $\mu\text{m}$   
Rtx-502.2 (cat.#10916)

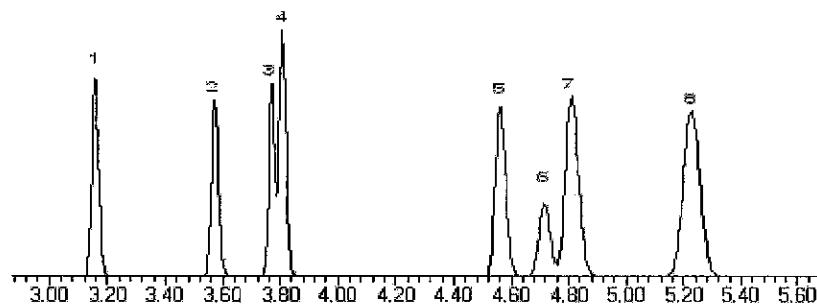
**Carrier Gas:**  
helium-constant flow 2.0 mL/min.

**Temp. Program:**  
40°C (hold 6 min.) to 100°C  
@ 6°C/min.

**Inj. Temp:**  
200°C

**Det. Temp:**  
250°C

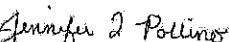
**Det. Type:**  
MSD



This chromatogram represents a general set of testing conditions chosen for product acceptance. For optimal results in your lab, conditions should be adjusted for your specific instrument, method, and application.

  
Kendra Swope - Mix Technician

Date Mixed: 08-Jan-2015 Balance: 1125113331

  
Jennifer L. Pollino - QC Analyst

Date Passed: 14-Jan-2015

Manufactured under Restek's ISO 9001:2008 Registered Quality System Certificate #FM 80397
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Reagent

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**VOA8260GAS1ST\_00095**



# CERTIFIED REFERENCE MATERIAL

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## Certificate of Analysis



### FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

*This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.*

**Catalog No. :** 569722

**Lot No.:** A0108198

**Description :** 8260 List 1 / Std #3 Gases (2015)

8260 List 1 / Std #3 Gases (2015) 2,000 ug/ml, P&T Methanol, 1 ml/ampul

**Container Size :** 2 mL

**Pkg Amt:** > 1 mL

**Expiration Date :** January 31, 2018

**Storage:** 0°C or colder

### C E R T I F I E D V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Dichlorodifluoromethane (CFC-12)	2,504.8 µg/mL	+/- 21.9788	µg/mL	Gravimetric
	CAS # 75-71-8	(Lot Q167-08)	+/- 32.6918	µg/mL	Unstressed
	Purity 99%		+/- 36.4326	µg/mL	Stressed
2	Chloromethane (methyl chloride)	2,509.8 µg/mL	+/- 19.6377	µg/mL	Gravimetric
	CAS # 74-87-3	(Lot SHBC8470V)	+/- 31.2039	µg/mL	Unstressed
	Purity 99%		+/- 35.1185	µg/mL	Stressed
3	Vinyl chloride	2,515.3 µg/mL	+/- 22.1368	µg/mL	Gravimetric
	CAS # 75-01-4	(Lot 17542)	+/- 32.8734	µg/mL	Unstressed
	Purity 99%		+/- 36.6254	µg/mL	Stressed
4	1,3-Butadiene	2,498.0 µg/mL	+/- 23.6713	µg/mL	Gravimetric
	CAS # 106-99-0	(Lot SHBD5808V)	+/- 33.8065	µg/mL	Unstressed
	Purity 99%		+/- 37.4176	µg/mL	Stressed
5	Bromomethane (methyl bromide)	2,503.7 µg/mL	+/- 30.8470	µg/mL	Gravimetric
	CAS # 74-83-9	(Lot 101604)	+/- 39.2011	µg/mL	Unstressed
	Purity 99%		+/- 42.3685	µg/mL	Stressed
6	Chloroethane (ethyl chloride)	2,507.7 µg/mL	+/- 21.9404	µg/mL	Gravimetric
	CAS # 75-00-3	(Lot SHBD1717V)	+/- 32.6873	µg/mL	Unstressed
	Purity 99%		+/- 36.4370	µg/mL	Stressed
7	Dichlorodifluoromethane (CFC-21)	2,500.7 µg/mL	+/- 26.0039	µg/mL	Gravimetric
	CAS # 75-43-4	(Lot Q9B-58)	+/- 35.4965	µg/mL	Unstressed
	Purity 99%		+/- 38.9583	µg/mL	Stressed

8	Trichlorofluoromethane (CFC-11)	2,501.9 $\mu\text{g/mL}$	+/- 21.5914 $\mu\text{g/mL}$	Gravimetric
CAS #	75-69-4	(Lot SHBD5121V)	+/- 32.4119 $\mu\text{g/mL}$	Unstressed
Purity	99%		+/- 36.1734 $\mu\text{g/mL}$	Stressed

**Solvent:** P&T Methanol  
**CAS #** 67-56-1  
**Purity** 99%

**Column:**  
60m x 0.25mm x 1.4 $\mu\text{m}$   
Rtx-502.2 (cat.#10916)

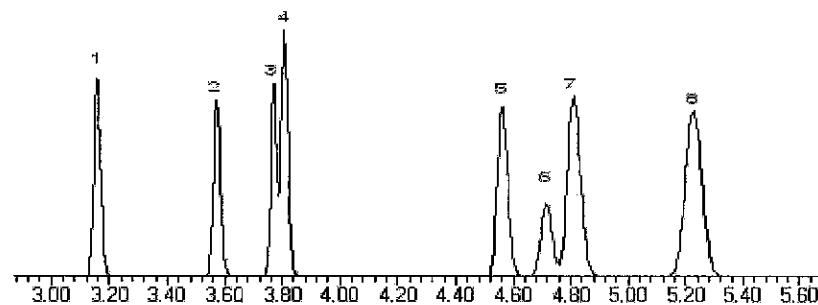
**Carrier Gas:**  
helium-constant flow 2.0 mL/min.

**Temp. Program:**  
40°C (hold 6 min.) to 100°C  
@ 6°C/min.

**Inj. Temp:**  
200°C

**Det. Temp:**  
250°C

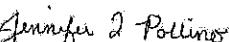
**Det. Type:**  
MSD



This chromatogram represents a general set of testing conditions chosen for product acceptance. For optimal results in your lab, conditions should be adjusted for your specific instrument, method, and application.

  
Kendra Swope - Mix Technician

Date Mixed: 08-Jan-2015 Balance: 1125113331

  
Jennifer L. Pollino - QC Analyst

Date Passed: 14-Jan-2015

Manufactured under Restek's ISO 9001:2008 Registered Quality System Certificate #FM 80397
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Reagent

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**VOA8260GAS2ND\_00093**



# CERTIFIED REFERENCE MATERIAL

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## Certificate of Analysis



### FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

*This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.*

**Catalog No. :** 569722.sec

**Lot No.:** A0108226

**Description :** 8260 List 1 / Std #3 Gases (2015)

8260 List 1 / Std #3 Gases (2015) 2,000 ug/ml, P&T Methanol, 1 ml/ampul

**Container Size :** 2 mL

**Pkg Amt:** > 1 mL

**Expiration Date :** January 31, 2018

**Storage:** 0°C or colder

### C E R T I F I E D V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)	Gravimetric
	Dichlorodifluoromethane (CFC-12) <b>CAS #</b> 75-71-8.SEC <b>Purity</b> 99%	2,494.8 µg/mL	+/- 23.5521 µg/mL	
			+/- 33.7009 µg/mL	Unstressed
			+/- 37.3133 µg/mL	Stressed
2	Chloromethane (methyl chloride) <b>CAS #</b> 74-87-3.SEC <b>Purity</b> 99%	2,505.6 µg/mL	+/- 26.4745 µg/mL	Gravimetric
			+/- 35.8743 µg/mL	Unstressed
			+/- 39.3156 µg/mL	Stressed
3	Vinyl chloride <b>CAS #</b> 75-01-4.SEC <b>Purity</b> 99%	2,499.8 µg/mL	+/- 25.3054 µg/mL	Gravimetric
			+/- 34.9816 µg/mL	Unstressed
			+/- 38.4872 µg/mL	Stressed
4	1,3-Butadiene <b>CAS #</b> 106-99-0.SEC <b>Purity</b> 99%	2,505.4 µg/mL	+/- 23.1450 µg/mL	Gravimetric
			+/- 33.4914 µg/mL	Unstressed
			+/- 37.1536 µg/mL	Stressed
5	Bromomethane (methyl bromide) <b>CAS #</b> 74-83-9.SEC <b>Purity</b> 99%	2,495.4 µg/mL	+/- 25.3762 µg/mL	Gravimetric
			+/- 35.0038 µg/mL	Unstressed
			+/- 38.4957 µg/mL	Stressed
6	Chloroethane (ethyl chloride) <b>CAS #</b> 75-00-3.SEC <b>Purity</b> 99%	2,499.5 µg/mL	+/- 21.8687 µg/mL	Gravimetric
			+/- 32.5806 µg/mL	Unstressed
			+/- 36.3180 µg/mL	Stressed
7	Dichlorofluoromethane (CFC-21) <b>CAS #</b> 75-43-4.SEC <b>Purity</b> 99%	2,511.0 µg/mL	+/- 21.9690 µg/mL	Gravimetric
			+/- 32.7299 µg/mL	Unstressed
			+/- 36.4846 µg/mL	Stressed

8	Trichlorofluoromethane (CFC-11)	2,504.4	µg/mL	+/- 25.2390	µg/mL	Gravimetric
	CAS # 75-69-4 SEC	(Lot Q158-102)		+/- 34.9647	µg/mL	Unstressed
	Purity 99%			+/- 38.4843	µg/mL	Stressed

**Solvent:** P&T Methanol  
**CAS #** 67-56-1  
**Purity** 99%

**Column:**  
60m x 0.25mm x 1.4µm  
Rtx-502.2 (cat.#10916)

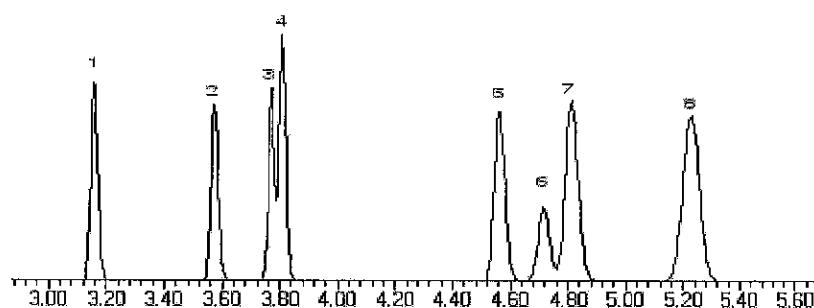
**Carrier Gas:**  
helium-constant flow 2.0 ml./min.

**Temp. Program:**  
40°C (hold 6 min.) to 100°C  
@ 6°C/min.

**Inj. Temp:**  
200°C

**Det. Temp:**  
250°C

**Det. Type:**  
MSD



This chromatogram represents a general set of testing conditions chosen for product acceptance. For optimal results in your lab, conditions should be adjusted for your specific instrument, method, and application.

*Michael J. Maye*

Date Mixed: 12-Jan-2015 Balance: 1127510105

*Jennifer L. Pollino*

Jennifer L. Pollino - QC Analyst

Date Passed: 14-Jan-2015

Manufactured under Restek's ISO 9001:2008 Registered Quality System Certificate #FM 80397
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Reagent

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**VOA8260INTRES\_00091**



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## Certificate of Analysis



### FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

*This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.*

**Catalog No. :** 567649

**Lot No.:** A0104742

**Description :** 8260 Internal Standard

8260 Internal Standard 250-5,000 ug/ml, P&T Methanol, 5 mL/ampul

**Container Size :** 5 mL

**Pkg Amt:** > 5 mL

**Expiration Date :** July 31, 2019

**Storage:** 0°C or colder

### C E R T I F I E D V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	tert-Butyl-d9-alcohol <b>CAS #</b> 25725-11-5 <b>Purity</b> 99%	5,003.0 µg/mL (Lot I201P5)	+/- 29.0879 µg/mL	+/- 106.1005 µg/mL	Gravimetric Unstressed
			+/- 106.5713 µg/mL	+/- 106.5713 µg/mL	Stressed
2	Fluorobenzene <b>CAS #</b> 462-06-6 <b>Purity</b> 99%	250.8 µg/mL (Lot 1380033)	+/- 1.4795 µg/mL	+/- 5.3247 µg/mL	Gravimetric Unstressed
			+/- 5.3483 µg/mL	+/- 5.3483 µg/mL	Stressed
3	1,4-Dioxane-d8 <b>CAS #</b> 17647-74-4 <b>Purity</b> 99%	5,009.6 µg/mL (Lot 11C-596)	+/- 29.1262 µg/mL	+/- 106.2405 µg/mL	Gravimetric Unstressed
			+/- 106.7119 µg/mL	+/- 106.7119 µg/mL	Stressed
4	Chlorobenzene-d5 <b>CAS #</b> 3114-55-4 <b>Purity</b> 99%	250.8 µg/mL (Lot PR-22736)	+/- 1.4795 µg/mL	+/- 5.3247 µg/mL	Gravimetric Unstressed
			+/- 5.3483 µg/mL	+/- 5.3483 µg/mL	Stressed
5	1,4-Dichlorobenzene-d4 <b>CAS #</b> 3855-82-1 <b>Purity</b> 99%	250.8 µg/mL (Lot PR-18488)	+/- 1.4795 µg/mL	+/- 5.3247 µg/mL	Gravimetric Unstressed
			+/- 5.3483 µg/mL	+/- 5.3483 µg/mL	Stressed

**Solvent:** P&T Methanol  
**CAS #** 67-56-1  
**Purity** 99%

Reagent

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**VOA8260KET1ST\_00036**



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## Certificate of Analysis

### FOR LABORATORY USE ONLY-READ MSDS PRIOR TO USE.

*This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.*

**Catalog No. :** 567642

**Lot No.:** A093365

**Description :** 8260 List 1 / Std #2 Ketones

8260 List 1 / Std #2 Ketones 10,000 ug/ml, P&T Methanol/Water (90:10),  
1 mL/ampul

**Container Size :** 2 mL

**Pkg Amt:** > 1 mL

**Expiration Date :** February 2016

**Storage:** 0°C or colder

### C E R T I F I E D V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Acetone	10,000.0 µg/mL	+/- 58.1378	µg/mL	Gravimetric
	CAS # 67-64-1		+/- 798.6896	µg/mL	Unstressed
	Purity 99%		+/- 799.0807	µg/mL	Stressed
2	2-Butanone (MEK)	10,000.0 µg/mL	+/- 58.1378	µg/mL	Gravimetric
	CAS # 78-93-3		+/- 798.6896	µg/mL	Unstressed
	Purity 99%		+/- 799.0807	µg/mL	Stressed
3	4-Methyl-2-pentanone (MIBK)	10,000.0 µg/mL	+/- 58.1378	µg/mL	Gravimetric
	CAS # 108-10-1		+/- 798.6896	µg/mL	Unstressed
	Purity 99%		+/- 799.0807	µg/mL	Stressed
4	2-Hexanone	10,000.0 µg/mL	+/- 58.1378	µg/mL	Gravimetric
	CAS # 591-78-6		+/- 798.6896	µg/mL	Unstressed
	Purity 99%		+/- 799.0807	µg/mL	Stressed

**Solvent:** P&T Methanol/Water (90:10)  
CAS # 67-56-1/7732-18-5  
Purity 99%

Reagent

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**VOA8260KET1ST\_00037**



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## Certificate of Analysis

### FOR LABORATORY USE ONLY-READ MSDS PRIOR TO USE.

*This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.*

**Catalog No. :** 567642

**Lot No.:** A093365

**Description :** 8260 List 1 / Std #2 Ketones

8260 List 1 / Std #2 Ketones 10,000 ug/ml, P&T Methanol/Water (90:10),  
1 mL/ampul

**Container Size :** 2 mL

**Pkg Amt:** > 1 mL

**Expiration Date :** February 2016

**Storage:** 0°C or colder

### C E R T I F I E D V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Acetone	10,000.0 µg/mL	+/- 58.1378	µg/mL	Gravimetric
	CAS # 67-64-1		+/- 798.6896	µg/mL	Unstressed
	Purity 99%		+/- 799.0807	µg/mL	Stressed
2	2-Butanone (MEK)	10,000.0 µg/mL	+/- 58.1378	µg/mL	Gravimetric
	CAS # 78-93-3		+/- 798.6896	µg/mL	Unstressed
	Purity 99%		+/- 799.0807	µg/mL	Stressed
3	4-Methyl-2-pentanone (MIBK)	10,000.0 µg/mL	+/- 58.1378	µg/mL	Gravimetric
	CAS # 108-10-1		+/- 798.6896	µg/mL	Unstressed
	Purity 99%		+/- 799.0807	µg/mL	Stressed
4	2-Hexanone	10,000.0 µg/mL	+/- 58.1378	µg/mL	Gravimetric
	CAS # 591-78-6		+/- 798.6896	µg/mL	Unstressed
	Purity 99%		+/- 799.0807	µg/mL	Stressed

**Solvent:** P&T Methanol/Water (90:10)  
**CAS #** 67-56-1/7732-18-5  
**Purity** 99%

Reagent

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**VOA8260KET1ST\_00039**



# CERTIFIED REFERENCE MATERIAL

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## Certificate of Analysis



### FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

*This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.*

**Catalog No.:** 569721

**Lot No.:** A0108151

**Description :** 8260 List 1/ Std #2 Ketones (2015)

8260 List 1/ Std #2 Ketones (2015) 12,500 µg/mL, P&T Methanol/Water (90:10), 1 mL/ampul

**Container Size :** 2 mL

**Pkg Amt:** > 1 mL

**Expiration Date :** January 31, 2018

**Storage:** 0°C or colder

### C E R T I F I E D   V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Acetone <b>CAS #</b> 67-64-1 <b>Purity</b> 99%	12,537.0 µg/mL (Lot 07196AK)	+/- 73.4069	µg/mL	Gravimetric
			+/- 667.2480	µg/mL	Unstressed
			+/- 667.9837	µg/mL	Stressed
2	2-Butanone (MEK) <b>CAS #</b> 78-93-3 <b>Purity</b> 99%	12,537.0 µg/mL (Lot BCBH7802V)	+/- 73.4069	µg/mL	Gravimetric
			+/- 667.2480	µg/mL	Unstressed
			+/- 667.9837	µg/mL	Stressed
3	4-Methyl-2-pentanone (MIBK) <b>CAS #</b> 108-10-1 <b>Purity</b> 99%	12,537.0 µg/mL (Lot SHBF5332V)	+/- 73.4069	µg/mL	Gravimetric
			+/- 667.2480	µg/mL	Unstressed
			+/- 667.9837	µg/mL	Stressed
4	2-Hexanone <b>CAS #</b> 591-78-6 <b>Purity</b> 99%	12,537.0 µg/mL (Lot MKBK8325V)	+/- 73.4069	µg/mL	Gravimetric
			+/- 667.2480	µg/mL	Unstressed
			+/- 667.9837	µg/mL	Stressed

**Solvent:** P&T Methanol/Water (90:10)  
**CAS #** 67-56-1/7732-18-5  
**Purity** 99%

Reagent

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**VOA8260KET2ND\_00042**



# CERTIFIED REFERENCE MATERIAL

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## Certificate of Analysis



### FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

*This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.*

Catalog No. :	<u>569721.SEC</u>	Lot No.:	<u>A0108157</u>
Description :	8260 List 1/ Std #2 Ketones (2015)		
8260 List 1/ Std #2 Ketones (2015) 12,500 µg/mL, P&T Methanol/Water (90:10), 1 mL/ampul			
Container Size :	<u>2 mL</u>	Pkg Amt:	<u>&gt; 1 mL</u>
Expiration Date :	<u>January 31, 2018</u>		
		Storage:	<u>0°C or colder</u>

### C E R T I F I E D V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Acetone	12,504.0 µg/mL	+/-	73.2137	µg/mL
	CAS # 67-64-1.SEC	(Lot 0902033)	+/-	665.4917	µg/mL
	Purity 99%		+/-	666.2255	µg/mL
2	2-Butanone (MEK)	12,506.0 µg/mL	+/-	73.2254	µg/mL
	CAS # 78-93-3.SEC	(Lot VEGGI)	+/-	665.5981	µg/mL
	Purity 99%		+/-	666.3320	µg/mL
3	4-Methyl-2-pentanone (MIBK)	12,537.3 µg/mL	+/-	73.4088	µg/mL
	CAS # 108-10-1.SEC	(Lot E29T040)	+/-	667.2658	µg/mL
	Purity 99%		+/-	668.0015	µg/mL
4	2-Hexanone	12,508.7 µg/mL	+/-	73.2410	µg/mL
	CAS # 591-78-6.SEC	(Lot ZSVCD-FF)	+/-	665.7401	µg/mL
	Purity 99%		+/-	666.4741	µg/mL

Solvent: P&T Methanol/Water (90:10)  
 CAS # 67-56-1/7732-18-5  
 Purity 99%

Reagent

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**VOA8260MEGA1\_00014**



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## Certificate of Analysis

### FOR LABORATORY USE ONLY-READ MSDS PRIOR TO USE.

*This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.*

**Catalog No. :** 567641

**Lot No.:** A093581

**Description :** 8260 List 1 / Std #1 MegaMix

8260 List 1 / Std #1 MegaMix 1000-50,000 µg/mL, P&T Methanol, 1 ml/ampul

**Container Size :** 2 mL

**Pkg Amt:** > 1 mL

**Expiration Date :** February 2016

**Storage:** 0°C or colder

### C E R T I F I E D V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Diethyl ether (ethyl ether) <b>CAS #</b> 60-29-7 <b>Purity</b> 99%	2,000.0 µg/mL	+/- 11.6282	µg/mL	Gravimetric
			+/- 44.2531	µg/mL	Unstressed
			+/- 44.4335	µg/mL	Stressed
2	1,1,2-Trichlorotrifluoroethane (CFC-113) <b>CAS #</b> 76-13-1 <b>Purity</b> 97%	1,999.9 µg/mL	+/- 11.6279	µg/mL	Gravimetric
			+/- 44.2519	µg/mL	Unstressed
			+/- 44.4323	µg/mL	Stressed
3	1,1-dichloroethene <b>CAS #</b> 75-35-4 <b>Purity</b> 98%	2,000.0 µg/mL	+/- 11.6281	µg/mL	Gravimetric
			+/- 44.2527	µg/mL	Unstressed
			+/- 44.4331	µg/mL	Stressed
4	tert-Butanol (TBA) <b>CAS #</b> 75-65-0 <b>Purity</b> 99%	20,000.0 µg/mL	+/- 116.2756	µg/mL	Gravimetric
			+/- 442.5291	µg/mL	Unstressed
			+/- 444.3332	µg/mL	Stressed
5	Iodomethane (methyl iodide) <b>CAS #</b> 74-88-4 <b>Purity</b> 99%	2,000.0 µg/mL	+/- 11.6282	µg/mL	Gravimetric
			+/- 44.2531	µg/mL	Unstressed
			+/- 44.4335	µg/mL	Stressed
6	Allyl chloride ( 3-chloropropene ) <b>CAS #</b> 107-05-1 <b>Purity</b> 98%	2,000.0 µg/mL	+/- 11.6281	µg/mL	Gravimetric
			+/- 44.2527	µg/mL	Unstressed
			+/- 44.4331	µg/mL	Stressed
7	Methyl acetate <b>CAS #</b> 79-20-9 <b>Purity</b> 99%	10,000.0 µg/mL	+/- 58.1378	µg/mL	Gravimetric
			+/- 221.2646	µg/mL	Unstressed
			+/- 222.1666	µg/mL	Stressed
8	Carbon disulfide <b>CAS #</b> 75-15-0 <b>Purity</b> 98%	2,000.0 µg/mL	+/- 11.6281	µg/mL	Gravimetric
			+/- 44.2527	µg/mL	Unstressed
			+/- 44.4331	µg/mL	Stressed
9	Methylene chloride (dichloromethane) <b>CAS #</b> 75-09-2 <b>Purity</b> 99%	2,000.0 µg/mL	+/- 11.6282	µg/mL	Gravimetric
			+/- 44.2531	µg/mL	Unstressed
			+/- 44.4335	µg/mL	Stressed

10	Acrylonitrile <b>CAS #</b> 107-13-1 <b>Purity</b> 99%	20,000.0	µg/mL	+/- 116.2756 +/- 442.5291 +/- 444.3332	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
11	Methyl-tert-butyl ether ( MTBE ) <b>CAS #</b> 1634-04-4 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
12	cis-1,2-Dichloroethene <b>CAS #</b> 156-59-2 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
13	n-Hexane (C6) <b>CAS #</b> 110-54-3 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
14	1,1-Dichloroethane <b>CAS #</b> 75-34-3 <b>Purity</b> 98%	2,000.0	µg/mL	+/- 11.6281 +/- 44.2527 +/- 44.4331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
15	2,2-Dichloropropane <b>CAS #</b> 594-20-7 <b>Purity</b> 98%	2,000.0	µg/mL	+/- 11.6281 +/- 44.2527 +/- 44.4331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
16	trans-1,2-Dichloroethene <b>CAS #</b> 156-60-5 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
17	chloroform <b>CAS #</b> 67-66-3 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
18	Isobutanol (2-Methyl-1-propanol) <b>CAS #</b> 78-83-1 <b>Purity</b> 99%	50,000.0	µg/mL	+/- 290.6891 +/- 1,106.3228 +/- 1,110.8331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
19	Bromochloromethane <b>CAS #</b> 74-97-5 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
20	Tetrahydrofuran <b>CAS #</b> 109-99-9 <b>Purity</b> 99%	4,000.0	µg/mL	+/- 23.2563 +/- 88.5061 +/- 88.8670	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
21	1,1,1-trichloroethane <b>CAS #</b> 71-55-6 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
22	Cyclohexane <b>CAS #</b> 110-82-7 <b>Purity</b> 98%	2,000.0	µg/mL	+/- 11.6281 +/- 44.2527 +/- 44.4331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
23	1,1-Dichloropropene <b>CAS #</b> 563-58-6 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
24	carbon tetrachloride <b>CAS #</b> 56-23-5 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
25	n-Heptane (C7) <b>CAS #</b> 142-82-5 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
26	Benzene <b>CAS #</b> 71-43-2 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
27	1,2-Dichloroethane <b>CAS #</b> 107-06-2 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
28	Trichloroethylene <b>CAS #</b> 79-01-6 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed

29	Methylcyclohexane CAS # 108-87-2 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
30	1,2-Dichloropropane CAS # 78-87-5 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
31	1,4-Dioxane CAS # 123-91-1 Purity 99%	40,000.0 µg/mL	+/- 232.5513 +/- 885.0582 +/- 888.6665	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
32	Dibromomethane CAS # 74-95-3 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
33	bromodichloromethane CAS # 75-27-4 Purity 97%	2,000.0 µg/mL	+/- 11.6284 +/- 44.2540 +/- 44.4344	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
34	cis-1,3-Dichloropropene CAS # 10061-01-5 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
35	Toluene CAS # 108-88-3 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
36	Ethyl methacrylate CAS # 97-63-2 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
37	trans-1,3-Dichloropropene CAS # 10061-02-6 Purity 97%	2,000.0 µg/mL	+/- 11.6284 +/- 44.2540 +/- 44.4344	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
38	1,1,2-Trichloroethane CAS # 79-00-5 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
39	1,3-Dichloropropane CAS # 142-28-9 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
40	Tetrachloroethene CAS # 127-18-4 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
41	dibromochloromethane CAS # 124-48-1 Purity 98%	2,000.0 µg/mL	+/- 11.6281 +/- 44.2527 +/- 44.4331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
42	1,2-Dibromoethane (EDB) CAS # 106-93-4 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
43	Chlorobenzene CAS # 108-90-7 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
44	1,1,1,2-Tetrachloroethane CAS # 630-20-6 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
45	m-Xylene CAS # 108-38-3 Purity 99%	1,000.0 µg/mL	+/- 5.8141 +/- 22.1265 +/- 22.2167	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
46	p-Xylene CAS # 106-42-3 Purity 99%	1,000.0 µg/mL	+/- 5.8141 +/- 22.1265 +/- 22.2167	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
47	o-Xylene CAS # 95-47-6 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed

48	Ethylbenzene <b>CAS #</b> 100-41-4 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
49	Styrene <b>CAS #</b> 100-42-5 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
50	Isopropylbenzene (cumene) <b>CAS #</b> 98-82-8 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
51	bromoform <b>CAS #</b> 75-25-2 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
52	1,1,2,2-Tetrachloroethane <b>CAS #</b> 79-34-5 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
53	1,2,3-Trichloropropane <b>CAS #</b> 96-18-4 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
54	trans-1,4-dichloro-2-butene <b>CAS #</b> 110-57-6 <b>Purity</b> 98%	2,000.0	µg/mL	+/- 11.6281 +/- 44.2527 +/- 44.4331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
55	n-Propylbenzene <b>CAS #</b> 103-65-1 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
56	Bromobenzene <b>CAS #</b> 108-86-1 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
57	1,3,5-Trimethylbenzene <b>CAS #</b> 108-67-8 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
58	2-Chlorotoluene <b>CAS #</b> 95-49-8 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
59	4-Chlorotoluene <b>CAS #</b> 106-43-4 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
60	tert-Butylbenzene <b>CAS #</b> 98-06-6 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
61	1,2,4-Trimethylbenzene <b>CAS #</b> 95-63-6 <b>Purity</b> 98%	2,000.0	µg/mL	+/- 11.6281 +/- 44.2527 +/- 44.4331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
62	sec-Butylbenzene <b>CAS #</b> 135-98-8 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
63	4-Isopropyltoluene (p-Cymene) <b>CAS #</b> 99-87-6 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
64	1,3-Dichlorobenzene <b>CAS #</b> 541-73-1 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
65	1,4-Dichlorobenzene <b>CAS #</b> 106-46-7 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
66	n-Butylbenzene <b>CAS #</b> 104-51-8 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed

67	1,2-Dichlorobenzene CAS # 95-50-1 Purity 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
68	1,2-Dibromo-3-chloropropane CAS # 96-12-8 Purity 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
69	1,2,4-Trichlorobenzene CAS # 120-82-1 Purity 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
70	Hexachlorobutadiene CAS # 87-68-3 Purity 97%	2,000.0	µg/mL	+/- 11.6284 +/- 44.2540 +/- 44.4344	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
71	Naphthalene CAS # 91-20-3 Purity 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
72	1,2,3-Trichlorobenzene CAS # 87-61-6 Purity 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed

**Solvent:** P&T Methanol  
**CAS #** 67-56-1  
**Purity** 99%

**Column:**  
 60m x .25mm x 1.4µm  
 Rtx-502.2 (cat.#10916)

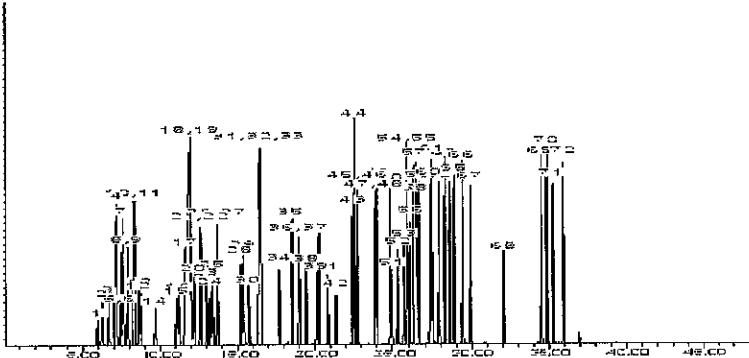
**Carrier Gas:**  
 helium-constant pressure 30 psi

**Temp. Program:**  
 40°C (hold 6 min.) to 240°C  
 @ 6°C/min. (hold 10 min.)

**Inj. Temp:**  
 200°C

**Det. Temp:**  
 250°C

**Det. Type:**  
 MSD



*Jennifer L. Pollino*  
 Jennifer L. Pollino - QC Analyst

Date Passed: 01-Mar-2013 Balance: B251644995

Manufactured under Restek's ISO 9001:2008  
 Registered Quality System  
 Certificate #FM 80397

Reagent

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**VOA8260MEGA1\_00027**



110 Benner Circle  
Bellefonte, PA 16823-8812  
Tel: (800)356-1688  
Fax: (814)353-1309

[www.restek.com](http://www.restek.com)



## Certificate of Analysis

### FOR LABORATORY USE ONLY-READ MSDS PRIOR TO USE.

*This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.*

**Catalog No. :** 567641

**Lot No.:** A093581

**Description :** 8260 List 1 / Std #1 MegaMix

8260 List 1 / Std #1 MegaMix 1000-50,000 µg/mL, P&T Methanol, 1 ml/ampul

**Container Size :** 2 mL

**Pkg Amt:** > 1 mL

**Expiration Date :** February 2016

**Storage:** 0°C or colder

### C E R T I F I E D V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Diethyl ether (ethyl ether) <b>CAS #</b> 60-29-7 <b>Purity</b> 99%	2,000.0 µg/mL	+/- 11.6282	µg/mL	Gravimetric
			+/- 44.2531	µg/mL	Unstressed
			+/- 44.4335	µg/mL	Stressed
2	1,1,2-Trichlorotrifluoroethane (CFC-113) <b>CAS #</b> 76-13-1 <b>Purity</b> 97%	1,999.9 µg/mL	+/- 11.6279	µg/mL	Gravimetric
			+/- 44.2519	µg/mL	Unstressed
			+/- 44.4323	µg/mL	Stressed
3	1,1-dichloroethene <b>CAS #</b> 75-35-4 <b>Purity</b> 98%	2,000.0 µg/mL	+/- 11.6281	µg/mL	Gravimetric
			+/- 44.2527	µg/mL	Unstressed
			+/- 44.4331	µg/mL	Stressed
4	tert-Butanol (TBA) <b>CAS #</b> 75-65-0 <b>Purity</b> 99%	20,000.0 µg/mL	+/- 116.2756	µg/mL	Gravimetric
			+/- 442.5291	µg/mL	Unstressed
			+/- 444.3332	µg/mL	Stressed
5	Iodomethane (methyl iodide) <b>CAS #</b> 74-88-4 <b>Purity</b> 99%	2,000.0 µg/mL	+/- 11.6282	µg/mL	Gravimetric
			+/- 44.2531	µg/mL	Unstressed
			+/- 44.4335	µg/mL	Stressed
6	Allyl chloride ( 3-chloropropene ) <b>CAS #</b> 107-05-1 <b>Purity</b> 98%	2,000.0 µg/mL	+/- 11.6281	µg/mL	Gravimetric
			+/- 44.2527	µg/mL	Unstressed
			+/- 44.4331	µg/mL	Stressed
7	Methyl acetate <b>CAS #</b> 79-20-9 <b>Purity</b> 99%	10,000.0 µg/mL	+/- 58.1378	µg/mL	Gravimetric
			+/- 221.2646	µg/mL	Unstressed
			+/- 222.1666	µg/mL	Stressed
8	Carbon disulfide <b>CAS #</b> 75-15-0 <b>Purity</b> 98%	2,000.0 µg/mL	+/- 11.6281	µg/mL	Gravimetric
			+/- 44.2527	µg/mL	Unstressed
			+/- 44.4331	µg/mL	Stressed
9	Methylene chloride (dichloromethane) <b>CAS #</b> 75-09-2 <b>Purity</b> 99%	2,000.0 µg/mL	+/- 11.6282	µg/mL	Gravimetric
			+/- 44.2531	µg/mL	Unstressed
			+/- 44.4335	µg/mL	Stressed

10	Acrylonitrile <b>CAS #</b> 107-13-1 <b>Purity</b> 99%	20,000.0	µg/mL	+/- 116.2756 +/- 442.5291 +/- 444.3332	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
11	Methyl-tert-butyl ether ( MTBE ) <b>CAS #</b> 1634-04-4 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
12	cis-1,2-Dichloroethene <b>CAS #</b> 156-59-2 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
13	n-Hexane (C6) <b>CAS #</b> 110-54-3 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
14	1,1-Dichloroethane <b>CAS #</b> 75-34-3 <b>Purity</b> 98%	2,000.0	µg/mL	+/- 11.6281 +/- 44.2527 +/- 44.4331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
15	2,2-Dichloropropane <b>CAS #</b> 594-20-7 <b>Purity</b> 98%	2,000.0	µg/mL	+/- 11.6281 +/- 44.2527 +/- 44.4331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
16	trans-1,2-Dichloroethene <b>CAS #</b> 156-60-5 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
17	chloroform <b>CAS #</b> 67-66-3 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
18	Isobutanol (2-Methyl-1-propanol) <b>CAS #</b> 78-83-1 <b>Purity</b> 99%	50,000.0	µg/mL	+/- 290.6891 +/- 1,106.3228 +/- 1,110.8331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
19	Bromochloromethane <b>CAS #</b> 74-97-5 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
20	Tetrahydrofuran <b>CAS #</b> 109-99-9 <b>Purity</b> 99%	4,000.0	µg/mL	+/- 23.2563 +/- 88.5061 +/- 88.8670	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
21	1,1,1-trichloroethane <b>CAS #</b> 71-55-6 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
22	Cyclohexane <b>CAS #</b> 110-82-7 <b>Purity</b> 98%	2,000.0	µg/mL	+/- 11.6281 +/- 44.2527 +/- 44.4331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
23	1,1-Dichloropropene <b>CAS #</b> 563-58-6 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
24	carbon tetrachloride <b>CAS #</b> 56-23-5 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
25	n-Heptane (C7) <b>CAS #</b> 142-82-5 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
26	Benzene <b>CAS #</b> 71-43-2 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
27	1,2-Dichloroethane <b>CAS #</b> 107-06-2 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
28	Trichloroethylene <b>CAS #</b> 79-01-6 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed

29	Methylcyclohexane CAS # 108-87-2 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
30	1,2-Dichloropropane CAS # 78-87-5 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
31	1,4-Dioxane CAS # 123-91-1 Purity 99%	40,000.0 µg/mL	+/- 232.5513 +/- 885.0582 +/- 888.6665	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
32	Dibromomethane CAS # 74-95-3 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
33	bromodichloromethane CAS # 75-27-4 Purity 97%	2,000.0 µg/mL	+/- 11.6284 +/- 44.2540 +/- 44.4344	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
34	cis-1,3-Dichloropropene CAS # 10061-01-5 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
35	Toluene CAS # 108-88-3 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
36	Ethyl methacrylate CAS # 97-63-2 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
37	trans-1,3-Dichloropropene CAS # 10061-02-6 Purity 97%	2,000.0 µg/mL	+/- 11.6284 +/- 44.2540 +/- 44.4344	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
38	1,1,2-Trichloroethane CAS # 79-00-5 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
39	1,3-Dichloropropane CAS # 142-28-9 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
40	Tetrachloroethene CAS # 127-18-4 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
41	dibromochloromethane CAS # 124-48-1 Purity 98%	2,000.0 µg/mL	+/- 11.6281 +/- 44.2527 +/- 44.4331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
42	1,2-Dibromoethane (EDB) CAS # 106-93-4 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
43	Chlorobenzene CAS # 108-90-7 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
44	1,1,1,2-Tetrachloroethane CAS # 630-20-6 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
45	m-Xylene CAS # 108-38-3 Purity 99%	1,000.0 µg/mL	+/- 5.8141 +/- 22.1265 +/- 22.2167	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
46	p-Xylene CAS # 106-42-3 Purity 99%	1,000.0 µg/mL	+/- 5.8141 +/- 22.1265 +/- 22.2167	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
47	o-Xylene CAS # 95-47-6 Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed

48	Ethylbenzene <b>CAS #</b> 100-41-4 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
49	Styrene <b>CAS #</b> 100-42-5 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
50	Isopropylbenzene (cumene) <b>CAS #</b> 98-82-8 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
51	bromoform <b>CAS #</b> 75-25-2 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
52	1,1,2,2-Tetrachloroethane <b>CAS #</b> 79-34-5 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
53	1,2,3-Trichloropropane <b>CAS #</b> 96-18-4 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
54	trans-1,4-dichloro-2-butene <b>CAS #</b> 110-57-6 <b>Purity</b> 98%	2,000.0	µg/mL	+/- 11.6281 +/- 44.2527 +/- 44.4331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
55	n-Propylbenzene <b>CAS #</b> 103-65-1 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
56	Bromobenzene <b>CAS #</b> 108-86-1 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
57	1,3,5-Trimethylbenzene <b>CAS #</b> 108-67-8 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
58	2-Chlorotoluene <b>CAS #</b> 95-49-8 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
59	4-Chlorotoluene <b>CAS #</b> 106-43-4 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
60	tert-Butylbenzene <b>CAS #</b> 98-06-6 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
61	1,2,4-Trimethylbenzene <b>CAS #</b> 95-63-6 <b>Purity</b> 98%	2,000.0	µg/mL	+/- 11.6281 +/- 44.2527 +/- 44.4331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
62	sec-Butylbenzene <b>CAS #</b> 135-98-8 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
63	4-Isopropyltoluene (p-Cymene) <b>CAS #</b> 99-87-6 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
64	1,3-Dichlorobenzene <b>CAS #</b> 541-73-1 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
65	1,4-Dichlorobenzene <b>CAS #</b> 106-46-7 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
66	n-Butylbenzene <b>CAS #</b> 104-51-8 <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed

67	1,2-Dichlorobenzene CAS # 95-50-1 Purity 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
68	1,2-Dibromo-3-chloropropane CAS # 96-12-8 Purity 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
69	1,2,4-Trichlorobenzene CAS # 120-82-1 Purity 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
70	Hexachlorobutadiene CAS # 87-68-3 Purity 97%	2,000.0	µg/mL	+/- 11.6284 +/- 44.2540 +/- 44.4344	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
71	Naphthalene CAS # 91-20-3 Purity 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
72	1,2,3-Trichlorobenzene CAS # 87-61-6 Purity 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed

**Solvent:** P&T Methanol  
**CAS #** 67-56-1  
**Purity** 99%

**Column:**  
 60m x .25mm x 1.4µm  
 Rtx-502.2 (cat.#10916)

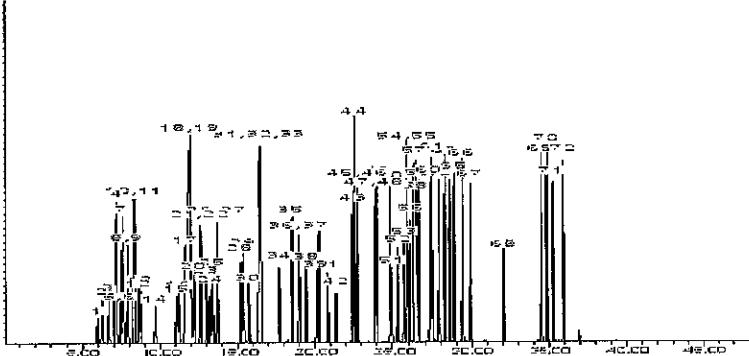
**Carrier Gas:**  
 helium-constant pressure 30 psi

**Temp. Program:**  
 40°C (hold 6 min.) to 240°C  
 @ 6°C/min. (hold 10 min.)

**Inj. Temp:**  
 200°C

**Det. Temp:**  
 250°C

**Det. Type:**  
 MSD



*Jennifer L. Pollino*  
 Jennifer L. Pollino - QC Analyst

Date Passed: 01-Mar-2013 Balance: B251644995

Manufactured under Restek's ISO 9001:2008  
 Registered Quality System  
 Certificate #FM 80397

Reagent

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**VOA8260MEGA2\_00011**



110 Benner Circle  
Bellefonte, PA 16823-8812  
Tel: (800)356-1688  
Fax: (814)353-1309  
[www.restek.com](http://www.restek.com)



## Certificate of Analysis

### FOR LABORATORY USE ONLY-READ MSDS PRIOR TO USE.

*This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.*

**Catalog No. :** 567641.sec

**Lot No.:** A093733

**Description :** 8260 List 1 / Std #1 MegaMix

8260 List 1 / Std #1 MegaMix 1,000-50,000 µg/ml, P&T Methanol, 1 ml/ampul

**Container Size :** 2 mL

**Pkg Amt:** > 1 mL

**Expiration Date :** February 2016

**Storage:** 0°C or colder

### C E R T I F I E D   V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Diethyl ether (ethyl ether)	2,000.0 µg/mL	+/- 11.6282	µg/mL	Gravimetric
	CAS # 60-29-7.SEC		+/- 44.2531	µg/mL	Unstressed
	Purity 99%		+/- 44.4335	µg/mL	Stressed
2	1,1,2-Trichlorotrifluoroethane (CFC-113)	2,000.0 µg/mL	+/- 11.6282	µg/mL	Gravimetric
	CAS # 76-13-1.SEC		+/- 44.2531	µg/mL	Unstressed
	Purity 99%		+/- 44.4335	µg/mL	Stressed
3	1,1-Dichloroethene	2,000.0 µg/mL	+/- 11.6282	µg/mL	Gravimetric
	CAS # 75-35-4.SEC		+/- 44.2531	µg/mL	Unstressed
	Purity 99%		+/- 44.4335	µg/mL	Stressed
4	tert-Butanol (TBA)	20,000.0 µg/mL	+/- 116.2756	µg/mL	Gravimetric
	CAS # 75-65-0.SEC		+/- 442.5291	µg/mL	Unstressed
	Purity 99%		+/- 444.3332	µg/mL	Stressed
5	Iodomethane (methyl iodide)	2,000.0 µg/mL	+/- 11.6284	µg/mL	Gravimetric
	CAS # 74-88-4.SEC		+/- 44.2540	µg/mL	Unstressed
	Purity 97%		+/- 44.4344	µg/mL	Stressed
6	Allyl chloride ( 3-chloropropene )	2,000.0 µg/mL	+/- 11.6281	µg/mL	Gravimetric
	CAS # 107-05-1.SEC		+/- 44.2527	µg/mL	Unstressed
	Purity 98%		+/- 44.4331	µg/mL	Stressed
7	Methyl acetate	10,000.0 µg/mL	+/- 58.1378	µg/mL	Gravimetric
	CAS # 79-20-9.SEC		+/- 221.2646	µg/mL	Unstressed
	Purity 99%		+/- 222.1666	µg/mL	Stressed
8	Carbon disulfide	2,000.0 µg/mL	+/- 11.6281	µg/mL	Gravimetric
	CAS # 75-15-0.SEC		+/- 44.2527	µg/mL	Unstressed
	Purity 98%		+/- 44.4331	µg/mL	Stressed
9	Methylene chloride (dichloromethane)	2,000.0 µg/mL	+/- 11.6282	µg/mL	Gravimetric
	CAS # 75-09-2.SEC		+/- 44.2531	µg/mL	Unstressed
	Purity 99%		+/- 44.4335	µg/mL	Stressed

10	Acrylonitrile <b>CAS #</b> 107-13-1.SEC <b>Purity</b> 99%	20,000.0	$\mu\text{g/mL}$	+/-	116.2756	$\mu\text{g/mL}$	Gravimetric
				+/-	442.5291	$\mu\text{g/mL}$	Unstressed
				+/-	444.3332	$\mu\text{g/mL}$	Stressed
11	Methyl-tert-butyl ether ( MTBE ) <b>CAS #</b> 1634-04-4.SEC <b>Purity</b> 99%	2,000.0	$\mu\text{g/mL}$	+/-	11.6282	$\mu\text{g/mL}$	Gravimetric
				+/-	44.2531	$\mu\text{g/mL}$	Unstressed
				+/-	44.4335	$\mu\text{g/mL}$	Stressed
12	cis-1,2-Dichloroethene <b>CAS #</b> 156-59-2.SEC <b>Purity</b> 99%	2,000.0	$\mu\text{g/mL}$	+/-	11.6282	$\mu\text{g/mL}$	Gravimetric
				+/-	44.2531	$\mu\text{g/mL}$	Unstressed
				+/-	44.4335	$\mu\text{g/mL}$	Stressed
13	n-Hexane (C6) <b>CAS #</b> 110-54-3.SEC <b>Purity</b> 98%	2,000.1	$\mu\text{g/mL}$	+/-	11.6286	$\mu\text{g/mL}$	Gravimetric
				+/-	44.2549	$\mu\text{g/mL}$	Unstressed
				+/-	44.4353	$\mu\text{g/mL}$	Stressed
14	1,1-Dichloroethane <b>CAS #</b> 75-34-3.SEC <b>Purity</b> 97%	2,000.0	$\mu\text{g/mL}$	+/-	11.6284	$\mu\text{g/mL}$	Gravimetric
				+/-	44.2540	$\mu\text{g/mL}$	Unstressed
				+/-	44.4344	$\mu\text{g/mL}$	Stressed
15	2,2-Dichloropropane <b>CAS #</b> 594-20-7.SEC <b>Purity</b> 99%	2,000.0	$\mu\text{g/mL}$	+/-	11.6282	$\mu\text{g/mL}$	Gravimetric
				+/-	44.2531	$\mu\text{g/mL}$	Unstressed
				+/-	44.4335	$\mu\text{g/mL}$	Stressed
16	trans-1,2-Dichloroethene <b>CAS #</b> 156-60-5.SEC <b>Purity</b> 97%	2,000.0	$\mu\text{g/mL}$	+/-	11.6284	$\mu\text{g/mL}$	Gravimetric
				+/-	44.2540	$\mu\text{g/mL}$	Unstressed
				+/-	44.4344	$\mu\text{g/mL}$	Stressed
17	Chloroform <b>CAS #</b> 67-66-3.SEC <b>Purity</b> 99%	2,000.0	$\mu\text{g/mL}$	+/-	11.6282	$\mu\text{g/mL}$	Gravimetric
				+/-	44.2531	$\mu\text{g/mL}$	Unstressed
				+/-	44.4335	$\mu\text{g/mL}$	Stressed
18	Isobutanol (2-Methyl-1-propanol) <b>CAS #</b> 78-83-1.SEC <b>Purity</b> 99%	50,000.0	$\mu\text{g/mL}$	+/-	290.6891	$\mu\text{g/mL}$	Gravimetric
				+/-	1,106.3228	$\mu\text{g/mL}$	Unstressed
				+/-	1,110.8331	$\mu\text{g/mL}$	Stressed
19	Bromochloromethane <b>CAS #</b> 74-97-5.SEC <b>Purity</b> 99%	2,000.0	$\mu\text{g/mL}$	+/-	11.6282	$\mu\text{g/mL}$	Gravimetric
				+/-	44.2531	$\mu\text{g/mL}$	Unstressed
				+/-	44.4335	$\mu\text{g/mL}$	Stressed
20	Tetrahydrofuran <b>CAS #</b> 109-99-9.SEC <b>Purity</b> 99%	4,000.0	$\mu\text{g/mL}$	+/-	23.2563	$\mu\text{g/mL}$	Gravimetric
				+/-	88.5061	$\mu\text{g/mL}$	Unstressed
				+/-	88.8670	$\mu\text{g/mL}$	Stressed
21	1,1,1-Trichloroethane <b>CAS #</b> 71-55-6.SEC <b>Purity</b> 99%	2,000.0	$\mu\text{g/mL}$	+/-	11.6282	$\mu\text{g/mL}$	Gravimetric
				+/-	44.2531	$\mu\text{g/mL}$	Unstressed
				+/-	44.4335	$\mu\text{g/mL}$	Stressed
22	Cyclohexane <b>CAS #</b> 110-82-7.SEC <b>Purity</b> 99%	2,000.0	$\mu\text{g/mL}$	+/-	11.6282	$\mu\text{g/mL}$	Gravimetric
				+/-	44.2531	$\mu\text{g/mL}$	Unstressed
				+/-	44.4335	$\mu\text{g/mL}$	Stressed
23	1,1-Dichloropropene <b>CAS #</b> 563-58-6.SEC <b>Purity</b> 98%	2,010.5	$\mu\text{g/mL}$	+/-	11.6890	$\mu\text{g/mL}$	Gravimetric
				+/-	44.4847	$\mu\text{g/mL}$	Unstressed
				+/-	44.6661	$\mu\text{g/mL}$	Stressed
24	Carbon tetrachloride <b>CAS #</b> 56-23-5.SEC <b>Purity</b> 98%	2,000.1	$\mu\text{g/mL}$	+/-	11.6286	$\mu\text{g/mL}$	Gravimetric
				+/-	44.2549	$\mu\text{g/mL}$	Unstressed
				+/-	44.4353	$\mu\text{g/mL}$	Stressed
25	n-Heptane (C7) <b>CAS #</b> 142-82-5.SEC <b>Purity</b> 99%	2,000.1	$\mu\text{g/mL}$	+/-	11.6288	$\mu\text{g/mL}$	Gravimetric
				+/-	44.2553	$\mu\text{g/mL}$	Unstressed
				+/-	44.4357	$\mu\text{g/mL}$	Stressed
26	Benzene <b>CAS #</b> 71-43-2.SEC <b>Purity</b> 99%	2,000.0	$\mu\text{g/mL}$	+/-	11.6282	$\mu\text{g/mL}$	Gravimetric
				+/-	44.2531	$\mu\text{g/mL}$	Unstressed
				+/-	44.4335	$\mu\text{g/mL}$	Stressed
27	1,2-Dichloroethane <b>CAS #</b> 107-06-2.SEC <b>Purity</b> 99%	2,000.0	$\mu\text{g/mL}$	+/-	11.6282	$\mu\text{g/mL}$	Gravimetric
				+/-	44.2531	$\mu\text{g/mL}$	Unstressed
				+/-	44.4335	$\mu\text{g/mL}$	Stressed
28	Trichloroethylene <b>CAS #</b> 79-01-6.SEC <b>Purity</b> 98%	2,000.1	$\mu\text{g/mL}$	+/-	11.6286	$\mu\text{g/mL}$	Gravimetric
				+/-	44.2549	$\mu\text{g/mL}$	Unstressed
				+/-	44.4353	$\mu\text{g/mL}$	Stressed

29	Methylcyclohexane CAS # 108-87-2.SEC Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
30	1,2-Dichloropropane CAS # 78-87-5.SEC Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
31	1,4-Dioxane CAS # 123-91-1.SEC Purity 99%	40,000.0 µg/mL	+/- 232.5513 +/- 885.0582 +/- 888.6665	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
32	Dibromomethane CAS # 74-95-3.SEC Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
33	Bromodichloromethane CAS # 75-27-4.SEC Purity 97%	2,000.1 µg/mL	+/- 11.6290 +/- 44.2562 +/- 44.4366	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
34	cis-1,3-Dichloropropene CAS # 10061-01-5.SEC Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
35	Toluene CAS # 108-88-3.SEC Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
36	Ethyl methacrylate CAS # 97-63-2.SEC Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
37	trans-1,3-Dichloropropene CAS # 10061-02-6.SEC Purity 98%	2,000.0 µg/mL	+/- 11.6281 +/- 44.2527 +/- 44.4331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
38	1,1,2-Trichloroethane CAS # 79-00-5.SEC Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
39	1,3-Dichloropropane CAS # 142-28-9.SEC Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
40	Tetrachloroethene CAS # 127-18-4.SEC Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
41	Dibromochloromethane CAS # 124-48-1.SEC Purity 97%	2,000.1 µg/mL	+/- 11.6290 +/- 44.2562 +/- 44.4366	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
42	1,2-Dibromoethane (EDB) CAS # 106-93-4.SEC Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
43	Chlorobenzene CAS # 108-90-7.SEC Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
44	1,1,1,2-Tetrachloroethane CAS # 630-20-6.SEC Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
45	m-Xylene CAS # 108-38-3.SEC Purity 99%	1,000.0 µg/mL	+/- 5.8141 +/- 22.1265 +/- 22.2167	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
46	p-Xylene CAS # 106-42-3.SEC Purity 99%	1,000.0 µg/mL	+/- 5.8141 +/- 22.1265 +/- 22.2167	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
47	o-Xylene CAS # 95-47-6.SEC Purity 99%	2,000.0 µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed

48	Ethylbenzene <b>CAS #</b> 100-41-4.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
49	Styrene <b>CAS #</b> 100-42-5.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
50	Isopropylbenzene (cumene) <b>CAS #</b> 98-82-8.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
51	Bromoform <b>CAS #</b> 75-25-2.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
52	1,1,2,2-Tetrachloroethane <b>CAS #</b> 79-34-5.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
53	1,2,3-Trichloropropane <b>CAS #</b> 96-18-4.SEC <b>Purity</b> 98%	2,000.0	µg/mL	+/- 11.6281 +/- 44.2527 +/- 44.4331	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
54	trans-1,4-Dichloro-2-butene <b>CAS #</b> 110-57-6.SEC <b>Purity</b> 97%	2,000.0	µg/mL	+/- 11.6284 +/- 44.2540 +/- 44.4344	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
55	n-Propylbenzene <b>CAS #</b> 103-65-1.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
56	Bromobenzene <b>CAS #</b> 108-86-1.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
57	1,3,5-Trimethylbenzene <b>CAS #</b> 108-67-8.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
58	2-Chlorotoluene <b>CAS #</b> 95-49-8.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
59	4-Chlorotoluene <b>CAS #</b> 106-43-4.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
60	tert-Butylbenzene <b>CAS #</b> 98-06-6.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
61	1,2,4-Trimethylbenzene <b>CAS #</b> 95-63-6.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
62	sec-Butylbenzene <b>CAS #</b> 135-98-8.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
63	4-Isopropyltoluene (p-cymene) <b>CAS #</b> 99-87-6.SEC <b>Purity</b> 96%	2,000.1	µg/mL	+/- 11.6285 +/- 44.2545 +/- 44.4349	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
64	1,3-Dichlorobenzene <b>CAS #</b> 541-73-1.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
65	1,4-Dichlorobenzene <b>CAS #</b> 106-46-7.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
66	n-Butylbenzene <b>CAS #</b> 104-51-8.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282 +/- 44.2531 +/- 44.4335	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed

67	1,2-Dichlorobenzene <b>CAS #</b> 95-50-1.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282	µg/mL	Gravimetric
				+/- 44.2531	µg/mL	Unstressed
				+/- 44.4335	µg/mL	Stressed
68	1,2-Dibromo-3-chloropropane <b>CAS #</b> 96-12-8.SEC <b>Purity</b> 97%	2,000.0	µg/mL	+/- 11.6284	µg/mL	Gravimetric
				+/- 44.2540	µg/mL	Unstressed
				+/- 44.4344	µg/mL	Stressed
69	1,2,4-Trichlorobenzene <b>CAS #</b> 120-82-1.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282	µg/mL	Gravimetric
				+/- 44.2531	µg/mL	Unstressed
				+/- 44.4335	µg/mL	Stressed
70	Hexachlorobutadiene <b>CAS #</b> 87-68-3.SEC <b>Purity</b> 97%	2,000.0	µg/mL	+/- 11.6284	µg/mL	Gravimetric
				+/- 44.2540	µg/mL	Unstressed
				+/- 44.4344	µg/mL	Stressed
71	Naphthalene <b>CAS #</b> 91-20-3.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282	µg/mL	Gravimetric
				+/- 44.2531	µg/mL	Unstressed
				+/- 44.4335	µg/mL	Stressed
72	1,2,3-Trichlorobenzene <b>CAS #</b> 87-61-6.SEC <b>Purity</b> 99%	2,000.0	µg/mL	+/- 11.6282	µg/mL	Gravimetric
				+/- 44.2531	µg/mL	Unstressed
				+/- 44.4335	µg/mL	Stressed

**Solvent:** P&T Methanol  
**CAS #** 67-56-1  
**Purity** 99%

**Column:**

60m x .25mm x 1.4µm  
Rtx-502.2 (cat.#10916)

**Carrier Gas:**

helium-constant pressure 30 psi

**Temp. Program:**

40°C (hold 6 min.) to 240°C  
@ 6°C/min. (hold 10 min.)

**Inj. Temp:**

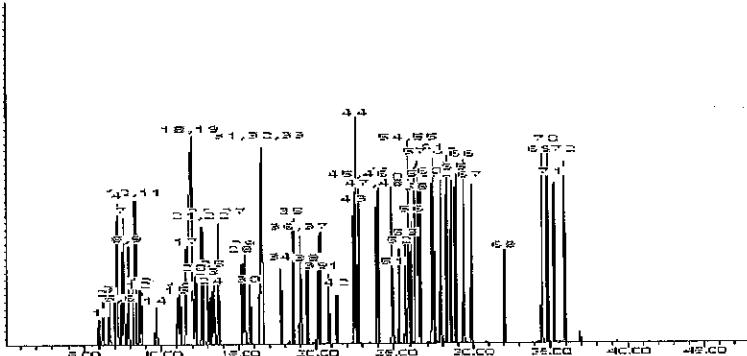
200°C

**Det. Temp:**

250°C

**Det. Type:**

MSD



*Jennifer L. Pollino*  
Jennifer L. Pollino - QC Analyst

Date Passed: 01-Mar-2013 Balance: 1127510105

Manufactured under Restek's ISO 9001:2008 Registered Quality System Certificate #FM 80397
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Reagent

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**VOA8260SURRES\_00063**



# CERTIFIED REFERENCE MATERIAL

110 Benner Circle  
Bellefonte, PA 16823-8812  
Tel: (800)356-1688  
Fax: (814)353-1309

[www.restek.com](http://www.restek.com)



## Certificate of Analysis



### FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

**Catalog No.:** 567650

**Lot No.:** A0100424

**Description :** 8260 Surrogate Standard

8260 Surrogate Standard 2,500 ug/ml, P&T Methanol, 5 mL/ampul

**Container Size :** 5 mL

**Pkg Amt:** > 5 mL

**Expiration Date :** January 31, 2019

**Storage:** 0°C or colder

### C E R T I F I E D V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Dibromofluoromethane <b>CAS #</b> 1868-53-7 <b>Purity</b> 99%	2,502.2 µg/mL	+/- 14.5480	µg/mL	Gravimetric
	(Lot 022012)		+/- 28.2159	µg/mL	Unstressed
			+/- 32.4683	µg/mL	Stressed
2	1,2-Dichloroethane-d4 <b>CAS #</b> 17060-07-0 <b>Purity</b> 99%	2,501.2 µg/mL	+/- 14.5422	µg/mL	Gravimetric
	(Lot 12K-027)		+/- 28.2046	µg/mL	Unstressed
			+/- 32.4554	µg/mL	Stressed
3	Toluene-d8 <b>CAS #</b> 2037-26-5 <b>Purity</b> 99%	2,500.8 µg/mL	+/- 14.5399	µg/mL	Gravimetric
	(Lot 13I-050)		+/- 28.2001	µg/mL	Unstressed
			+/- 32.4502	µg/mL	Stressed
4	1-Bromo-4-fluorobenzene (BFB) <b>CAS #</b> 460-00-4 <b>Purity</b> 99%	2,501.4 µg/mL	+/- 14.5434	µg/mL	Gravimetric
	(Lot 01127COV)		+/- 28.2069	µg/mL	Unstressed
			+/- 32.4580	µg/mL	Stressed

**Solvent:** P&T Methanol  
**CAS #** 67-56-1  
**Purity** 99%

Reagent

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**VOA8260SURRES\_00087**



# CERTIFIED REFERENCE MATERIAL

110 Benner Circle  
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## Certificate of Analysis



### FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

**Catalog No.:** 567650

**Lot No.:** A0102817

**Description :** 8260 Surrogate Standard

8260 Surrogate Standard 2,500 ug/ml, P&T Methanol, 5 mL/ampul

**Container Size :** 5 mL

**Pkg Amt:** > 5 mL

**Expiration Date :** April 30, 2019

**Storage:** 0°C or colder

### C E R T I F I E D V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Dibromofluoromethane <b>CAS #</b> 1868-53-7 <b>Purity</b> 99%	2,503.8 µg/mL	+/- 14.5573	µg/mL	Gravimetric
	(Lot 022012)		+/- 28.2339	µg/mL	Unstressed
			+/- 32.4891	µg/mL	Stressed
2	1,2-Dichloroethane-d4 <b>CAS #</b> 17060-07-0 <b>Purity</b> 99%	2,502.4 µg/mL	+/- 14.5492	µg/mL	Gravimetric
	(Lot 13J-483)		+/- 28.2182	µg/mL	Unstressed
			+/- 32.4709	µg/mL	Stressed
3	Toluene-d8 <b>CAS #</b> 2037-26-5 <b>Purity</b> 99%	2,500.0 µg/mL	+/- 14.5352	µg/mL	Gravimetric
	(Lot 13I-050)		+/- 28.1911	µg/mL	Unstressed
			+/- 32.4398	µg/mL	Stressed
4	1-Bromo-4-fluorobenzene (BFB) <b>CAS #</b> 460-00-4 <b>Purity</b> 99%	2,503.6 µg/mL	+/- 14.5561	µg/mL	Gravimetric
	(Lot 01127COV)		+/- 28.2317	µg/mL	Unstressed
			+/- 32.4865	µg/mL	Stressed

**Solvent:** P&T Methanol  
**CAS #** 67-56-1  
**Purity** 99%

Reagent

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**VOA8260VARES\_00050**



# CERTIFIED REFERENCE MATERIAL

110 Benner Circle  
Bellefonte, PA 16823-8812  
Tel: (800)356-1688  
Fax: (814)353-1309

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## Certificate of Analysis



### FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

*This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.*

**Catalog No.:** 569724

**Lot No.:** A0108225

**Description :** 8260 List 1 / Std #6 Vinyl Acetate (2015)

8260 List 1 / Std #6 Vinyl Acetate (2015) 5000 ug/ml, P&T Methanol, 1 ml/ampul

**Container Size :** 2 mL

**Pkg Amt:** > 1 mL

**Expiration Date :** July 31, 2015

**Storage:** 0°C or colder

### C E R T I F I E D V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Vinyl acetate <b>CAS #</b> 108-05-4 <b>Purity</b> 99%	5,000.0 µg/mL	+/- 29.3428	µg/mL	Gravimetric

**Solvent:** P&T Methanol  
**CAS #** 67-56-1  
**Purity** 99%

#### Tech Tips:

Vinyl acetate is a volatile organic ester included in the target lists of several US EPA and other methods. Under acidic conditions, esters react with alcohols to form new esters (transesterification). Methanol-based mixes containing halogenated compounds are slightly acidic, so it is important to minimize exposure of vinyl acetate to mixes of halogenated compounds in methanol. For this reason, we offer vinyl acetate in individual solution, and suggest that it be introduced into the working level calibration solution immediately before use. This will minimize problems and ensure more consistent results.

Reagent

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**VOAACRORES\_00064**



# CERTIFIED REFERENCE MATERIAL

110 Benner Circle  
Bellefonte, PA 16823-8812  
Tel: (800)356-1688  
Fax: (814)353-1309

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## Certificate of Analysis

### FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

*This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.*

**Catalog No.:** 568720      **Lot No.:** A0107338  
**Description :** 8260 List 1/Std #5 Acrolein High  
                 8260 List 1/Std #5 Acrolein High 19,750 µg/mL, Water, 1 mL/ampul  
**Container Size :** 2 mL      **Pkg Amt:** > 1 mL  
**Expiration Date :** March 31, 2015      **Storage:** 10°C or colder  
**Handling:** This product is photosensitive.

### C E R T I F I E D   V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Acrolein CAS # 107-02-8 Purity 99%	19,759.0 µg/mL (Lot 140429JLM)	+/- 115.6933 µg/mL	+/- 633.5357 µg/mL	Gravimetric Unstressed Stressed

**Solvent:** Water  
CAS # 7732-18-5  
Purity 99%

Reagent

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**VOARESEE1ST\_00008**

110 Benner Circle  
 Bellefonte, PA 16823-8812  
 Tel: (800)356-1688  
 Fax: (814)353-1309

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## Certificate of Analysis

### FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

*This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.*

**Catalog No. :** 568363-FL

**Lot No.:** A097285

**Description :** Custom EE Standard

Custom EE Standard 5,000 $\mu$ g/mL, P&T Methanol, 1mL/ampul

**Container Size :** 2 mL

**Pkg Amt:** > 1 mL

**Expiration Date :** February 28, 2015

**Storage:** 0°C or colder

### C E R T I F I E D V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	3-Chlorobenzotrifluoride	5,001.0 $\mu$ g/mL	+/- 29.3487	$\mu$ g/mL	Gravimetric
	CAS # 98-15-7	(Lot 21324DO)	+/- 53.0822	$\mu$ g/mL	Unstressed
	Purity 99%		+/- 61.7282	$\mu$ g/mL	Stressed
2	4-Chlorobenzotrifluoride	5,003.0 $\mu$ g/mL	+/- 29.3604	$\mu$ g/mL	Gravimetric
	CAS # 98-56-6	(Lot 08507BO)	+/- 53.1034	$\mu$ g/mL	Unstressed
	Purity 99%		+/- 61.7529	$\mu$ g/mL	Stressed
3	2-Chlorobenzotrifluoride	5,005.0 $\mu$ g/mL	+/- 29.3721	$\mu$ g/mL	Gravimetric
	CAS # 88-16-4	(Lot I0316DQ)	+/- 53.1247	$\mu$ g/mL	Unstressed
	Purity 99%		+/- 61.7775	$\mu$ g/mL	Stressed
4	3-Chlorotoluene	5,000.0 $\mu$ g/mL	+/- 29.3428	$\mu$ g/mL	Gravimetric
	CAS # 108-41-8	(Lot 13528LX)	+/- 53.0716	$\mu$ g/mL	Unstressed
	Purity 99%		+/- 61.7158	$\mu$ g/mL	Stressed
5	2,4-Dichlorobenzotrifluoride	5,002.0 $\mu$ g/mL	+/- 29.3545	$\mu$ g/mL	Gravimetric
	CAS # 320-60-5	(Lot MKBL3552V)	+/- 53.0928	$\mu$ g/mL	Unstressed
	Purity 99%		+/- 61.7405	$\mu$ g/mL	Stressed
6	3,4-Dichlorobenzotrifluoride	5,000.0 $\mu$ g/mL	+/- 29.3428	$\mu$ g/mL	Gravimetric
	CAS # 328-84-7	(Lot 11105EJV)	+/- 53.0716	$\mu$ g/mL	Unstressed
	Purity 99%		+/- 61.7158	$\mu$ g/mL	Stressed
7	2,5-Dichlorobenzotrifluoride	5,000.0 $\mu$ g/mL	+/- 29.3428	$\mu$ g/mL	Gravimetric
	CAS # 320-50-3	(Lot 04415DSV)	+/- 53.0716	$\mu$ g/mL	Unstressed
	Purity 99%		+/- 61.7158	$\mu$ g/mL	Stressed
8	2,4-Dichlorotoluene	5,002.0 $\mu$ g/mL	+/- 29.3545	$\mu$ g/mL	Gravimetric
	CAS # 95-73-8	(Lot 07715JS)	+/- 53.0928	$\mu$ g/mL	Unstressed
	Purity 99%		+/- 61.7405	$\mu$ g/mL	Stressed

9	2,5-Dichlorotoluene <b>CAS #</b> 19398-61-9 <b>Purity</b> 99%	(Lot 10119CU)	5,000.0	µg/mL	+/- 29.3428 +/- 53.0716 +/- 61.7158	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
10	2,6-Dichlorotoluene <b>CAS #</b> 118-69-4 <b>Purity</b> 99%	(Lot 16921JS)	5,001.0	µg/mL	+/- 29.3487 +/- 53.0822 +/- 61.7282	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
11	3,4-Dichlorotoluene <b>CAS #</b> 95-75-0 <b>Purity</b> 99%	(Lot 09419AS)	5,003.0	µg/mL	+/- 29.3604 +/- 53.1034 +/- 61.7529	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
12	2,3-Dichlorotoluene <b>CAS #</b> 32768-54-0 <b>Purity</b> 99%	(Lot 00317)	5,008.0	µg/mL	+/- 29.3897 +/- 53.1565 +/- 61.8146	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
13	2,4,5-Trichlorotoluene <b>CAS #</b> 6639-30-1 <b>Purity</b> 99%	(Lot 1767300)	5,001.0	µg/mL	+/- 29.3487 +/- 53.0822 +/- 61.7282	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
14	2,3,6-Trichlorotoluene <b>CAS #</b> 2077-46-5 <b>Purity</b> 99%	(Lot RM01250)	5,001.0	µg/mL	+/- 29.3487 +/- 53.0822 +/- 61.7282	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed

**Solvent:** P&T Methanol  
**CAS #** 67-56-1  
**Purity** 99%

# **Method 8260C Low Level**

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**Volatile Organic Compounds (GC/MS)  
by Method 8260C Low Level**

FORM II  
GC/MS VOA SURROGATE RECOVERY

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Matrix: Water Level: Low  
GC Column (1): DB-624 ID: 0.18 (mm)

Client Sample ID	Lab Sample ID	DBFM #	DCA #	TOL #	BFB #
HD-MW-165-0/1-0	180-42975-1	110	118	100	96
HD-MW-162-0/1-0	180-42975-2	116	115	98	91
HD-MW-162-0/1-0 DL	180-42975-2 DL	110	107	102	98
HD-MW-169-0/1-0	180-42975-3	109	115	98	92
HD-QC2-0/1-2	180-42975-4	112	114	103	97
	MB 180-138583/5	107	113	101	93
	MB 180-138685/5	111	116	100	97
	LCS 180-138583/8	91	95	99	94
	LCS 180-138685/8	93	93	101	94
	LCSD 180-138685/10	89	92	97	92
HD-MW-165-0/1-0 MS	180-42975-1 MS	92	90	99	89
HD-MW-165-0/1-0 MSD	180-42975-1 MSD	92	96	96	93

DBFM = Dibromofluoromethane (Surr)  
DCA = 1,2-Dichloroethane-d4 (Surr)  
TOL = Toluene-d8 (Surr)  
BFB = 4-Bromofluorobenzene (Surr)

QC LIMITS
70-128
64-135
71-118
70-118

# Column to be used to flag recovery values

FORM II 8260C

FORM III  
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Matrix: Water Level: Low Lab File ID: 50415008.D  
Lab ID: LCS 180-138583/8 Client ID: \_\_\_\_\_

COMPOUND	SPIKE ADDED (ug/L)	LCS CONCENTRATION (ug/L)	LCS % REC	QC LIMITS REC	#
Chloromethane	10.0	10.1	101	50-139	
Vinyl chloride	10.0	10.0	100	53-138	
Bromomethane	10.0	8.98	90	33-150	
Chloroethane	10.0	10.5	105	36-142	
1,1-Dichloroethene	10.0	7.04	70	65-136	
Acetone	20.0	19.0	95	22-150	
Carbon disulfide	10.0	3.87	39	54-132	*
Methylene Chloride	10.0	7.62	76	63-129	
trans-1,2-Dichloroethene	10.0	7.74	77	73-126	
Methyl tert-butyl ether	10.0	7.94	79	64-123	
1,1-Dichloroethane	10.0	8.31	83	73-126	
cis-1,2-Dichloroethene	10.0	7.88	79	70-120	
Bromochloromethane	10.0	8.32	83	70-127	
2-Butanone (MEK)	20.0	16.7	83	39-138	
Chloroform	10.0	9.11	91	72-127	
1,1,1-Trichloroethane	10.0	8.87	89	63-133	
Carbon tetrachloride	10.0	9.33	93	55-150	
Benzene	10.0	8.86	89	80-120	
1,2-Dichloroethane	10.0	8.84	88	68-132	
Trichloroethene	10.0	8.40	84	73-120	
1,2-Dichloropropane	10.0	9.27	93	76-124	
Bromodichloromethane	10.0	8.85	88	66-130	
cis-1,3-Dichloropropene	10.0	9.60	96	66-120	
4-Methyl-2-pentanone (MIBK)	20.0	16.8	84	45-145	
Toluene	10.0	9.81	98	80-123	
trans-1,3-Dichloropropene	10.0	11.6	116	65-125	
1,1,2-Trichloroethane	10.0	11.2	112	77-127	
Tetrachloroethene	10.0	9.33	93	70-135	
2-Hexanone	20.0	15.3	76	25-132	
Dibromochloromethane	10.0	10.1	101	60-140	
1,2-Dibromoethane (EDB)	10.0	9.99	100	74-123	
Chlorobenzene	10.0	10.2	102	80-120	
1,1,1,2-Tetrachloroethane	10.0	11.8	118	63-140	
Ethylbenzene	10.0	9.72	97	72-126	
Xylenes, Total	20.0	19.0	95	76-128	
Styrene	10.0	9.70	97	71-127	
Bromoform	10.0	9.35	93	46-150	
1,1,2,2-Tetrachloroethane	10.0	10.9	109	62-125	
1,4-Dioxane	200	179 J	90	10-160	

# Column to be used to flag recovery and RPD values

FORM III  
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Matrix: Water Level: Low Lab File ID: 50416008.D  
Lab ID: LCS 180-138685/8 Client ID: \_\_\_\_\_

COMPOUND	SPIKE ADDED (ug/L)	LCS CONCENTRATION (ug/L)	LCS % REC	QC LIMITS REC	#
Chloromethane	10.0	10.0	100	50-139	
Vinyl chloride	10.0	10.5	105	53-138	
Bromomethane	10.0	9.71	97	33-150	
Chloroethane	10.0	11.0	110	36-142	
1,1-Dichloroethene	10.0	9.09	91	65-136	
Acetone	20.0	21.0	105	22-150	
Carbon disulfide	10.0	6.75	68	54-132	
Methylene Chloride	10.0	9.92	99	63-129	
trans-1,2-Dichloroethene	10.0	9.42	94	73-126	
Methyl tert-butyl ether	10.0	9.22	92	64-123	
1,1-Dichloroethane	10.0	9.66	97	73-126	
cis-1,2-Dichloroethene	10.0	9.13	91	70-120	
Bromochloromethane	10.0	9.67	97	70-127	
2-Butanone (MEK)	20.0	17.2	86	39-138	
Chloroform	10.0	10.1	101	72-127	
1,1,1-Trichloroethane	10.0	10.6	106	63-133	
Carbon tetrachloride	10.0	11.0	110	55-150	
Benzene	10.0	10.2	102	80-120	
1,2-Dichloroethane	10.0	10.0	100	68-132	
Trichloroethene	10.0	9.34	93	73-120	
1,2-Dichloropropane	10.0	9.73	97	76-124	
Bromodichloromethane	10.0	9.17	92	66-130	
cis-1,3-Dichloropropene	10.0	9.80	98	66-120	
4-Methyl-2-pentanone (MIBK)	20.0	16.2	81	45-145	
Toluene	10.0	10.5	105	80-123	
trans-1,3-Dichloropropene	10.0	10.8	108	65-125	
1,1,2-Trichloroethane	10.0	10.7	107	77-127	
Tetrachloroethene	10.0	10.2	102	70-135	
2-Hexanone	20.0	16.8	84	25-132	
Dibromochloromethane	10.0	9.51	95	60-140	
1,2-Dibromoethane (EDB)	10.0	10.2	102	74-123	
Chlorobenzene	10.0	10.3	103	80-120	
1,1,1,2-Tetrachloroethane	10.0	11.1	111	63-140	
Ethylbenzene	10.0	9.73	97	72-126	
Xylenes, Total	20.0	19.8	99	76-128	
Styrene	10.0	9.71	97	71-127	
Bromoform	10.0	9.19	92	46-150	
1,1,2,2-Tetrachloroethane	10.0	10.0	100	62-125	
1,4-Dioxane	200	167 J	84	10-160	

# Column to be used to flag recovery and RPD values

FORM III  
GC/MS VOA LAB CONTROL SAMPLE DUPLICATE RECOVERY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Matrix: Water Level: Low Lab File ID: 50416010.D

Lab ID: LCSD 180-138685/10 Client ID: \_\_\_\_\_

COMPOUND	SPIKE ADDED (ug/L)	LCSD CONCENTRATION (ug/L)	LCSD % REC	% RPD	QC LIMITS		#
					RPD	REC	
Chloromethane	10.0	10.8	108	8	35	50-139	
Vinyl chloride	10.0	10.3	103	1	35	53-138	
Bromomethane	10.0	8.47	85	14	35	33-150	
Chloroethane	10.0	10.7	107	3	35	36-142	
1,1-Dichloroethene	10.0	9.04	90	1	35	65-136	
Acetone	20.0	20.2	101	4	35	22-150	
Carbon disulfide	10.0	6.12	61	10	35	54-132	
Methylene Chloride	10.0	10.2	102	3	35	63-129	
trans-1,2-Dichloroethene	10.0	9.27	93	2	35	73-126	
Methyl tert-butyl ether	10.0	9.56	96	4	35	64-123	
1,1-Dichloroethane	10.0	9.47	95	2	35	73-126	
cis-1,2-Dichloroethene	10.0	9.45	95	3	35	70-120	
Bromoform	10.0	9.04	90	7	35	70-127	
2-Butanone (MEK)	20.0	17.7	89	3	35	39-138	
Chloroform	10.0	9.80	98	3	35	72-127	
1,1,1-Trichloroethane	10.0	10.2	102	3	35	63-133	
Carbon tetrachloride	10.0	10.3	103	7	35	55-150	
Benzene	10.0	10.1	101	1	32	80-120	
1,2-Dichloroethane	10.0	9.65	96	4	32	68-132	
Trichloroethene	10.0	9.24	92	1	35	73-120	
1,2-Dichloropropane	10.0	9.60	96	1	34	76-124	
Bromodichloromethane	10.0	8.86	89	3	35	66-130	
cis-1,3-Dichloropropene	10.0	9.88	99	1	35	66-120	
4-Methyl-2-pentanone (MIBK)	20.0	16.9	84	4	35	45-145	
Toluene	10.0	10.4	104	2	35	80-123	
trans-1,3-Dichloropropene	10.0	10.3	103	5	35	65-125	
1,1,2-Trichloroethane	10.0	10.4	104	2	35	77-127	
Tetrachloroethene	10.0	10.0	100	2	35	70-135	
2-Hexanone	20.0	17.3	86	3	35	25-132	
Dibromochloromethane	10.0	9.55	95	0	35	60-140	
1,2-Dibromoethane (EDB)	10.0	10.1	101	1	35	74-123	
Chlorobenzene	10.0	10.2	102	1	29	80-120	
1,1,1,2-Tetrachloroethane	10.0	10.7	107	4	34	63-140	
Ethylbenzene	10.0	9.71	97	0	33	72-126	
Xylenes, Total	20.0	19.2	96	3	32	76-128	
Styrene	10.0	9.79	98	1	34	71-127	
Bromoform	10.0	8.77	88	5	35	46-150	
1,1,2,2-Tetrachloroethane	10.0	9.87	99	2	35	62-125	
1,4-Dioxane	200	193 J	96	14	35	10-160	

# Column to be used to flag recovery and RPD values

FORM III  
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Matrix: Water Level: Low Lab File ID: 50415009.D  
Lab ID: 180-42975-1 MS Client ID: HD-MW-165-0/1-0 MS

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE CONCENTRATION (ug/L)	MS CONCENTRATION (ug/L)	MS % REC	QC LIMITS REC	#
Chloromethane	10.0	1.0 U	10.5	105	50-139	
Vinyl chloride	10.0	1.0 U	9.56	96	53-138	
Bromomethane	10.0	1.0 U	9.20	92	33-150	
Chloroethane	10.0	1.0 U	10.3	103	36-142	
1,1-Dichloroethene	10.0	1.0 U	6.75	67	65-136	
Acetone	20.0	5.0 U	20.2	101	22-150	
Carbon disulfide	10.0	1.0 U	3.81	38	54-132	F1
Methylene Chloride	10.0	1.0 U	7.73	77	63-129	
trans-1,2-Dichloroethene	10.0	1.0 U	7.54	75	73-126	
Methyl tert-butyl ether	10.0	1.0 U	8.13	81	64-123	
1,1-Dichloroethane	10.0	1.0 U	8.29	83	73-126	
cis-1,2-Dichloroethene	10.0	1.0 U	8.26	83	70-120	
Bromoform	10.0	1.0 U	7.70	77	70-127	
2-Butanone (MEK)	20.0	5.0 U	16.6	83	39-138	
Chloroform	10.0	1.0 U	9.04	90	72-127	
1,1,1-Trichloroethane	10.0	1.0 U	8.69	87	63-133	
Carbon tetrachloride	10.0	1.0 U	8.96	90	55-150	
Benzene	10.0	1.0 U	8.62	86	80-120	
1,2-Dichloroethane	10.0	1.0 U	8.54	85	68-132	
Trichloroethene	10.0	16	19.5	37	73-120	F1
1,2-Dichloropropane	10.0	1.0 U	8.78	88	76-124	
Bromodichloromethane	10.0	1.0 U	8.29	83	66-130	
cis-1,3-Dichloropropene	10.0	1.0 U	8.80	88	66-120	
4-Methyl-2-pentanone (MIBK)	20.0	5.0 U	16.8	84	45-145	
Toluene	10.0	1.0 U	9.59	96	80-123	
trans-1,3-Dichloropropene	10.0	1.0 U	11.0	110	65-125	
1,1,2-Trichloroethane	10.0	1.0 U	10.7	107	77-127	
Tetrachloroethene	10.0	7.4	15.2	78	70-135	
2-Hexanone	20.0	5.0 U	15.4	77	25-132	
Dibromochloromethane	10.0	1.0 U	9.09	91	60-140	
1,2-Dibromoethane (EDB)	10.0	1.0 U	9.76	98	74-123	
Chlorobenzene	10.0	1.0 U	9.83	98	80-120	
1,1,1,2-Tetrachloroethane	10.0	1.0 U	10.7	107	63-140	
Ethylbenzene	10.0	1.0 U	9.18	92	72-126	
Xylenes, Total	20.0	3.0 U	18.5	93	76-128	
Styrene	10.0	1.0 U	9.29	93	71-127	
Bromoform	10.0	1.0 U	8.75	87	46-150	
1,1,2,2-Tetrachloroethane	10.0	1.0 U	10.1	101	62-125	
1,4-Dioxane	200	200 U	175 J	87	10-160	

# Column to be used to flag recovery and RPD values

FORM III  
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Matrix: Water Level: Low Lab File ID: 50415010.D

Lab ID: 180-42975-1 MSD Client ID: HD-MW-165-0/1-0 MSD

COMPOUND	SPIKE ADDED (ug/L)	MSD CONCENTRATION (ug/L)	MSD % REC	% RPD	QC LIMITS		#
					RPD	REC	
Chloromethane	10.0	10.8	108	3	35	50-139	
Vinyl chloride	10.0	9.37	94	2	35	53-138	
Bromomethane	10.0	10.1	101	9	35	33-150	
Chloroethane	10.0	10.4	104	0	35	36-142	
1,1-Dichloroethene	10.0	6.67	67	1	35	65-136	
Acetone	20.0	19.1	96	5	35	22-150	
Carbon disulfide	10.0	3.88	39	2	35	54-132	F1
Methylene Chloride	10.0	8.19	82	6	35	63-129	
trans-1,2-Dichloroethene	10.0	7.67	77	2	35	73-126	
Methyl tert-butyl ether	10.0	8.73	87	7	35	64-123	
1,1-Dichloroethane	10.0	8.39	84	1	35	73-126	
cis-1,2-Dichloroethene	10.0	8.61	86	4	35	70-120	
Bromoform	10.0	9.01	90	16	35	70-127	
2-Butanone (MEK)	20.0	18.7	94	12	35	39-138	
Chloroform	10.0	9.15	92	1	35	72-127	
1,1,1-Trichloroethane	10.0	8.67	87	0	35	63-133	
Carbon tetrachloride	10.0	8.65	86	4	35	55-150	
Benzene	10.0	8.79	88	2	32	80-120	
1,2-Dichloroethane	10.0	9.38	94	9	32	68-132	
Trichloroethene	10.0	19.7	38	1	35	73-120	F1
1,2-Dichloropropane	10.0	9.80	98	11	34	76-124	
Bromodichloromethane	10.0	9.01	90	8	35	66-130	
cis-1,3-Dichloropropene	10.0	9.60	96	9	35	66-120	
4-Methyl-2-pentanone (MIBK)	20.0	18.0	90	7	35	45-145	
Toluene	10.0	9.38	94	2	35	80-123	
trans-1,3-Dichloropropene	10.0	11.3	113	3	35	65-125	
1,1,2-Trichloroethane	10.0	10.4	104	3	35	77-127	
Tetrachloroethene	10.0	14.2	68	7	35	70-135	F1
2-Hexanone	20.0	16.2	81	5	35	25-132	
Dibromochloromethane	10.0	9.15	91	1	35	60-140	
1,2-Dibromoethane (EDB)	10.0	10.2	102	4	35	74-123	
Chlorobenzene	10.0	9.90	99	1	29	80-120	
1,1,1,2-Tetrachloroethane	10.0	10.8	108	0	34	63-140	
Ethylbenzene	10.0	9.33	93	2	33	72-126	
Xylenes, Total	20.0	18.3	91	1	32	76-128	
Styrene	10.0	9.54	95	3	34	71-127	
Bromoform	10.0	8.52	85	3	35	46-150	
1,1,2,2-Tetrachloroethane	10.0	10.5	105	4	35	62-125	
1,4-Dioxane	200	191 J	95	9	35	10-160	

# Column to be used to flag recovery and RPD values

FORM IV  
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Lab File ID: 50415005.D Lab Sample ID: MB 180-138583/5  
Matrix: Water Heated Purge: (Y/N) N  
Instrument ID: CHHP5 Date Analyzed: 04/15/2015 14:33  
GC Column: DB-624 ID: 0.18 (mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
HD-MW-165-0/1-0	180-42975-1	50415006.D	04/15/2015 15:09
HD-QC2-0/1-2	180-42975-4	50415007.D	04/15/2015 15:33
	LCS 180-138583/8	50415008.D	04/15/2015 15:57
HD-MW-165-0/1-0 MS	180-42975-1 MS	50415009.D	04/15/2015 16:21
HD-MW-165-0/1-0 MSD	180-42975-1 MSD	50415010.D	04/15/2015 16:46
HD-MW-162-0/1-0	180-42975-2	50415020.D	04/15/2015 20:47

FORM IV  
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Lab File ID: 50416005.D Lab Sample ID: MB 180-138685/5  
Matrix: Water Heated Purge: (Y/N) N  
Instrument ID: CHHP5 Date Analyzed: 04/16/2015 11:32  
GC Column: DB-624 ID: 0.18 (mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
	LCS 180-138685/8	50416008.D	04/16/2015 13:05
	LCSD 180-138685/10	50416010.D	04/16/2015 13:53
HD-MW-162-0/1-0 DL	180-42975-2 DL	50416016.D	04/16/2015 16:29
HD-MW-169-0/1-0	180-42975-3	50416017.D	04/16/2015 16:53

FORM V  
GC/MS VOA INSTRUMENT PERFORMANCE CHECK  
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Lab File ID: 50316001.D BFB Injection Date: 03/16/2015  
Instrument ID: CHHP5 BFB Injection Time: 10:49  
Analysis Batch No.: 135593

M/E	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0 % of mass 95	22.7
75	30.0 - 60.0 % of mass 95	54.1
95	Base Peak, 100% relative abundance	100.0
96	5.0 - 9.0 % of mass 95	6.9
173	Less than 2.0 % of mass 174	0.8 (0.9)1
174	50.0 - 120.00 % of mass 95	85.5
175	5.0 - 9.0 % of mass 174	6.4 (7.5)1
176	95.0 - 101.0 % of mass 174	83.4 (97.4)1
177	5.0 - 9.0 % of mass 176	4.9 (5.8)2

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
	IC 180-135593/4	50316004.D	03/16/2015	12:41
	ICIS 180-135593/5	50316005.D	03/16/2015	13:05
	IC 180-135593/6	50316006.D	03/16/2015	13:29
	IC 180-135593/7	50316007.D	03/16/2015	13:53
	IC 180-135593/8	50316008.D	03/16/2015	14:17
	IC 180-135593/9	50316009.D	03/16/2015	14:41
	IC 180-135593/10	50316010.D	03/16/2015	15:05
	IC 180-135593/13	50316013.D	03/16/2015	16:17

FORM V  
GC/MS VOA INSTRUMENT PERFORMANCE CHECK  
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Lab File ID: 50415001.D BFB Injection Date: 04/15/2015  
Instrument ID: CHHP5 BFB Injection Time: 12:36  
Analysis Batch No.: 138583

M/E	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0 % of mass 95	23.1
75	30.0 - 60.0 % of mass 95	58.9
95	Base Peak, 100% relative abundance	100.0
96	5.0 - 9.0 % of mass 95	7.2
173	Less than 2.0 % of mass 174	0.6 (0.8)1
174	50.0 - 120.00 % of mass 95	81.5
175	5.0 - 9.0 % of mass 174	5.7 (7.0)1
176	95.0 - 101.0 % of mass 174	81.7 (100.2)1
177	5.0 - 9.0 % of mass 176	5.7 (6.9)2

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
	CCVIS 180-138583/2	50415002.D	04/15/2015	13:21
	CCV 180-138583/3	50415003.D	04/15/2015	13:45
	MB 180-138583/5	50415005.D	04/15/2015	14:33
HD-MW-165-0/1-0	180-42975-1	50415006.D	04/15/2015	15:09
HD-QC2-0/1-2	180-42975-4	50415007.D	04/15/2015	15:33
	LCS 180-138583/8	50415008.D	04/15/2015	15:57
HD-MW-165-0/1-0 MS	180-42975-1 MS	50415009.D	04/15/2015	16:21
HD-MW-165-0/1-0 MSD	180-42975-1 MSD	50415010.D	04/15/2015	16:46
HD-MW-162-0/1-0	180-42975-2	50415020.D	04/15/2015	20:47

FORM V  
GC/MS VOA INSTRUMENT PERFORMANCE CHECK  
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Lab File ID: 50416001.D BFB Injection Date: 04/16/2015  
Instrument ID: CHHP5 BFB Injection Time: 09:31  
Analysis Batch No.: 138685

M/E	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0 % of mass 95	20.9
75	30.0 - 60.0 % of mass 95	49.2
95	Base Peak, 100% relative abundance	100.0
96	5.0 - 9.0 % of mass 95	6.0
173	Less than 2.0 % of mass 174	0.5 (0.6)1
174	50.0 - 120.00 % of mass 95	81.0
175	5.0 - 9.0 % of mass 174	6.4 (7.9)1
176	95.0 - 101.0 % of mass 174	79.6 (98.3)1
177	5.0 - 9.0 % of mass 176	4.8 (6.0)2

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
	CCVIS 180-138685/2	50416002.D	04/16/2015	10:19
	MB 180-138685/5	50416005.D	04/16/2015	11:32
	LCS 180-138685/8	50416008.D	04/16/2015	13:05
	LCSD 180-138685/10	50416010.D	04/16/2015	13:53
HD-MW-162-0/1-0 DL	180-42975-2 DL	50416016.D	04/16/2015	16:29
HD-MW-169-0/1-0	180-42975-3	50416017.D	04/16/2015	16:53

FORM VIII  
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Sample No.: CCVIS 180-138583/2 Date Analyzed: 04/15/2015 13:21  
Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm)  
Lab File ID (Standard): 50415002.D Heated Purge: (Y/N) N  
Calibration ID: 22514

	TBA		FB		CBZ		
	AREA #	RT #	AREA #	RT #	AREA #	RT #	
12/24 HOUR STD	144875	4.31	546873	7.27	119760	10.36	
UPPER LIMIT	289750	4.81	1093746	7.77	239520	10.86	
LOWER LIMIT	72438	3.81	273437	6.77	59880	9.86	
LAB SAMPLE ID	CLIENT SAMPLE ID						
CCV 180-138583/3		158736	4.30	503999	7.27	110941	10.36
MB 180-138583/5		174312	4.31	460612	7.27	107539	10.36
180-42975-1	HD-MW-165-0/1-0	172081	4.29	443957	7.27	107798	10.36
180-42975-4	HD-QC2-0/1-2	167449	4.30	422314	7.27	101006	10.36
LCS 180-138583/8		163019	4.32	552275	7.27	123275	10.36
180-42975-1 MS	HD-MW-165-0/1-0 MS	171044	4.32	564634	7.27	124370	10.36
180-42975-1 MSD	HD-MW-165-0/1-0 MSD	208690	4.31	561630	7.27	131185	10.36
180-42975-2	HD-MW-162-0/1-0	168713	4.30	429853	7.28	110118	10.36

TBA = TBA-d9 (IS)

FB = Fluorobenzene (IS)

CBZ = Chlorobenzene-d5

Area Limit = 50%-200% of internal standard area

RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII  
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Sample No.: CCVIS 180-138583/2 Date Analyzed: 04/15/2015 13:21  
Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm)  
Lab File ID (Standard): 50415002.D Heated Purge: (Y/N) N  
Calibration ID: 22514

	DCB		AREA #	RT #	AREA #	RT #	AREA #	RT #
	AREA #	RT #						
12/24 HOUR STD	175806	12.68						
UPPER LIMIT	351612	13.18						
LOWER LIMIT	87903	12.18						
LAB SAMPLE ID	CLIENT SAMPLE ID							
CCV 180-138583/3		129092	12.68					
MB 180-138583/5		150393	12.68					
180-42975-1	HD-MW-165-0/1-0	154656	12.69					
180-42975-4	HD-QC2-0/1-2	149655	12.68					
LCS 180-138583/8		185676	12.68					
180-42975-1 MS	HD-MW-165-0/1-0 MS	179202	12.68					
180-42975-1 MSD	HD-MW-165-0/1-0 MSD	190234	12.68					
180-42975-2	HD-MW-162-0/1-0	142429	12.68					

DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area  
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII  
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Sample No.: CCVIS 180-138685/2 Date Analyzed: 04/16/2015 10:19  
Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm)  
Lab File ID (Standard): 50416002.D Heated Purge: (Y/N) N  
Calibration ID: 22514

	TBA		FB		CBZ	
	AREA #	RT #	AREA #	RT #	AREA #	RT #
12/24 HOUR STD	154447	4.32	501487	7.27	113593	10.36
UPPER LIMIT	308894	4.82	1002974	7.77	227186	10.86
LOWER LIMIT	77224	3.82	250744	6.77	56797	9.86
LAB SAMPLE ID	CLIENT SAMPLE ID					
MB 180-138685/5		194715	4.30	448391	7.28	108906
LCS 180-138685/8		158617	4.32	554542	7.27	125698
LCSD 180-138685/10		173426	4.32	571009	7.27	129134
180-42975-2 DL	HD-MW-162-0/1-0 DL	175233	4.30	442739	7.27	107268
180-42975-3	HD-MW-169-0/1-0	178588	4.30	432619	7.27	105105

TBA = TBA-d9 (IS)

FB = Fluorobenzene (IS)

CBZ = Chlorobenzene-d5

Area Limit = 50%-200% of internal standard area

RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII  
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Sample No.: CCVIS 180-138685/2 Date Analyzed: 04/16/2015 10:19  
Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm)  
Lab File ID (Standard): 50416002.D Heated Purge: (Y/N) N  
Calibration ID: 22514

	DCB		AREA #	RT #	AREA #	RT #	AREA #	RT #
	AREA #	RT #						
12/24 HOUR STD	167378	12.68						
UPPER LIMIT	334756	13.18						
LOWER LIMIT	83689	12.18						
LAB SAMPLE ID	CLIENT SAMPLE ID							
MB 180-138685/5		157942	12.69					
LCS 180-138685/8		186690	12.68					
LCSD 180-138685/10		189386	12.68					
180-42975-2 DL	HD-MW-162-0/1-0 DL	156120	12.68					
180-42975-3	HD-MW-169-0/1-0	144245	12.68					

DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area  
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: HD-MW-165-0/1-0 Lab Sample ID: 180-42975-1  
Matrix: Water Lab File ID: 50415006.D  
Analysis Method: 8260C Date Collected: 04/10/2015 08:40  
Sample wt/vol: 5 (mL) Date Analyzed: 04/15/2015 15:09  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138583 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
74-87-3	Chloromethane	1.0	U	1.0	0.28
75-01-4	Vinyl chloride	1.0	U	1.0	0.23
74-83-9	Bromomethane	1.0	U	1.0	0.31
75-00-3	Chloroethane	1.0	U	1.0	0.21
75-35-4	1,1-Dichloroethene	1.0	U	1.0	0.30
67-64-1	Acetone	5.0	U	5.0	2.5
75-15-0	Carbon disulfide	1.0	U F1 *	1.0	0.21
75-09-2	Methylene Chloride	1.0	U	1.0	0.13
156-60-5	trans-1,2-Dichloroethene	1.0	U	1.0	0.17
1634-04-4	Methyl tert-butyl ether	1.0	U	1.0	0.18
75-34-3	1,1-Dichloroethane	1.0	U	1.0	0.12
156-59-2	cis-1,2-Dichloroethene	1.0	U	1.0	0.24
74-97-5	Bromochloromethane	1.0	U	1.0	0.18
78-93-3	2-Butanone (MEK)	5.0	U	5.0	0.55
67-66-3	Chloroform	1.0	U	1.0	0.17
71-55-6	1,1,1-Trichloroethane	1.0	U	1.0	0.29
56-23-5	Carbon tetrachloride	1.0	U	1.0	0.14
71-43-2	Benzene	1.0	U	1.0	0.11
107-06-2	1,2-Dichloroethane	1.0	U	1.0	0.21
79-01-6	Trichloroethene	16	F1	1.0	0.14
78-87-5	1,2-Dichloropropane	1.0	U	1.0	0.095
75-27-4	Bromodichloromethane	1.0	U	1.0	0.13
10061-01-5	cis-1,3-Dichloropropene	1.0	U	1.0	0.19
108-10-1	4-Methyl-2-pentanone (MIBK)	5.0	U	5.0	0.53
108-88-3	Toluene	1.0	U	1.0	0.15
10061-02-6	trans-1,3-Dichloropropene	1.0	U	1.0	0.15
79-00-5	1,1,2-Trichloroethane	1.0	U	1.0	0.20
127-18-4	Tetrachloroethene	7.4	F1	1.0	0.15
591-78-6	2-Hexanone	5.0	U	5.0	0.16
124-48-1	Dibromochloromethane	1.0	U	1.0	0.14
106-93-4	1,2-Dibromoethane (EDB)	1.0	U	1.0	0.18
108-90-7	Chlorobenzene	1.0	U	1.0	0.14
630-20-6	1,1,1,2-Tetrachloroethane	1.0	U	1.0	0.28
100-41-4	Ethylbenzene	1.0	U	1.0	0.23
1330-20-7	Xylenes, Total	3.0	U	3.0	0.49

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: HD-MW-165-01/1-0 Lab Sample ID: 180-42975-1  
Matrix: Water Lab File ID: 50415006.D  
Analysis Method: 8260C Date Collected: 04/10/2015 08:40  
Sample wt/vol: 5 (mL) Date Analyzed: 04/15/2015 15:09  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138583 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
100-42-5	Styrene	1.0	U	1.0	0.097
75-25-2	Bromoform	1.0	U	1.0	0.19
79-34-5	1,1,2,2-Tetrachloroethane	1.0	U	1.0	0.20
107-13-1	Acrylonitrile	20	U	20	0.55
123-91-1	1,4-Dioxane	200	U	200	34

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	118		64-135
2037-26-5	Toluene-d8 (Surr)	100		71-118
460-00-4	4-Bromofluorobenzene (Surr)	96		70-118
1868-53-7	Dibromofluoromethane (Surr)	110		70-128

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415006.D  
 Lims ID: 180-42975-A-1 Lab Sample ID: 180-42975-1  
 Client ID: HD-MW-165-01-0  
 Sample Type: Client  
 Inject. Date: 15-Apr-2015 15:09:30 ALS Bottle#: 6 Worklist Smp#: 6  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: 180-42975-A-1  
 Misc. Info.: 180-0006480-006  
 Operator ID: 001562 Instrument ID: CHHP5  
 Method: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 07:33:03 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICAL File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 16-Apr-2015 07:33:03

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.290	4.302	-0.012	0	172081	1000.0	
* 2 Fluorobenzene (IS)	96	7.270	7.271	-0.001	98	443957	50.0	
* 3 Chlorobenzene-d5	119	10.361	10.361	0.000	88	107798	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.685	12.679	0.006	97	154656	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.528	6.528	0.000	93	111201	55.1	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.893	6.899	-0.006	0	156783	58.9	
\$ 7 Toluene-d8 (Surr)	98	8.919	8.919	0.000	94	431545	50.2	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.529	11.529	0.000	87	149192	48.2	
11 Dichlorodifluoromethane	85		1.619				ND	
12 Chloromethane	50		1.789				ND	
13 Vinyl chloride	62		1.917				ND	
14 Butadiene	39		1.959				ND	
15 Bromomethane	94		2.264				ND	
16 Chloroethane	64		2.416				ND	
17 Dichlorofluoromethane	67		2.671				ND	
18 Trichlorofluoromethane	101		2.726				ND	
19 Ethanol	45		3.018				ND	
20 Ethyl ether	59		3.091				ND	
21 Acrolein	56		3.261				ND	
22 1,1-Dichloroethene	96		3.395				ND	
23 1,1,2-Trichloro-1,2,2-trif	101		3.450				ND	
24 Acetone	43		3.499				ND	
25 Iodomethane	142		3.626				ND	
26 Carbon disulfide	76		3.675				ND	
27 Isopropyl alcohol	45		3.785				ND	
29 Acetonitrile	40		3.931				ND	
28 3-Chloro-1-propene	76		3.943				ND	
30 Methyl acetate	43		4.022				ND	
31 Methylene Chloride	84	4.150	4.143	0.007	10	1658	0.5600	M
32 2-Methyl-2-propanol	59		4.441				ND	
33 Acrylonitrile	53		4.551				ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ng	Flags
34 trans-1,2-Dichloroethene	96		4.563				ND	
35 Methyl tert-butyl ether	73		4.600				ND	
36 Hexane	57		4.983				ND	
37 1,1-Dichloroethane	63		5.165				ND	
38 Vinyl acetate	43		5.299				ND	
39 2-Chloro-1,3-butadiene	53		5.312				ND	
41 Isopropyl ether	45		5.324				ND	
40 Isopropyl ether TIC	45		5.409				ND	
42 Tert-butyl ethyl ether	59		5.799				ND	
44 2,2-Dichloropropane	77		5.926				ND	
45 cis-1,2-Dichloroethene	96	5.938	5.938	0.000	31	2849	1.02	
43 Tert-butyl ethyl ether (TI)	59		5.961				ND	
46 2-Butanone (MEK)	43		5.987				ND	
47 Propionitrile	54		6.060				ND	
48 Ethyl acetate	43		6.084				ND	
49 Chlorobromomethane	128		6.224				ND	
50 Methacrylonitrile	41		6.237				ND	
51 Tetrahydrofuran	42		6.279				ND	
52 Chloroform	83		6.339				ND	
53 1,1,1-Trichloroethane	97		6.528				ND	
54 Cyclohexane	56		6.583				ND	
56 Carbon tetrachloride	117		6.723				ND	
55 1,1-Dichloropropene	75		6.723				ND	
57 Isobutyl alcohol	41		6.942				ND	
58 Benzene	78		6.954				ND	
59 1,2-Dichloroethane	62		6.984				ND	
61 Tert-amyl methyl ether	73		7.113				ND	
60 Tert-amyl methyl ether (TI)	73	7.331	7.262	0.069	1	187	0.0211	
62 n-Heptane	43		7.276				ND	
63 n-Butanol	56		7.654				ND	
64 Trichloroethene	130	7.666	7.666	0.000	97	209048	79.3	
65 Ethyl acrylate	55		7.818				ND	
66 Methylcyclohexane	83		7.860				ND	
67 1,2-Dichloropropane	63		7.897				ND	
68 Dibromomethane	93		8.025				ND	
70 1,4-Dioxane	88		8.049				ND	
69 Methyl methacrylate	69		8.049				ND	
71 Dichlorobromomethane	83		8.195				ND	
72 2-Nitropropane	41		8.439				ND	
73 2-Chloroethyl vinyl ether	63		8.517				ND	
74 cis-1,3-Dichloropropene	75		8.651				ND	
75 4-Methyl-2-pentanone (MIBK)	43		8.822				ND	
76 Toluene	91		8.986				ND	
77 trans-1,3-Dichloropropene	75		9.217				ND	
78 Ethyl methacrylate	69		9.314				ND	
79 1,1,2-Trichloroethane	97		9.399				ND	
80 Tetrachloroethene	164	9.533	9.533	0.000	96	79762	36.9	
81 1,3-Dichloropropane	76		9.558				ND	
82 2-Hexanone	43		9.655				ND	
83 n-Butyl acetate	43		9.783				ND	
84 Chlorodibromomethane	129		9.789				ND	
85 Ethylene Dibromide	107		9.898				ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ng	Flags
86 3-Chlorobenzotrifluoride	180		10.373				ND	
87 Chlorobenzene	112		10.391				ND	
88 4-Chlorobenzotrifluoride	180		10.428				ND	
89 1,1,1,2-Tetrachloroethane	131		10.470				ND	
90 Ethylbenzene	106		10.501				ND	
91 m-Xylene & p-Xylene	106		10.616				ND	
92 o-Xylene	106		11.012				ND	
93 Styrene	104		11.024				ND	
94 Bromoform	173		11.212				ND	
95 Cyclohexanol	57		11.231				ND	
96 2-Chlorobenzotrifluoride	180		11.273				ND	
97 Isopropylbenzene	105		11.377				ND	
98 Cyclohexanone	55		11.481				ND	
99 1,1,2,2-Tetrachloroethane	83		11.675				ND	
100 Bromobenzene	156		11.681				ND	
101 1,2,3-Trichloropropane	110		11.717				ND	
102 trans-1,4-Dichloro-2-buten	53		11.729				ND	
103 N-Propylbenzene	120		11.784				ND	
104 2-Chlorotoluene	126		11.869				ND	
105 3-Chlorotoluene	126		11.930				ND	
106 1,3,5-Trimethylbenzene	105		11.961				ND	
107 4-Chlorotoluene	126		11.985				ND	
108 tert-Butylbenzene	119		12.283				ND	
109 Pentachloroethane	167		12.302				ND	
110 1,2,4-Trimethylbenzene	105		12.332				ND	
111 1,2-dichloro-4-(trifluoromethyl)	214		12.399				ND	
112 sec-Butylbenzene	105		12.502				ND	
113 1,3-Dichlorobenzene	146		12.618				ND	
114 4-Isopropyltoluene	119		12.648				ND	
115 1,4-Dichlorobenzene	146	12.709	12.703	0.006	17	1049	0.1948	
116 2,4-Dichloro-1-(trifluoromethyl)	214		12.758				ND	
117 1,2,3-Trimethylbenzene	105		12.758				ND	
118 2,5-Dichlorobenzotrifluoride	214		12.806				ND	
119 Benzyl chloride	91		12.843				ND	
120 n-Butylbenzene	91		13.062				ND	
121 1,2-Dichlorobenzene	146		13.080				ND	
122 1,2-Dibromo-3-Chloropropan	75		13.853				ND	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125		14.005				ND	
124 1,3,5-Trichlorobenzene	180		14.072				ND	
125 2,3- & 3,4- Dichlorotoluene	125		14.424				ND	
126 1,2,4-Trichlorobenzene	180		14.686				ND	
127 Hexachlorobutadiene	225		14.856				ND	
128 Naphthalene	128		14.942				ND	
129 1,2,3-Trichlorobenzene	180		15.185				ND	
131 2,4,5-Trichlorotoluene	159		15.964				ND	
130 2,3,6-Trichlorotoluene	159		16.061				ND	
132 2-Methylnaphthalene	142		16.074				ND	
151 Isooctane	57		0.000				ND	
149 3,4-Dichlorotoluene	1		0.000				ND	
148 2,3-Dichlorotoluene	1		0.000				ND	
147 2,4-Dichlorotoluene	1		0.000				ND	
152 Formaldehyde TIC	1		0.000				ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ng	Flags
146 2,5-Dichlorotoluene	1		0.000				ND	
150 2,6-Dichlorotoluene	1		0.000				ND	
S 133 Xylenes, Total	106		1.000				ND	
S 134 1,2-Dichloroethene, Total	96			0			1.02	
S 135 1,3-Dichloropropene, Total	1		0.000				ND	
T 153 1,2 Epoxybutane TIC	42		0.000				ND	
T 136 Mesityl oxide TIC	83		0.000				ND	
T 137 Tetrahydrofuran TIC	42		0.000				ND	
T 138 Methyl n-amyl ketone TIC	43		0.000				ND	

**QC Flag Legend**

Review Flags

M - Manually Integrated

**Reagents:**

VOA8260INT\_00031  
VOA8260SURR\_00033

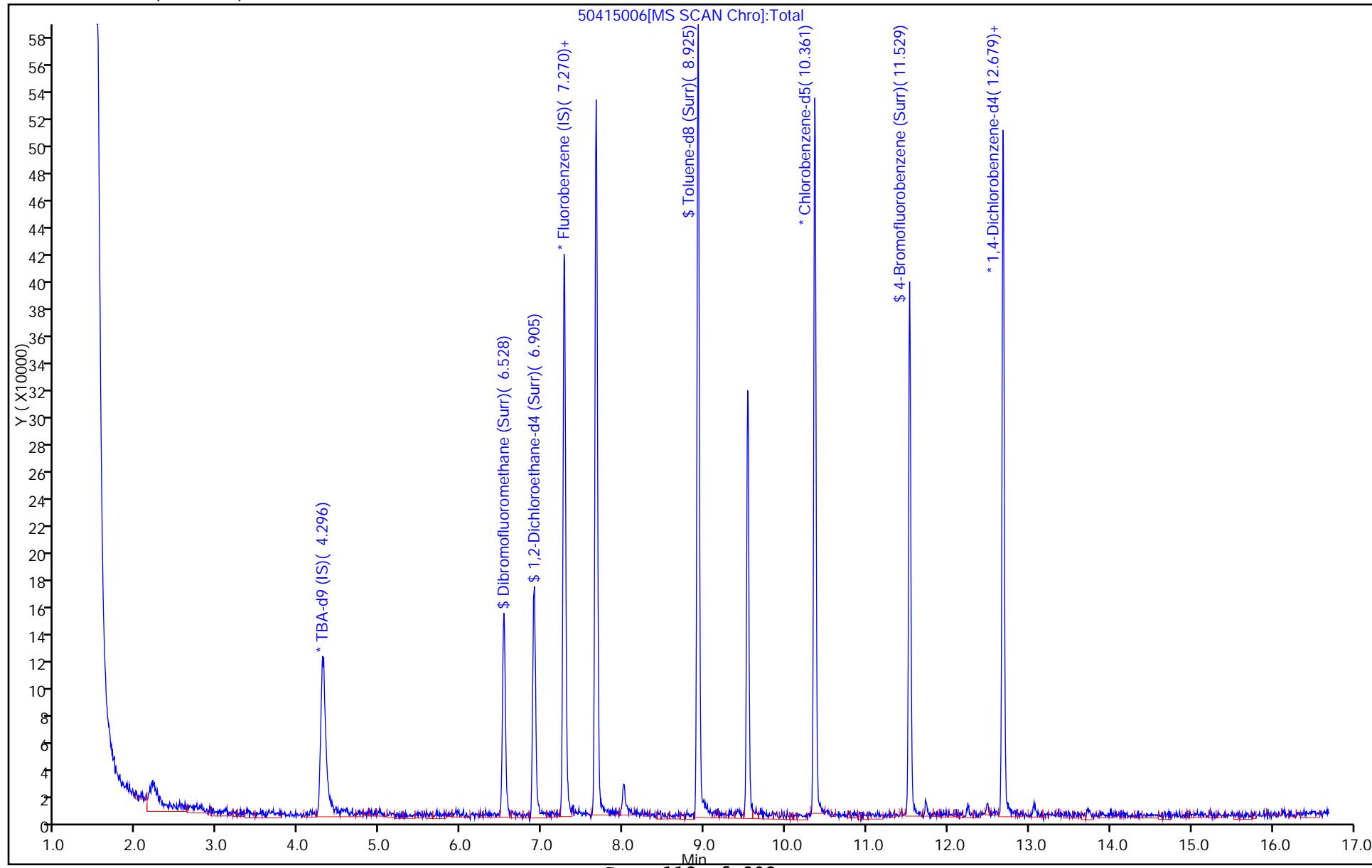
Amount Added: 2.00 Units: uL Run Reagent  
Amount Added: 2.00 Units: uL Run Reagent

Report Date: 16-Apr-2015 07:33:04

Chrom Revision: 2.2 13-Mar-2015 11:20:44

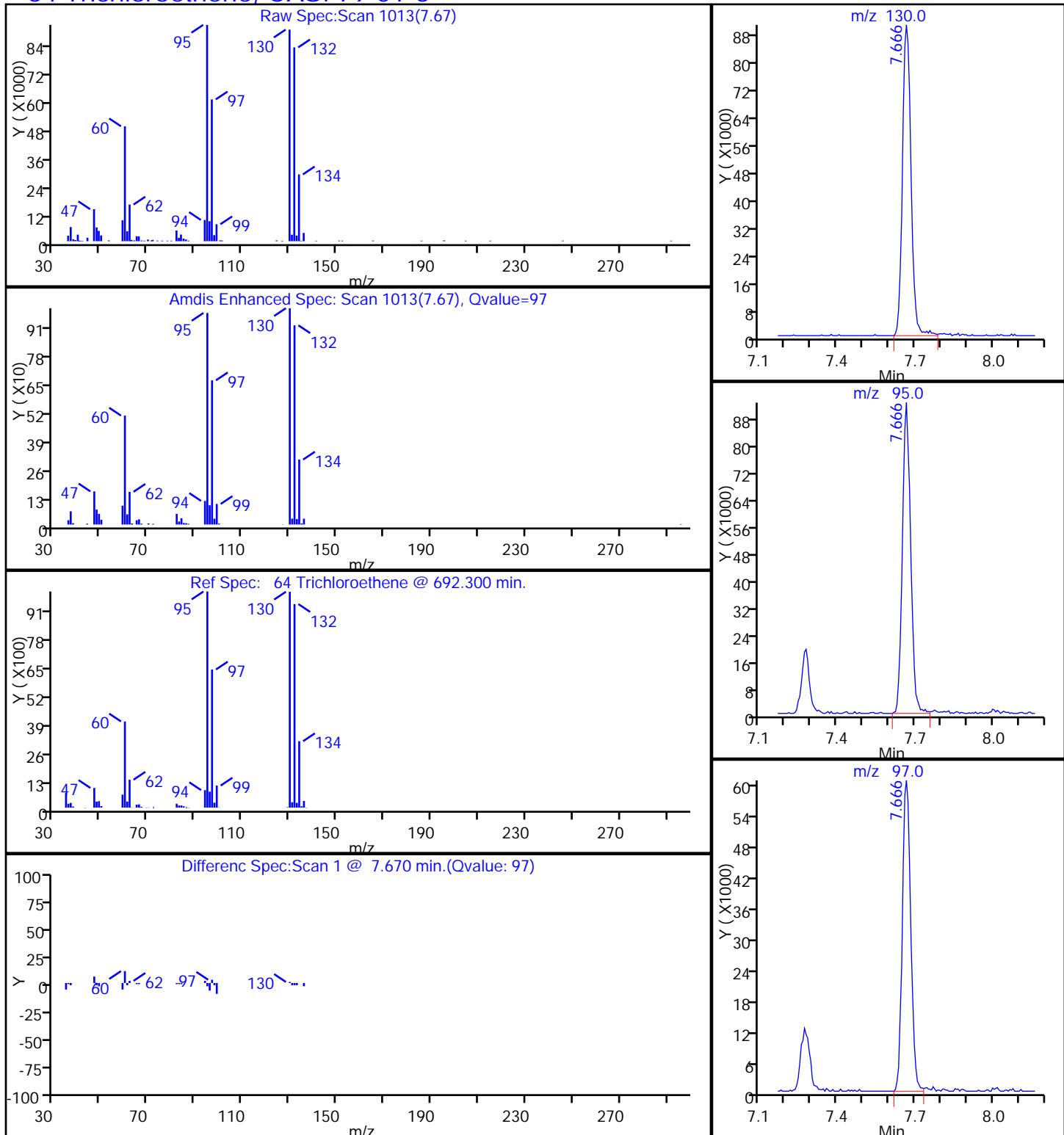
## TestAmerica Pittsburgh

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150415-6480.b\\50415006.D  
Injection Date: 15-Apr-2015 15:09:30 Instrument ID: CHHP5 Operator ID: 001562  
Lims ID: 180-42975-A-1 Lab Sample ID: 180-42975-1 Worklist Smp#: 6  
Client ID: HD-MW-165-0/1-0  
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 6  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)



## TestAmerica Pittsburgh

Data File: \PITCHROM\ChromData\CHHP5\20150415-6480.b\50415006.D  
 Injection Date: 15-Apr-2015 15:09:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-A-1 Lab Sample ID: 180-42975-1  
 Client ID: HD-MW-165-0/1-0  
 Operator ID: 001562 ALS Bottle#: 6 Worklist Smp#: 6  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 ( 0.18 mm) Detector: MS SCAN

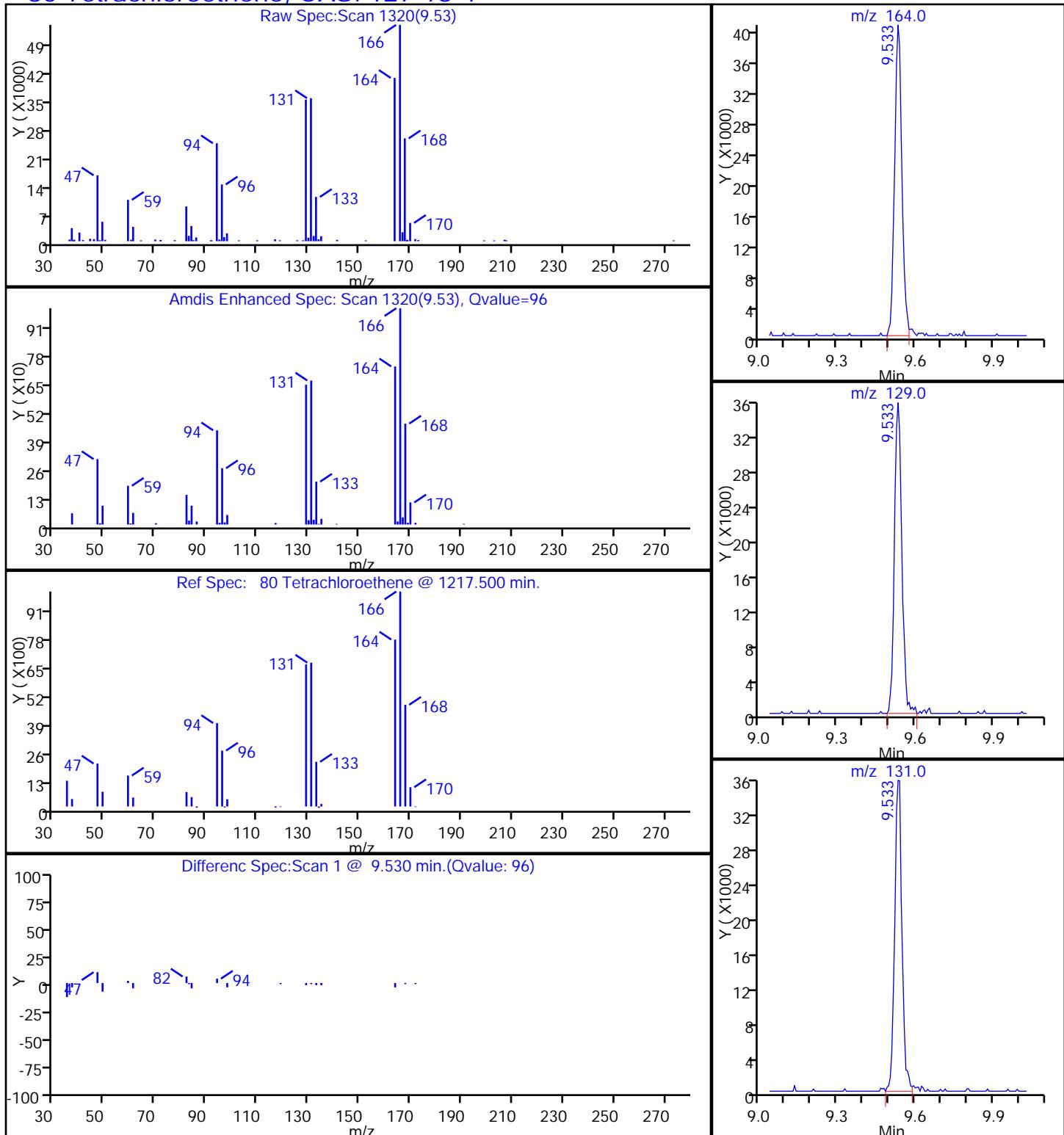
**64 Trichloroethene, CAS: 79-01-6**

## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415006.D  
 Injection Date: 15-Apr-2015 15:09:30  
 Lims ID: 180-42975-A-1  
 Client ID: HD-MW-165-0/1-0  
 Operator ID: 001562  
 Purge Vol: 5.000 mL  
 Method: MSVOA\_LL\_CHHP5  
 Column: DB-624 ( 0.18 mm)

Instrument ID:	CHHP5
Lab Sample ID:	180-42975-1
ALS Bottle#:	6
Dil. Factor:	1.0000
Limit Group:	VOA 8260C ICAL
Detector	MS SCAN
Worklist Smp#:	6

## 80 Tetrachloroethene, CAS: 127-18-4



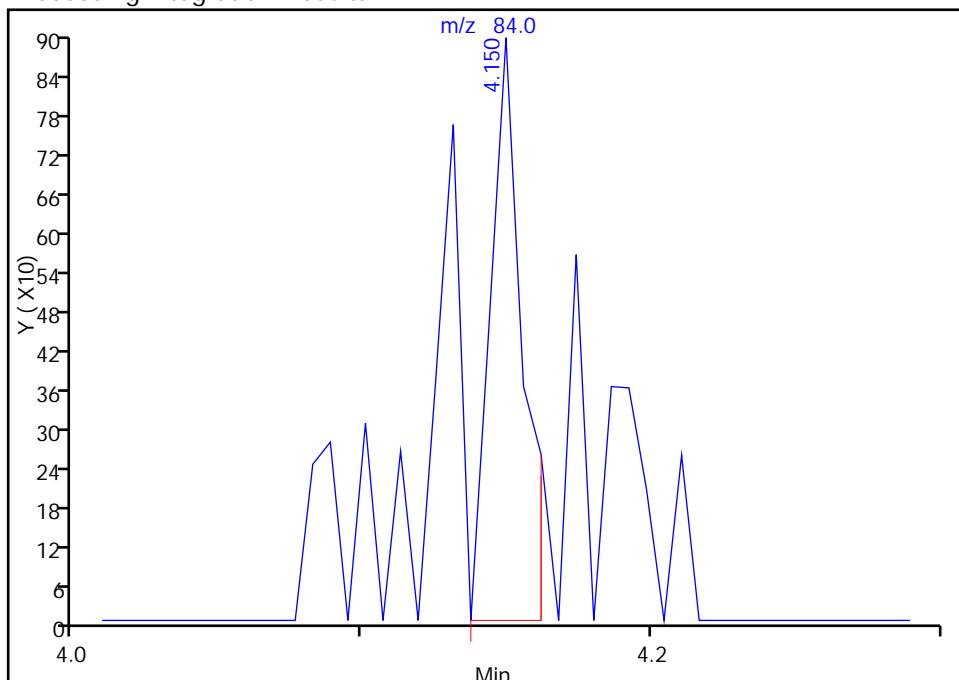
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415006.D  
 Injection Date: 15-Apr-2015 15:09:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-A-1 Lab Sample ID: 180-42975-1  
 Client ID: HD-MW-165-0/1-0  
 Operator ID: 001562 ALS Bottle#: 6 Worklist Smp#: 6  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 31 Methylene Chloride, CAS: 75-09-2

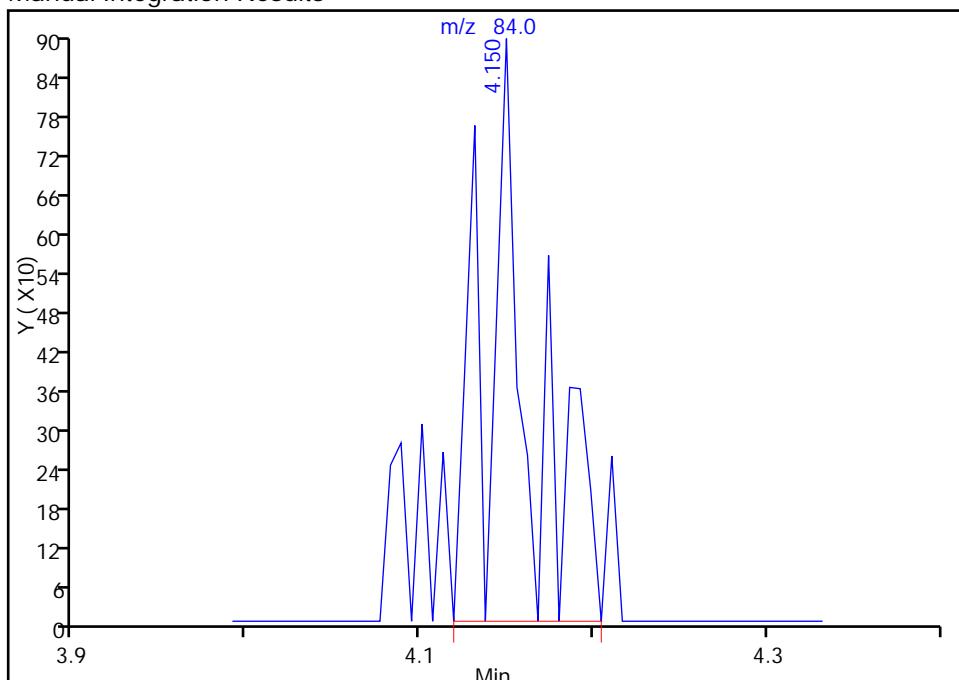
RT: 4.15  
 Area: 708  
 Amount: 0.239128  
 Amount Units: ng

## Processing Integration Results



RT: 4.15  
 Area: 1658  
 Amount: 0.559991  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 16-Apr-2015 07:33:03

Audit Action: Manually Integrated

Audit Reason: Split Peak

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: HD-MW-162-0/1-0 Lab Sample ID: 180-42975-2  
Matrix: Water Lab File ID: 50415020.D  
Analysis Method: 8260C Date Collected: 04/10/2015 09:50  
Sample wt/vol: 5 (mL) Date Analyzed: 04/15/2015 20:47  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 4  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138583 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
74-87-3	Chloromethane	4.0	U	4.0	1.1
75-01-4	Vinyl chloride	4.0	U	4.0	0.91
74-83-9	Bromomethane	4.0	U	4.0	1.3
75-00-3	Chloroethane	4.0	U	4.0	0.86
75-35-4	1,1-Dichloroethene	1.2	J	4.0	1.2
67-64-1	Acetone	20	U	20	10
75-15-0	Carbon disulfide	4.0	U *	4.0	0.85
75-09-2	Methylene Chloride	1.4	J	4.0	0.50
156-60-5	trans-1,2-Dichloroethene	4.0	U	4.0	0.68
1634-04-4	Methyl tert-butyl ether	4.0	U	4.0	0.73
75-34-3	1,1-Dichloroethane	4.0	U	4.0	0.47
156-59-2	cis-1,2-Dichloroethene	4.0	U	4.0	0.95
74-97-5	Bromochloromethane	4.0	U	4.0	0.72
78-93-3	2-Butanone (MEK)	20	U	20	2.2
67-66-3	Chloroform	4.0	U	4.0	0.68
71-55-6	1,1,1-Trichloroethane	4.0	U	4.0	1.1
56-23-5	Carbon tetrachloride	4.0	U	4.0	0.55
71-43-2	Benzene	4.0	U	4.0	0.42
107-06-2	1,2-Dichloroethane	4.0	U	4.0	0.85
79-01-6	Trichloroethene	190		4.0	0.57
78-87-5	1,2-Dichloropropane	4.0	U	4.0	0.38
75-27-4	Bromodichloromethane	4.0	U	4.0	0.52
10061-01-5	cis-1,3-Dichloropropene	4.0	U	4.0	0.75
108-10-1	4-Methyl-2-pentanone (MIBK)	20	U	20	2.1
108-88-3	Toluene	4.0	U	4.0	0.60
10061-02-6	trans-1,3-Dichloropropene	4.0	U	4.0	0.59
79-00-5	1,1,2-Trichloroethane	4.0	U	4.0	0.81
127-18-4	Tetrachloroethene	560	E	4.0	0.59
591-78-6	2-Hexanone	20	U	20	0.64
124-48-1	Dibromochloromethane	4.0	U	4.0	0.55
106-93-4	1,2-Dibromoethane (EDB)	4.0	U	4.0	0.72
108-90-7	Chlorobenzene	4.0	U	4.0	0.54
630-20-6	1,1,1,2-Tetrachloroethane	4.0	U	4.0	1.1
100-41-4	Ethylbenzene	4.0	U	4.0	0.91
1330-20-7	Xylenes, Total	12	U	12	2.0
100-42-5	Styrene	4.0	U	4.0	0.39

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: HD-MW-162-0/1-0 Lab Sample ID: 180-42975-2  
Matrix: Water Lab File ID: 50415020.D  
Analysis Method: 8260C Date Collected: 04/10/2015 09:50  
Sample wt/vol: 5 (mL) Date Analyzed: 04/15/2015 20:47  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 4  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138583 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-25-2	Bromoform	4.0	U	4.0	0.77
79-34-5	1,1,2,2-Tetrachloroethane	4.0	U	4.0	0.80
107-13-1	Acrylonitrile	80	U	80	2.2
123-91-1	1,4-Dioxane	800	U	800	140

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	115		64-135
2037-26-5	Toluene-d8 (Surr)	98		71-118
460-00-4	4-Bromofluorobenzene (Surr)	91		70-118
1868-53-7	Dibromofluoromethane (Surr)	116		70-128

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415020.D  
 Lims ID: 180-42975-A-2 Lab Sample ID: 180-42975-2  
 Client ID: HD-MW-162-01-0  
 Sample Type: Client  
 Inject. Date: 15-Apr-2015 20:47:30 ALS Bottle#: 20 Worklist Smp#: 20  
 Purge Vol: 5.000 mL Dil. Factor: 4.0000  
 Sample Info: 180-42975-A-2, 4x  
 Misc. Info.: 180-0006480-020  
 Operator ID: 001562 Instrument ID: CHHP5  
 Method: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 08:55:53 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICAL File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 16-Apr-2015 08:55:53

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.295	4.302	-0.007	0	168713	1000.0	
* 2 Fluorobenzene (IS)	96	7.276	7.271	0.005	98	429853	50.0	
* 3 Chlorobenzene-d5	119	10.360	10.361	-0.001	88	110118	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.684	12.679	0.005	97	142429	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.534	6.528	0.006	93	113014	57.8	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.905	6.899	0.006	0	148490	57.6	
\$ 7 Toluene-d8 (Surr)	98	8.925	8.919	0.006	94	429410	48.9	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.528	11.529	-0.001	87	143361	45.3	
12 Chloromethane	50		1.789				ND	
13 Vinyl chloride	62		1.917				ND	
15 Bromomethane	94		2.264				ND	
16 Chloroethane	64		2.416				ND	
22 1,1-Dichloroethene	96	3.401	3.395	0.006	39	3707	1.50	M
24 Acetone	43	3.517	3.499	0.017	66	3429	3.89	
26 Carbon disulfide	76		3.675				ND	
31 Methylene Chloride	84	4.149	4.143	0.006	76	4853	1.69	M
33 Acrylonitrile	53		4.551				ND	
34 trans-1,2-Dichloroethene	96		4.563				ND	
35 Methyl tert-butyl ether	73		4.600				ND	
37 1,1-Dichloroethane	63		5.165				ND	
45 cis-1,2-Dichloroethene	96		5.938				ND	
46 2-Butanone (MEK)	43		5.987				ND	
49 Chlorobromomethane	128		6.224				ND	
52 Chloroform	83	6.339	6.339	0.000	26	2037	0.4900	
53 1,1,1-Trichloroethane	97	6.528	6.528	0.000	1	580	0.2185	M
56 Carbon tetrachloride	117		6.723				ND	
58 Benzene	78		6.954				ND	
59 1,2-Dichloroethane	62		6.984				ND	
64 Trichloroethene	130	7.665	7.666	-0.001	97	590933	231.5	
67 1,2-Dichloropropane	63		7.897				ND	
70 1,4-Dioxane	88		8.049				ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ng	Flags
71 Dichlorobromomethane	83		8.195				ND	
74 cis-1,3-Dichloropropene	75		8.651				ND	
75 4-Methyl-2-pentanone (MIBK)	43		8.822				ND	
76 Toluene	91		8.986				ND	
77 trans-1,3-Dichloropropene	75		9.217				ND	
79 1,1,2-Trichloroethane	97		9.399				ND	
80 Tetrachloroethene	164	9.533	9.533	0.000	93	1535152	695.5	E
82 2-Hexanone	43		9.655				ND	
84 Chlorodibromomethane	129		9.789				ND	
85 Ethylene Dibromide	107		9.898				ND	
87 Chlorobenzene	112		10.391				ND	
89 1,1,1,2-Tetrachloroethane	131		10.470				ND	
90 Ethylbenzene	106		10.501				ND	
91 m-Xylene & p-Xylene	106		10.616				ND	
92 o-Xylene	106		11.012				ND	
93 Styrene	104		11.024				ND	
94 Bromoform	173		11.212				ND	
99 1,1,2,2-Tetrachloroethane	83		11.675				ND	
S 133 Xylenes, Total	106		1.000				ND	

**QC Flag Legend**

Processing Flags

E - Exceeded Maximum Amount

Review Flags

M - Manually Integrated

**Reagents:**

VOA8260INT\_00031

Amount Added: 2.00

Units: uL

Run Reagent

VOA8260SURR\_00033

Amount Added: 2.00

Units: uL

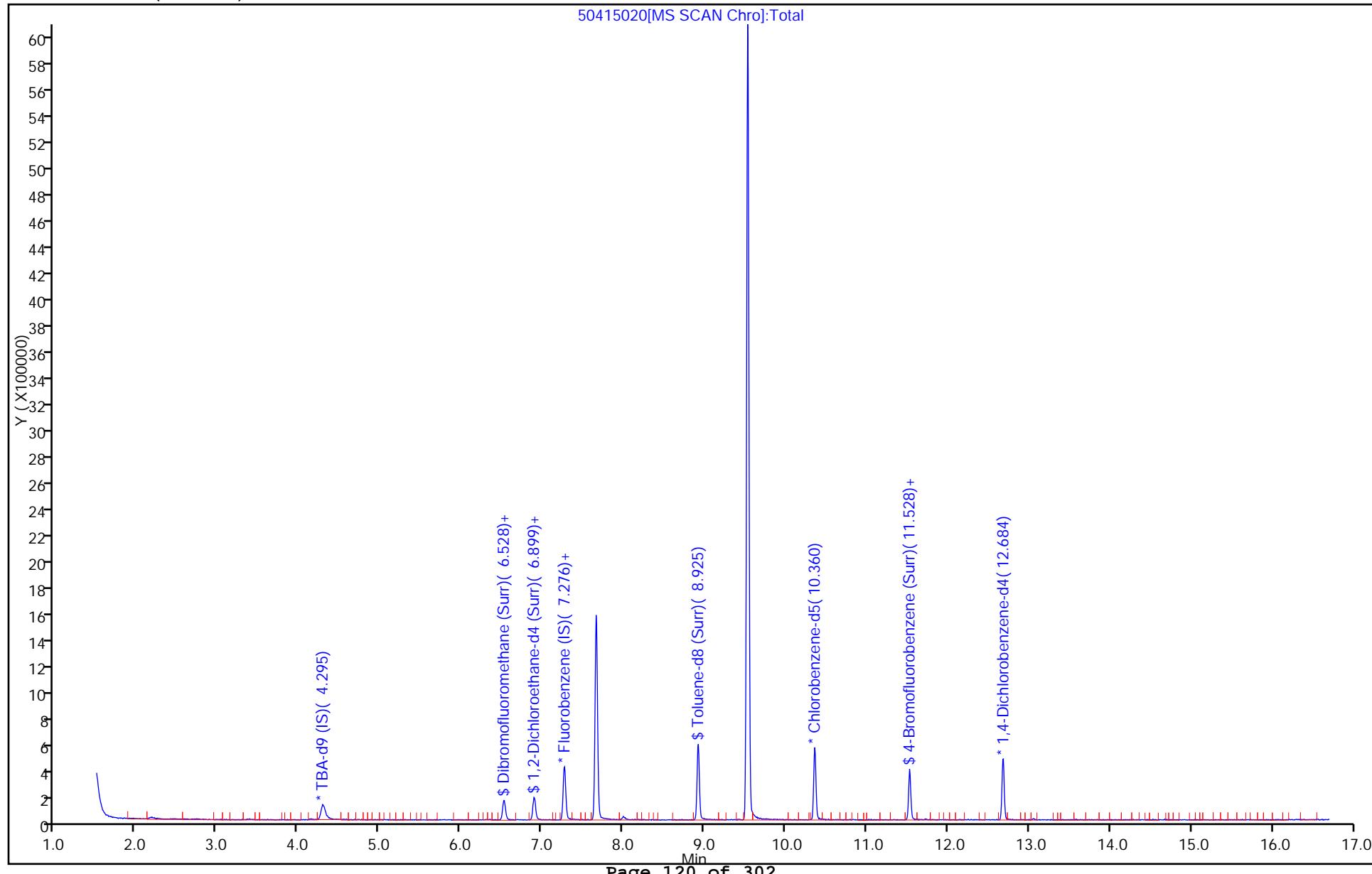
Run Reagent

Report Date: 16-Apr-2015 08:55:54

Chrom Revision: 2.2 13-Mar-2015 11:20:44

## TestAmerica Pittsburgh

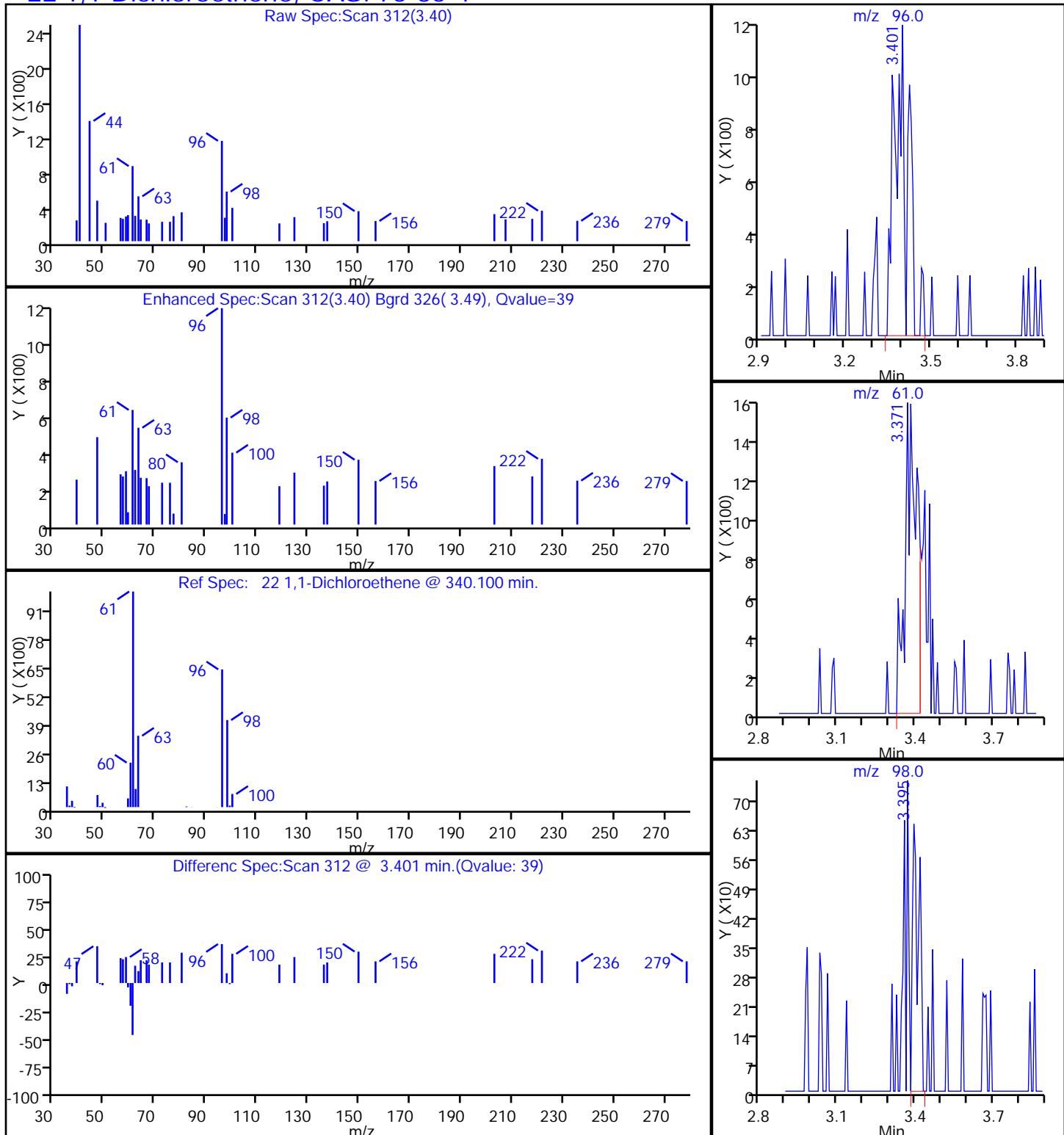
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Injection Date: 15-Apr-2015 20:47:30 Instrument ID: CHHP5 Operator ID: 001562  
Lims ID: 180-42975-A-2 Lab Sample ID: 180-42975-2 Worklist Smp#: 20  
Client ID: HD-MW-162-0/1-0  
Purge Vol: 5.000 mL Dil. Factor: 4.0000 ALS Bottle#: 20  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)



## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415020.D  
 Injection Date: 15-Apr-2015 20:47:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-A-2 Lab Sample ID: 180-42975-2  
 Client ID: HD-MW-162-0/1-0  
 Operator ID: 001562 ALS Bottle#: 20 Worklist Smp#: 20  
 Purge Vol: 5.000 mL Dil. Factor: 4.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 ( 0.18 mm) Detector: MS SCAN

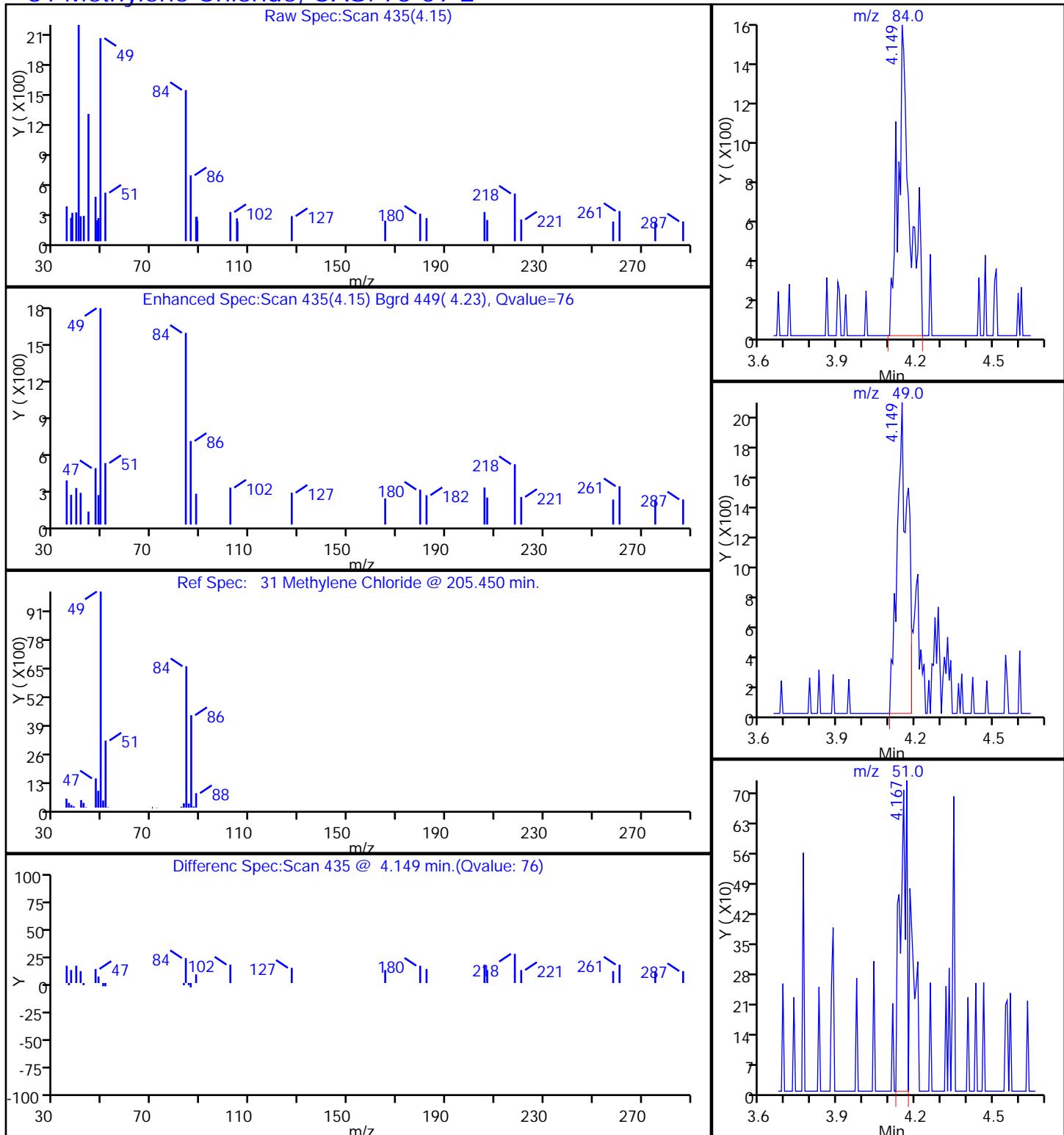
## 22 1,1-Dichloroethene, CAS: 75-35-4



## TestAmerica Pittsburgh

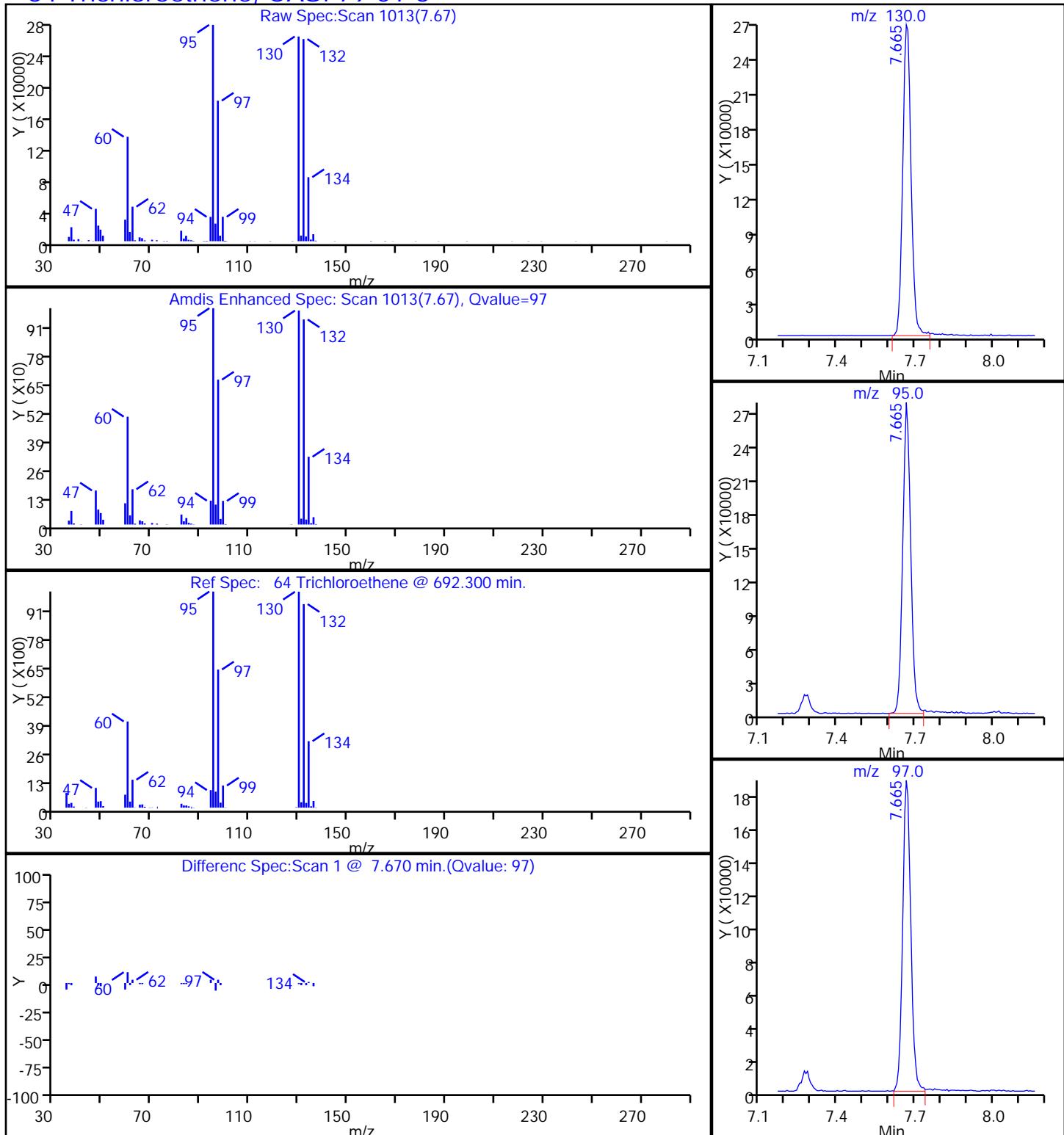
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 Injection Date: 15-Apr-2015 20:47:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-A-2 Lab Sample ID: 180-42975-2  
 Client ID: HD-MW-162-0/1-0  
 Operator ID: 001562 ALS Bottle#: 20 Worklist Smp#: 20  
 Purge Vol: 5.000 mL Dil. Factor: 4.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 ( 0.18 mm) Detector: MS SCAN

## 31 Methylene Chloride, CAS: 75-09-2



## TestAmerica Pittsburgh

Data File: \PITCHROM\ChromData\CHHP5\20150415-6480.b\50415020.D  
 Injection Date: 15-Apr-2015 20:47:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-A-2 Lab Sample ID: 180-42975-2  
 Client ID: HD-MW-162-0/1-0  
 Operator ID: 001562 ALS Bottle#: 20 Worklist Smp#: 20  
 Purge Vol: 5.000 mL Dil. Factor: 4.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 ( 0.18 mm) Detector: MS SCAN

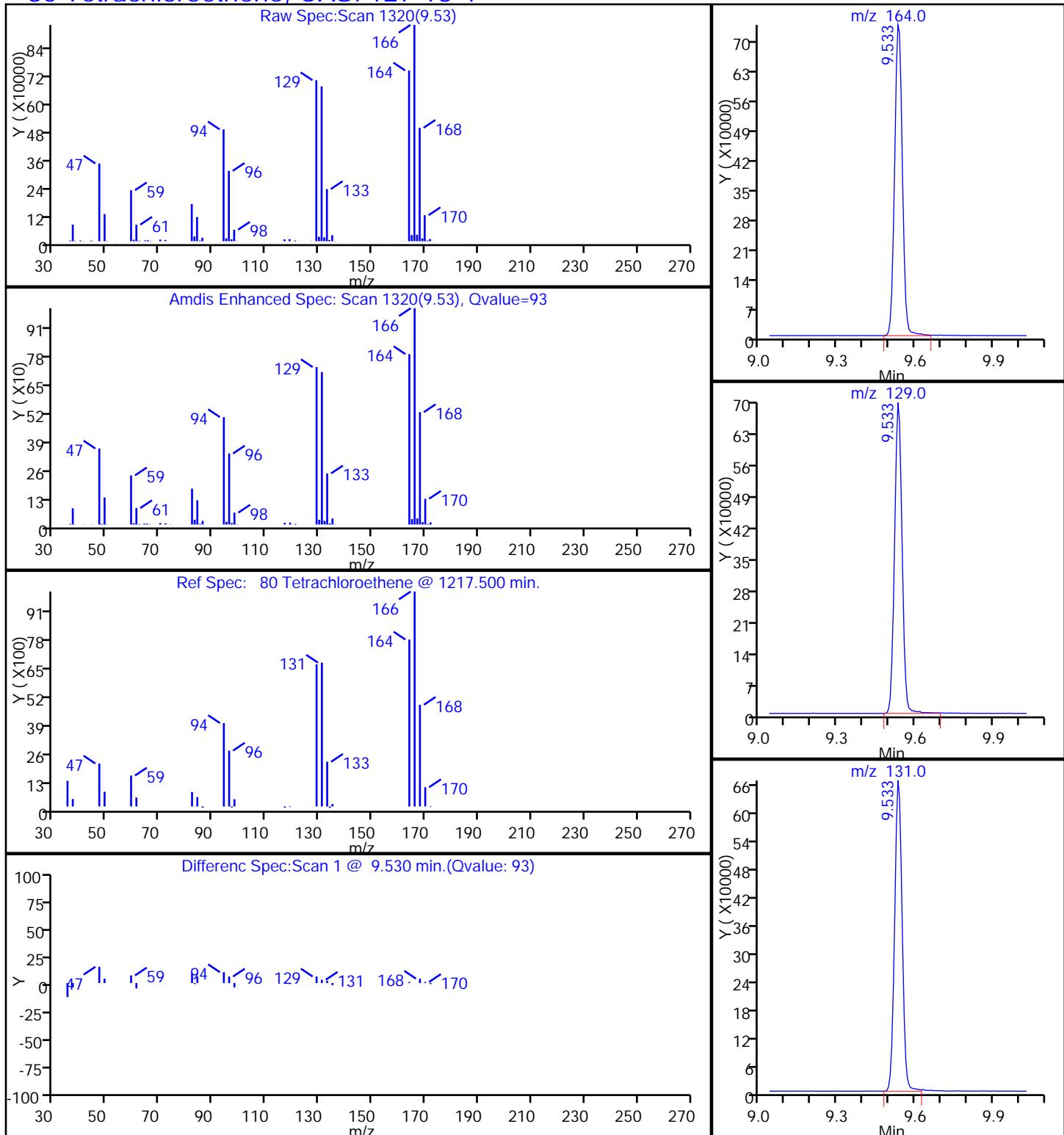
**64 Trichloroethene, CAS: 79-01-6**

## TestAmerica Pittsburgh

Data File: \PITCHROM\ChromData\CHHP5\20150415-6480.b\50415020.D  
 Injection Date: 15-Apr-2015 20:47:30  
 Lims ID: 180-42975-A-2  
 Client ID: HD-MW-162-0/1-0  
 Operator ID: 001562  
 Purge Vol: 5.000 mL  
 Method: MSVOA\_LL\_CHHP5  
 Column: DB-624 ( 0.18 mm)

Instrument ID:	CHHP5
Lab Sample ID:	180-42975-2
ALS Bottle#:	20
Dil. Factor:	4.0000
Limit Group:	VOA 8260C ICAL
Detector	MS SCAN
Worklist Smp#:	20

## 80 Tetrachloroethene, CAS: 127-18-4



## TestAmerica Pittsburgh

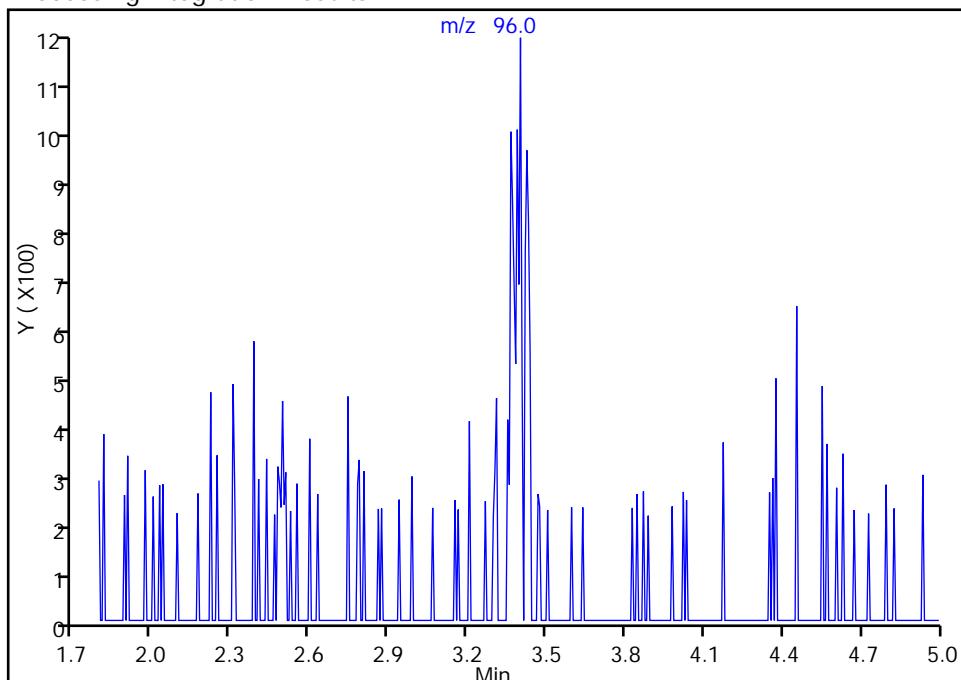
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 Injection Date: 15-Apr-2015 20:47:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-A-2 Lab Sample ID: 180-42975-2  
 Client ID: HD-MW-162-0/1-0  
 Operator ID: 001562 ALS Bottle#: 20 Worklist Smp#: 20  
 Purge Vol: 5.000 mL Dil. Factor: 4.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 22 1,1-Dichloroethene, CAS: 75-35-4

Not Detected

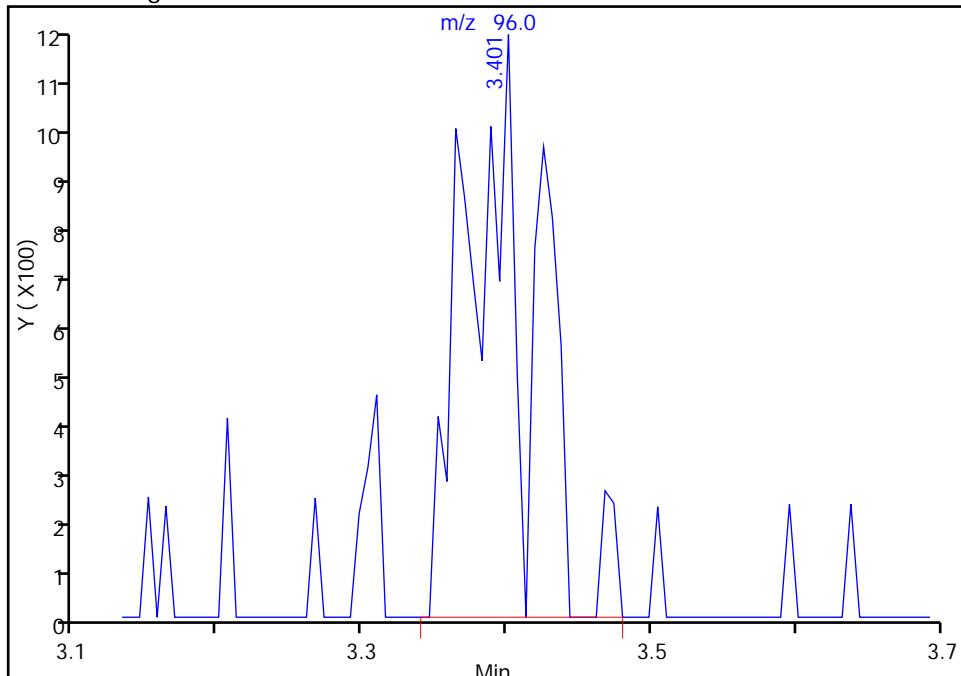
Expected RT: 3.40

## Processing Integration Results



## Manual Integration Results

RT: 3.40  
 Area: 3707  
 Amount: 1.495495  
 Amount Units: ng



Reviewer: fergusond, 16-Apr-2015 08:55:53

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

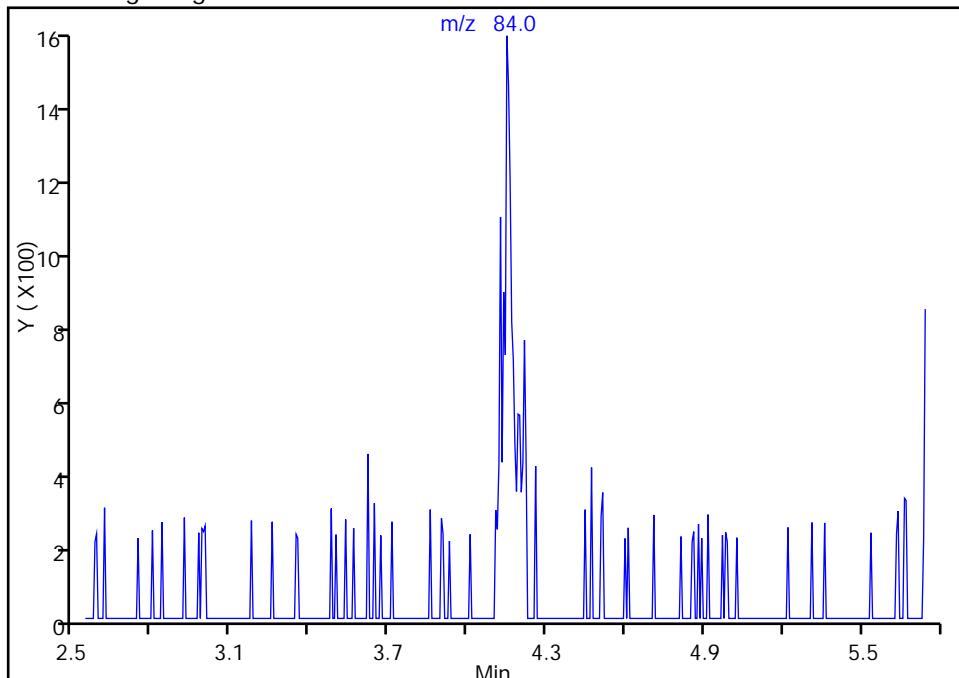
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415020.D  
 Injection Date: 15-Apr-2015 20:47:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-A-2 Lab Sample ID: 180-42975-2  
 Client ID: HD-MW-162-0/1-0  
 Operator ID: 001562 ALS Bottle#: 20 Worklist Smp#: 20  
 Purge Vol: 5.000 mL Dil. Factor: 4.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 31 Methylene Chloride, CAS: 75-09-2

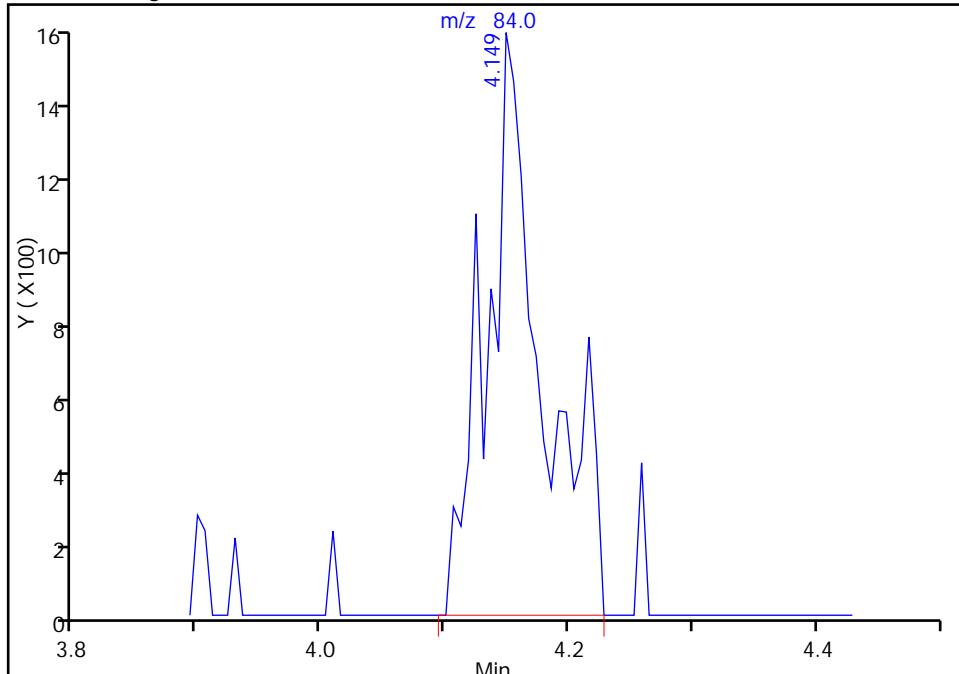
Not Detected  
 Expected RT: 4.14

## Processing Integration Results



## Manual Integration Results

RT: 4.15  
 Area: 4853  
 Amount: 1.692887  
 Amount Units: ng



Reviewer: fergusond, 16-Apr-2015 08:55:53

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

## TestAmerica Pittsburgh

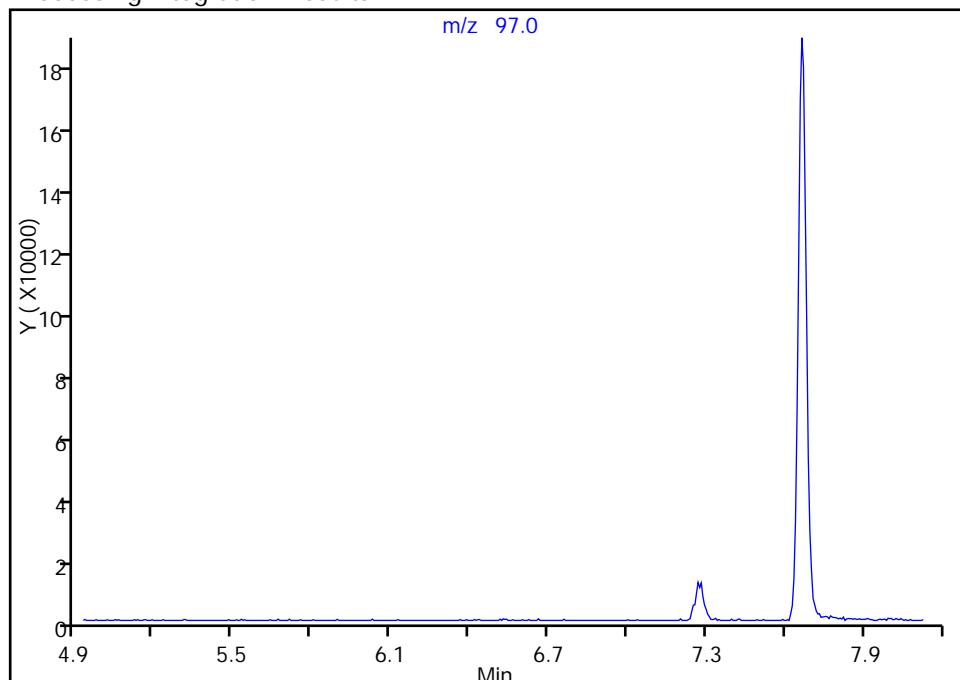
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 Injection Date: 15-Apr-2015 20:47:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-A-2 Lab Sample ID: 180-42975-2  
 Client ID: HD-MW-162-0/1-0  
 Operator ID: 001562 ALS Bottle#: 20 Worklist Smp#: 20  
 Purge Vol: 5.000 mL Dil. Factor: 4.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 53 1,1,1-Trichloroethane, CAS: 71-55-6

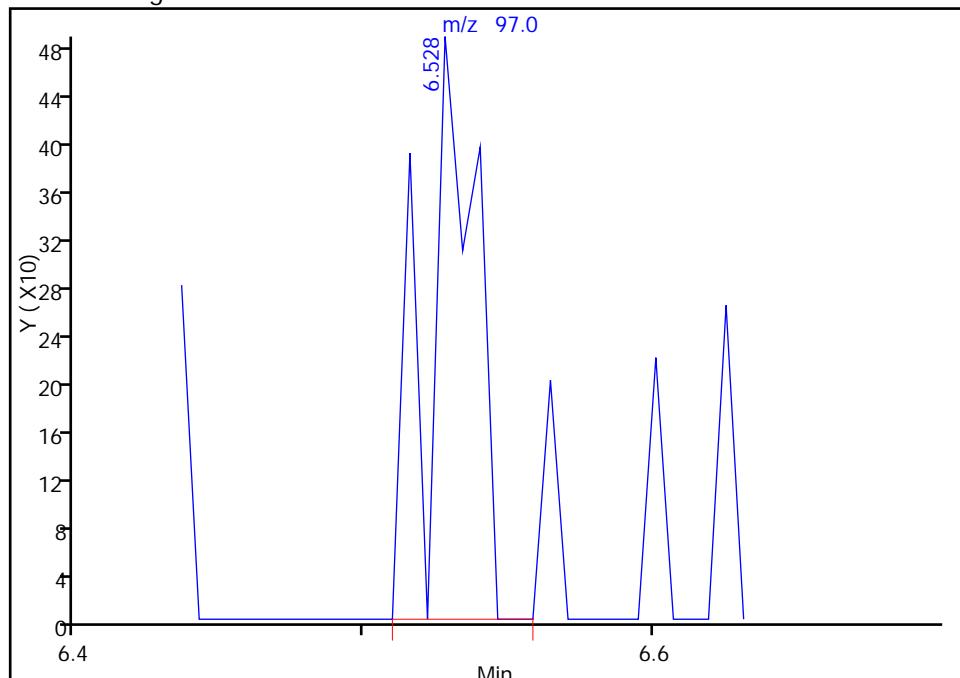
Not Detected

Expected RT: 6.53

## Processing Integration Results



## Manual Integration Results



Reviewer: fergusond, 16-Apr-2015 08:55:53

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Client Sample ID: HD-MW-162-0/1-0 DL Lab Sample ID: 180-42975-2 DL

Matrix: Water Lab File ID: 50416016.D

Analysis Method: 8260C Date Collected: 04/10/2015 09:50

Sample wt/vol: 5 (mL) Date Analyzed: 04/16/2015 16:29

Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 40

Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)

% Moisture: \_\_\_\_\_ Level: (low/med) Low

Analysis Batch No.: 138685 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
74-87-3	<i>Chloromethane</i>	40	U	40	11
75-01-4	<i>Vinyl chloride</i>	40	U	40	9.1
74-83-9	<i>Bromomethane</i>	40	U	40	13
75-00-3	<i>Chloroethane</i>	40	U	40	8.6
75-35-4	<i>1,1-Dichloroethene</i>	40	U	40	12
67-64-1	<i>Acetone</i>	200	U	200	100
75-15-0	<i>Carbon disulfide</i>	40	U	40	8.5
75-09-2	<i>Methylene Chloride</i>	22	J B	40	5.0
156-60-5	<i>trans-1,2-Dichloroethene</i>	40	U	40	6.8
1634-04-4	<i>Methyl tert-butyl ether</i>	40	U	40	7.3
75-34-3	<i>1,1-Dichloroethane</i>	40	U	40	4.7
156-59-2	<i>cis-1,2-Dichloroethene</i>	40	U	40	9.5
74-97-5	<i>Bromochloromethane</i>	40	U	40	7.2
78-93-3	<i>2-Butanone (MEK)</i>	200	U	200	22
67-66-3	<i>Chloroform</i>	40	U	40	6.8
71-55-6	<i>1,1,1-Trichloroethane</i>	40	U	40	11
56-23-5	<i>Carbon tetrachloride</i>	40	U	40	5.5
71-43-2	<i>Benzene</i>	40	U	40	4.2
107-06-2	<i>1,2-Dichloroethane</i>	40	U	40	8.5
79-01-6	<i>Trichloroethene</i>	180		40	5.7
78-87-5	<i>1,2-Dichloropropane</i>	40	U	40	3.8
75-27-4	<i>Bromodichloromethane</i>	40	U	40	5.2
10061-01-5	<i>cis-1,3-Dichloropropene</i>	40	U	40	7.5
108-10-1	<i>4-Methyl-2-pentanone (MIBK)</i>	200	U	200	21
108-88-3	<i>Toluene</i>	40	U	40	6.0
10061-02-6	<i>trans-1,3-Dichloropropene</i>	40	U	40	5.9
79-00-5	<i>1,1,2-Trichloroethane</i>	40	U	40	8.1
127-18-4	<i>Tetrachloroethene</i>	700		40	5.9
591-78-6	<i>2-Hexanone</i>	200	U	200	6.4
124-48-1	<i>Dibromochloromethane</i>	40	U	40	5.5
106-93-4	<i>1,2-Dibromoethane (EDB)</i>	40	U	40	7.2
108-90-7	<i>Chlorobenzene</i>	40	U	40	5.4
630-20-6	<i>1,1,1,2-Tetrachloroethane</i>	40	U	40	11
100-41-4	<i>Ethylbenzene</i>	40	U	40	9.1
1330-20-7	<i>Xylenes, Total</i>	120	U	120	20
100-42-5	<i>Styrene</i>	40	U	40	3.9

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: HD-MW-162-0/1-0 DL Lab Sample ID: 180-42975-2 DL  
Matrix: Water Lab File ID: 50416016.D  
Analysis Method: 8260C Date Collected: 04/10/2015 09:50  
Sample wt/vol: 5 (mL) Date Analyzed: 04/16/2015 16:29  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 40  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138685 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-25-2	Bromoform	40	U	40	7.7
79-34-5	1,1,2,2-Tetrachloroethane	40	U	40	8.0
107-13-1	Acrylonitrile	800	U	800	22
123-91-1	1,4-Dioxane	8000	U	8000	1400

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	107		64-135
2037-26-5	Toluene-d8 (Surr)	102		71-118
460-00-4	4-Bromofluorobenzene (Surr)	98		70-118
1868-53-7	Dibromofluoromethane (Surr)	110		70-128

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\50416016.D  
 Lims ID: 180-42975-B-2 Lab Sample ID: 180-42975-2  
 Client ID: HD-MW-162-01-0  
 Sample Type: Client  
 Inject. Date: 16-Apr-2015 16:29:30 ALS Bottle#: 16 Worklist Smp#: 16  
 Purge Vol: 5.000 mL Dil. Factor: 40.0000  
 Sample Info: 180-42975-B-2, 40x  
 Misc. Info.: 180-0006494-016  
 Operator ID: 001562 Instrument ID: CHHP5  
 Method: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 16:58:40 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICAL File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 16-Apr-2015 16:58:40

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.298	4.317	-0.019	0	175233	1000.0	
* 2 Fluorobenzene (IS)	96	7.273	7.268	0.005	98	442739	50.0	
* 3 Chlorobenzene-d5	119	10.364	10.358	0.006	89	107268	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.682	12.682	0.000	97	156120	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.525	6.525	0.000	94	110666	55.0	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.908	6.897	0.011	0	142020	53.5	
\$ 7 Toluene-d8 (Surr)	98	8.928	8.916	0.012	94	434202	50.8	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.526	11.532	-0.006	87	150567	48.9	
12 Chloromethane	50		1.786				ND	
13 Vinyl chloride	62		1.914				ND	
15 Bromomethane	94		2.267				ND	
16 Chloroethane	64		2.413				ND	
22 1,1-Dichloroethene	96		3.386				ND	
24 Acetone	43		3.496				ND	
26 Carbon disulfide	76		3.691				ND	
31 Methylene Chloride	84	4.128	4.147	-0.019	64	8147	2.76	M
33 Acrylonitrile	53		4.548				ND	
34 trans-1,2-Dichloroethene	96		4.561				ND	
35 Methyl tert-butyl ether	73		4.591				ND	
37 1,1-Dichloroethane	63		5.175				ND	
45 cis-1,2-Dichloroethene	96		5.935				ND	
46 2-Butanone (MEK)	43		5.990				ND	
49 Chlorobromomethane	128		6.227				ND	
52 Chloroform	83		6.337				ND	
53 1,1,1-Trichloroethane	97		6.525				ND	
56 Carbon tetrachloride	117		6.720				ND	
58 Benzene	78		6.951				ND	
59 1,2-Dichloroethane	62		6.982				ND	
64 Trichloroethene	130	7.669	7.669	0.000	95	60035	22.8	
67 1,2-Dichloropropane	63		7.900				ND	
70 1,4-Dioxane	88		8.052				ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ng	Flags
71 Dichlorobromomethane	83		8.192				ND	
74 cis-1,3-Dichloropropene	75		8.655				ND	
75 4-Methyl-2-pentanone (MIBK)	43		8.819				ND	
76 Toluene	91		8.989				ND	
77 trans-1,3-Dichloropropene	75		9.220				ND	
79 1,1,2-Trichloroethane	97		9.397				ND	
80 Tetrachloroethene	164	9.536	9.531	0.005	96	188822	87.8	
82 2-Hexanone	43		9.652				ND	
84 Chlorodibromomethane	129		9.786				ND	
85 Ethylene Dibromide	107		9.902				ND	
87 Chlorobenzene	112		10.388				ND	
89 1,1,1,2-Tetrachloroethane	131		10.474				ND	
90 Ethylbenzene	106		10.498				ND	
91 m-Xylene & p-Xylene	106		10.614				ND	
92 o-Xylene	106		11.009				ND	
93 Styrene	104		11.021				ND	
94 Bromoform	173		11.204				ND	
99 1,1,2,2-Tetrachloroethane	83		11.672				ND	
S 133 Xylenes, Total	106		1.000				ND	

**QC Flag Legend**

Review Flags

M - Manually Integrated

**Reagents:**

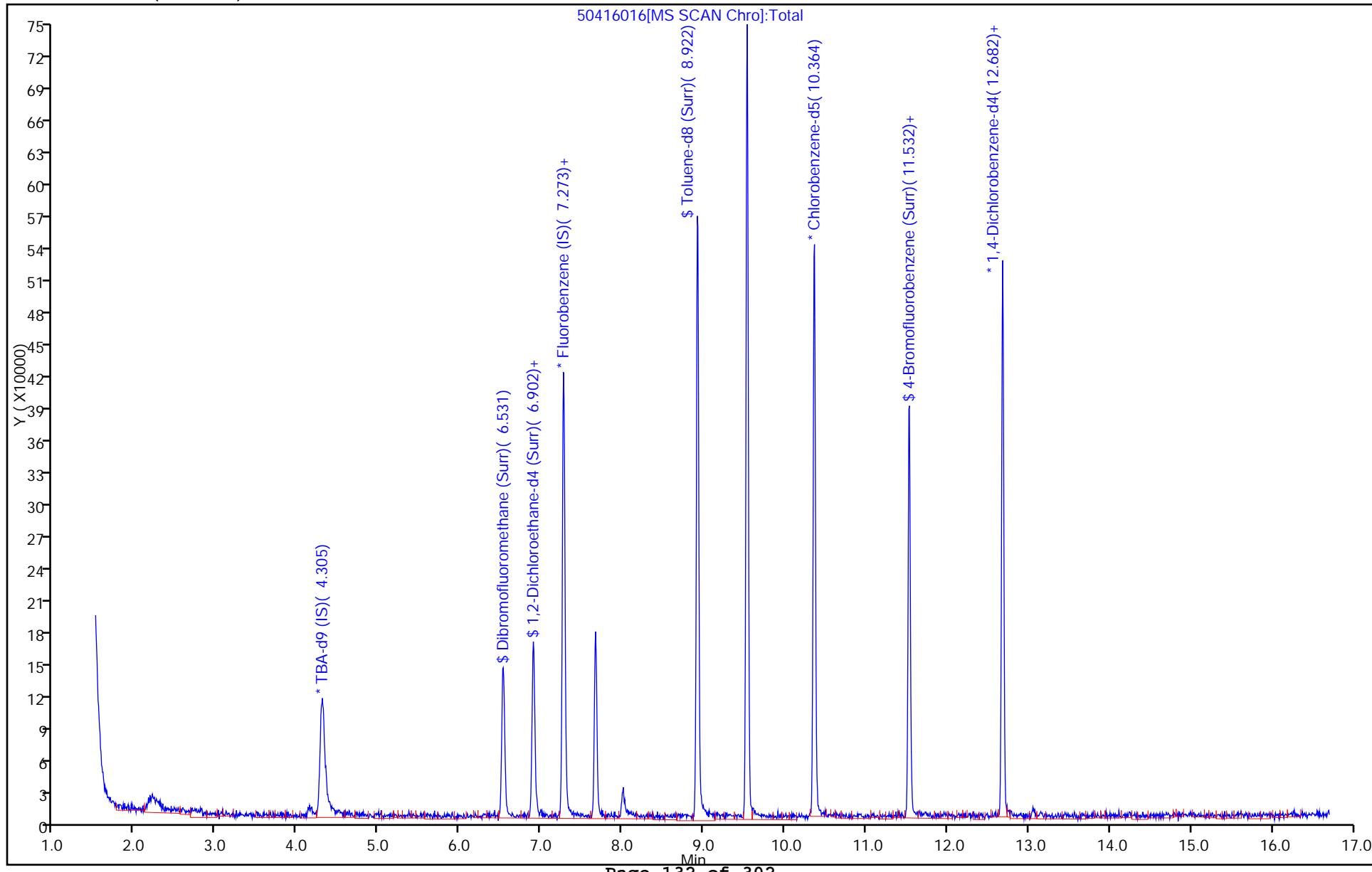
VOA8260INT_00031	Amount Added: 2.00	Units: uL	Run Reagent
VOA8260SURR_00033	Amount Added: 2.00	Units: uL	Run Reagent

Report Date: 16-Apr-2015 16:58:40

Chrom Revision: 2.2 13-Mar-2015 11:20:44

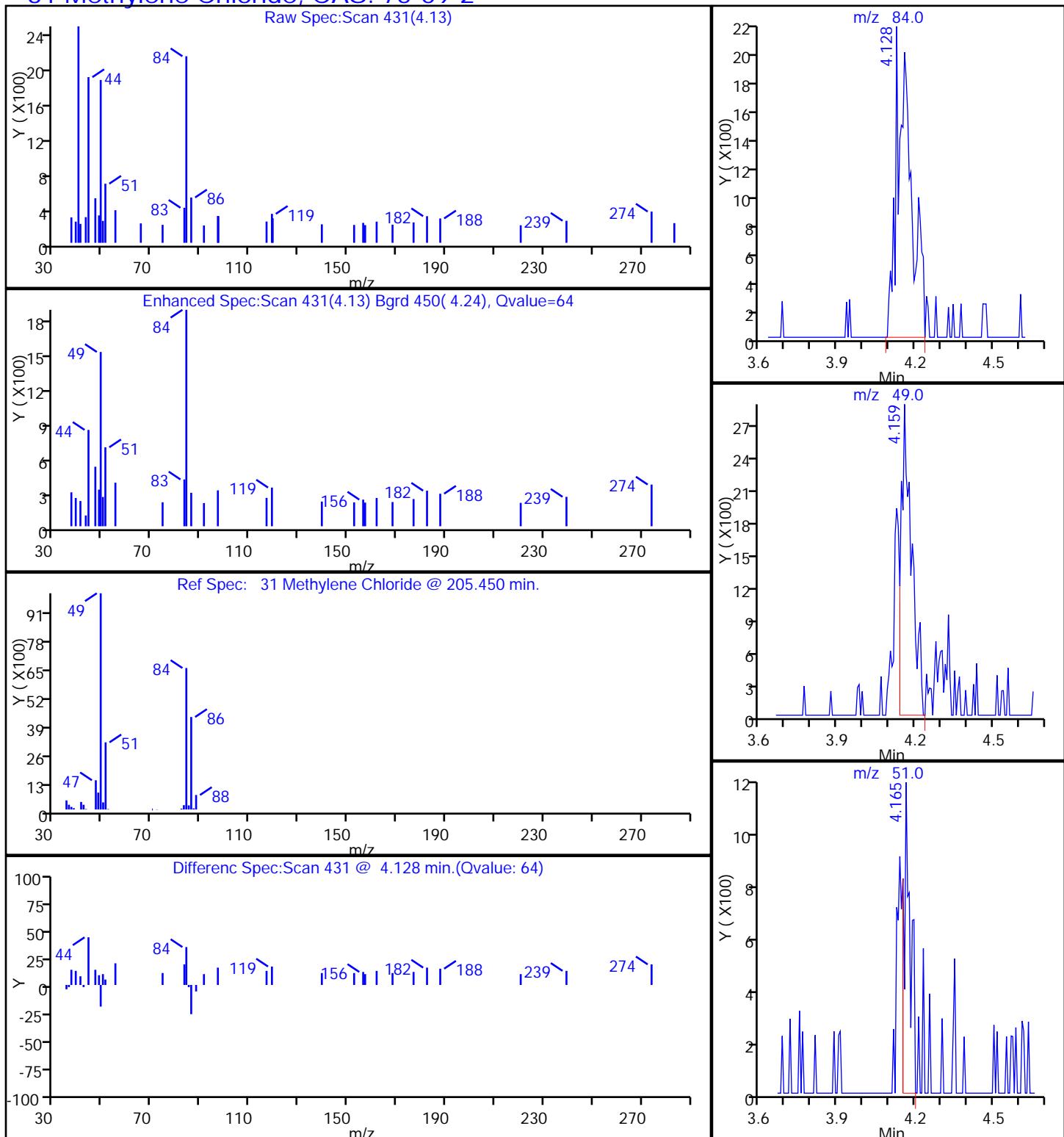
TestAmerica Pittsburgh

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150416-6494.b\\50416016.D  
Injection Date: 16-Apr-2015 16:29:30 Instrument ID: CHHP5 Operator ID: 001562  
Lims ID: 180-42975-B-2 Lab Sample ID: 180-42975-2 Worklist Smp#: 16  
Client ID: HD-MW-162-0/1-0  
Purge Vol: 5.000 mL Dil. Factor: 40.0000 ALS Bottle#: 16  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)

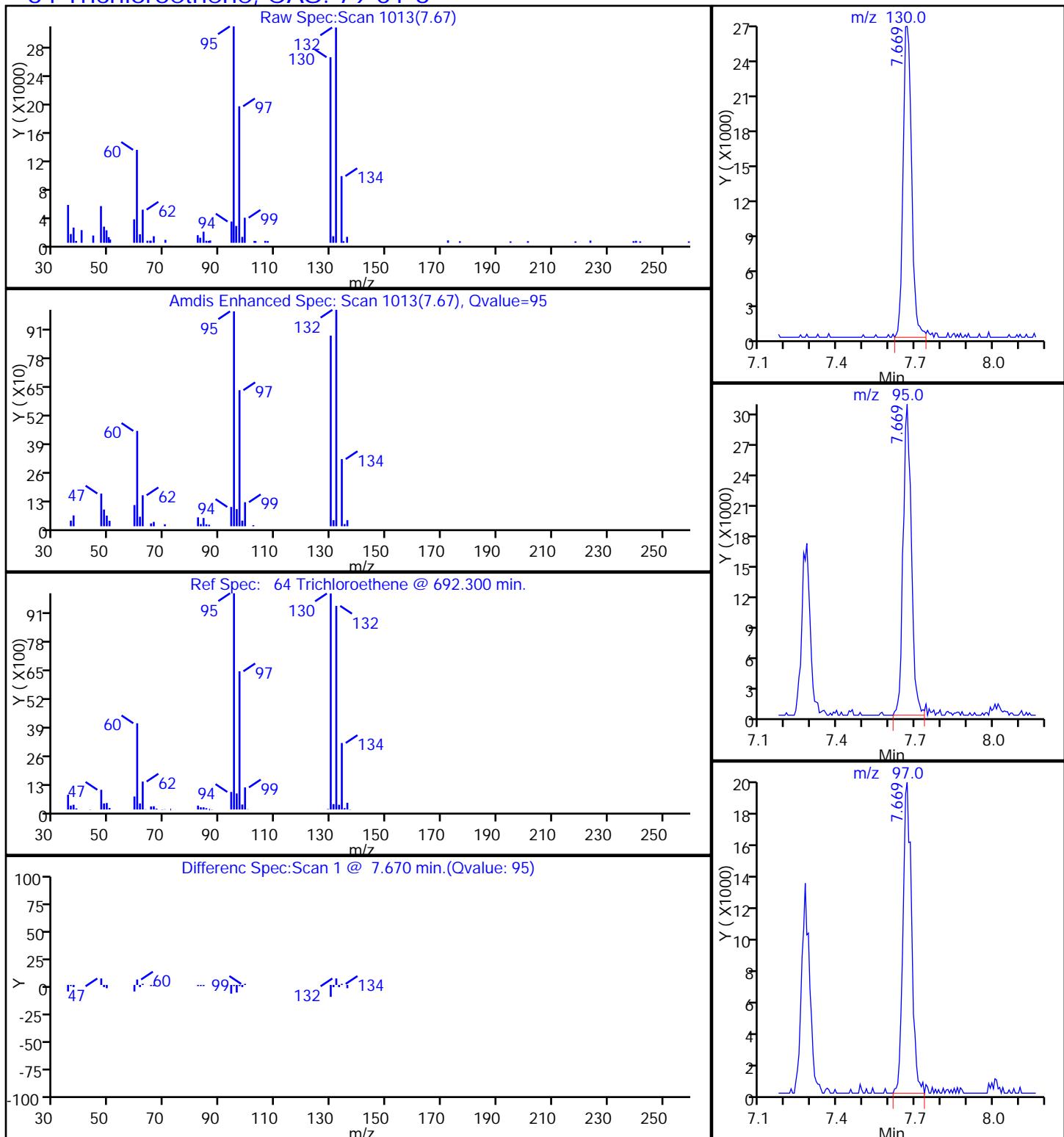


TestAmerica Pittsburgh  
 Data File: \PITCHROM\ChromData\CHHP5\20150416-6494.b\50416016.D  
 Injection Date: 16-Apr-2015 16:29:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-B-2 Lab Sample ID: 180-42975-2  
 Client ID: HD-MW-162-0/1-0  
 Operator ID: 001562 ALS Bottle#: 16 Worklist Smp#: 16  
 Purge Vol: 5.000 mL Dil. Factor: 40.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 ( 0.18 mm) Detector: MS SCAN

### 31 Methylene Chloride, CAS: 75-09-2

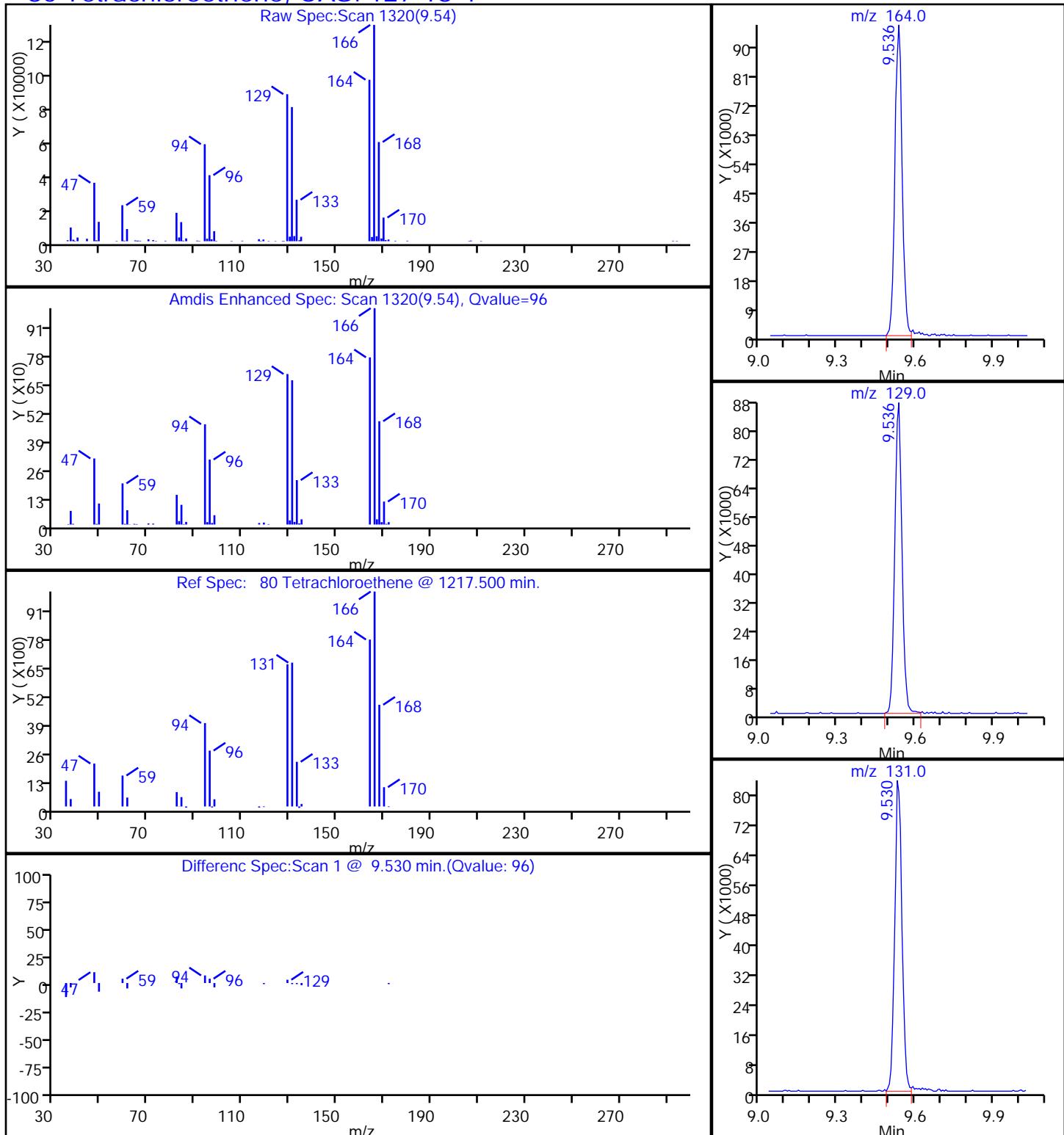


TestAmerica Pittsburgh  
 Data File: \PITCHROM\ChromData\CHHP5\20150416-6494.b\50416016.D  
 Injection Date: 16-Apr-2015 16:29:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-B-2 Lab Sample ID: 180-42975-2  
 Client ID: HD-MW-162-0/1-0  
 Operator ID: 001562 ALS Bottle#: 16 Worklist Smp#: 16  
 Purge Vol: 5.000 mL Dil. Factor: 40.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 ( 0.18 mm) Detector: MS SCAN

**64 Trichloroethene, CAS: 79-01-6**

TestAmerica Pittsburgh  
 Data File: \PITCHROM\ChromData\CHHP5\20150416-6494.b\50416016.D  
 Injection Date: 16-Apr-2015 16:29:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-B-2 Lab Sample ID: 180-42975-2  
 Client ID: HD-MW-162-0/1-0  
 Operator ID: 001562 ALS Bottle#: 16 Worklist Smp#: 16  
 Purge Vol: 5.000 mL Dil. Factor: 40.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 ( 0.18 mm) Detector: MS SCAN

### 80 Tetrachloroethene, CAS: 127-18-4



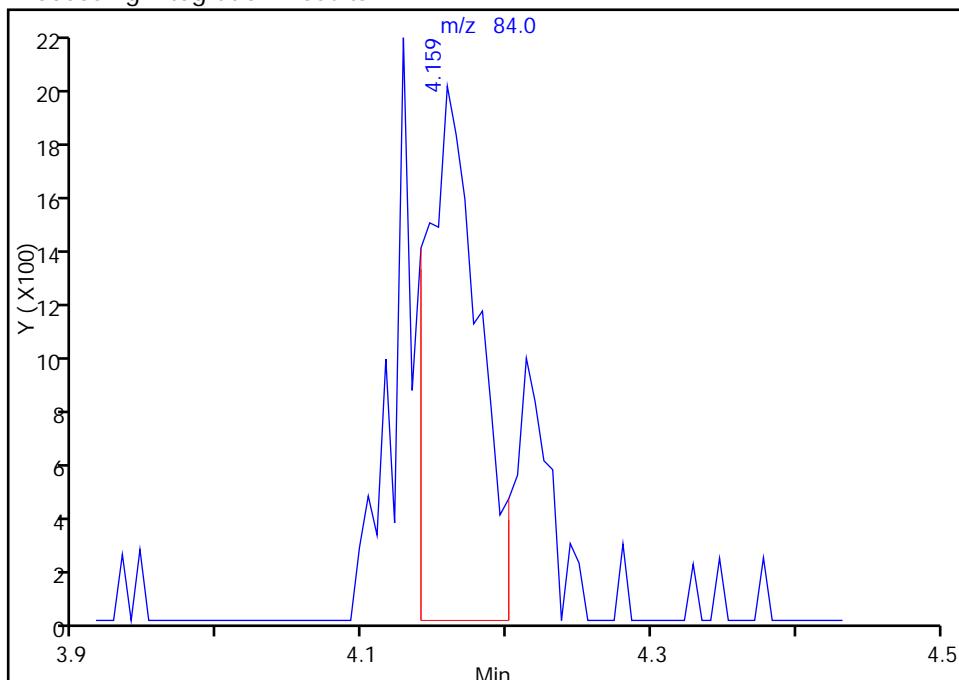
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\50416016.D  
 Injection Date: 16-Apr-2015 16:29:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-B-2 Lab Sample ID: 180-42975-2  
 Client ID: HD-MW-162-0/1-0  
 Operator ID: 001562 ALS Bottle#: 16 Worklist Smp#: 16  
 Purge Vol: 5.000 mL Dil. Factor: 40.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 31 Methylene Chloride, CAS: 75-09-2

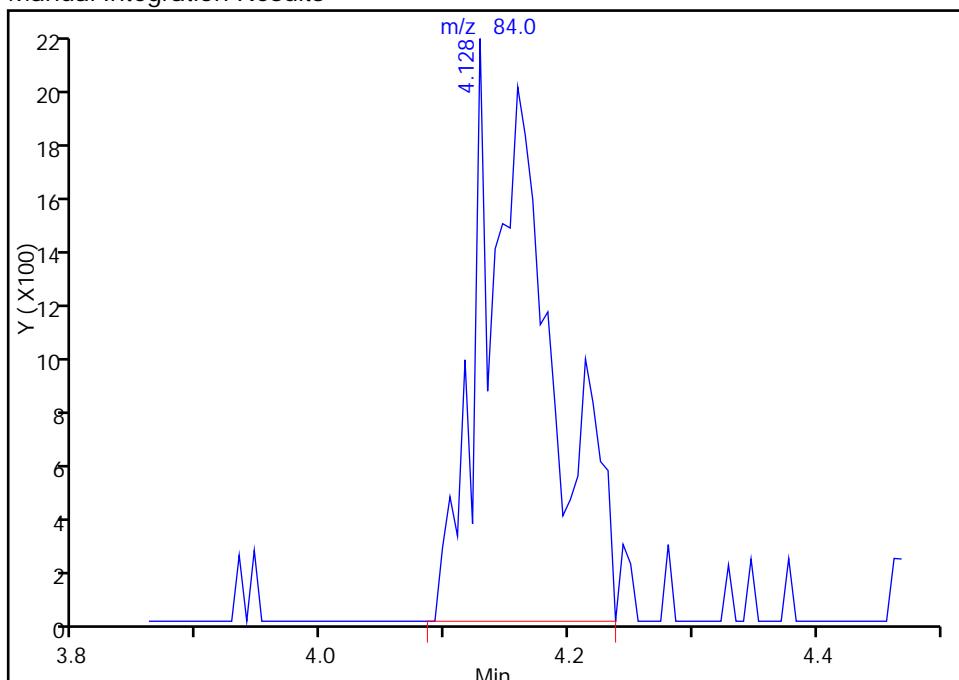
RT: 4.16  
 Area: 4922  
 Amount: 1.666984  
 Amount Units: ng

## Processing Integration Results



RT: 4.13  
 Area: 8147  
 Amount: 2.759227  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 16-Apr-2015 16:58:40

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: HD-MW-169-0/1-0 Lab Sample ID: 180-42975-3  
Matrix: Water Lab File ID: 50416017.D  
Analysis Method: 8260C Date Collected: 04/10/2015 12:32  
Sample wt/vol: 5 (mL) Date Analyzed: 04/16/2015 16:53  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138685 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
74-87-3	Chloromethane	1.0	U	1.0	0.28
75-01-4	Vinyl chloride	1.0	U	1.0	0.23
74-83-9	Bromomethane	1.0	U	1.0	0.31
75-00-3	Chloroethane	1.0	U	1.0	0.21
75-35-4	1,1-Dichloroethene	1.0	U	1.0	0.30
67-64-1	Acetone	5.0	U	5.0	2.5
75-15-0	Carbon disulfide	1.0	U	1.0	0.21
75-09-2	Methylene Chloride	0.14	J B	1.0	0.13
156-60-5	trans-1,2-Dichloroethene	1.0	U	1.0	0.17
1634-04-4	Methyl tert-butyl ether	1.0	U	1.0	0.18
75-34-3	1,1-Dichloroethane	1.0	U	1.0	0.12
156-59-2	cis-1,2-Dichloroethene	1.0	U	1.0	0.24
74-97-5	Bromochloromethane	1.0	U	1.0	0.18
78-93-3	2-Butanone (MEK)	5.0	U	5.0	0.55
67-66-3	Chloroform	1.0	U	1.0	0.17
71-55-6	1,1,1-Trichloroethane	1.0	U	1.0	0.29
56-23-5	Carbon tetrachloride	1.0	U	1.0	0.14
71-43-2	Benzene	1.0	U	1.0	0.11
107-06-2	1,2-Dichloroethane	1.0	U	1.0	0.21
79-01-6	Trichloroethene	1.0	U	1.0	0.14
78-87-5	1,2-Dichloropropane	1.0	U	1.0	0.095
75-27-4	Bromodichloromethane	1.0	U	1.0	0.13
10061-01-5	cis-1,3-Dichloropropene	1.0	U	1.0	0.19
108-10-1	4-Methyl-2-pentanone (MIBK)	5.0	U	5.0	0.53
108-88-3	Toluene	1.0	U	1.0	0.15
10061-02-6	trans-1,3-Dichloropropene	1.0	U	1.0	0.15
79-00-5	1,1,2-Trichloroethane	1.0	U	1.0	0.20
127-18-4	Tetrachloroethene	1.0	U	1.0	0.15
591-78-6	2-Hexanone	5.0	U	5.0	0.16
124-48-1	Dibromochloromethane	1.0	U	1.0	0.14
106-93-4	1,2-Dibromoethane (EDB)	1.0	U	1.0	0.18
108-90-7	Chlorobenzene	1.0	U	1.0	0.14
630-20-6	1,1,1,2-Tetrachloroethane	1.0	U	1.0	0.28
100-41-4	Ethylbenzene	1.0	U	1.0	0.23
1330-20-7	Xylenes, Total	3.0	U	3.0	0.49
100-42-5	Styrene	1.0	U	1.0	0.097

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: HD-MW-169-01/0 Lab Sample ID: 180-42975-3  
Matrix: Water Lab File ID: 50416017.D  
Analysis Method: 8260C Date Collected: 04/10/2015 12:32  
Sample wt/vol: 5 (mL) Date Analyzed: 04/16/2015 16:53  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138685 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-25-2	Bromoform	1.0	U	1.0	0.19
79-34-5	1,1,2,2-Tetrachloroethane	1.0	U	1.0	0.20
107-13-1	Acrylonitrile	20	U	20	0.55
123-91-1	1,4-Dioxane	200	U	200	34

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	115		64-135
2037-26-5	Toluene-d8 (Surr)	98		71-118
460-00-4	4-Bromofluorobenzene (Surr)	92		70-118
1868-53-7	Dibromofluoromethane (Surr)	109		70-128

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\50416017.D  
 Lims ID: 180-42975-C-3 Lab Sample ID: 180-42975-3  
 Client ID: HD-MW-169-01-0  
 Sample Type: Client  
 Inject. Date: 16-Apr-2015 16:53:30 ALS Bottle#: 17 Worklist Smp#: 17  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: 180-42975-C-3  
 Misc. Info.: 180-0006494-017  
 Operator ID: 001562 Instrument ID: CHHP5  
 Method: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 17-Apr-2015 07:48:32 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICAL File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK047

First Level Reviewer: fergusond Date: 17-Apr-2015 07:48:32

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.296	4.317	-0.021	0	178588	1000.0	
* 2 Fluorobenzene (IS)	96	7.271	7.268	0.003	98	432619	50.0	
* 3 Chlorobenzene-d5	119	10.362	10.358	0.004	89	105105	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.679	12.682	-0.003	98	144245	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.529	6.525	0.004	93	107053	54.4	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.900	6.897	0.003	0	148921	57.4	
\$ 7 Toluene-d8 (Surr)	98	8.920	8.916	0.004	94	409374	48.9	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.524	11.532	-0.008	88	139456	46.2	
12 Chloromethane	50		1.786				ND	
13 Vinyl chloride	62		1.914				ND	
15 Bromomethane	94		2.267				ND	
16 Chloroethane	64		2.413				ND	
22 1,1-Dichloroethene	96		3.386				ND	
24 Acetone	43	3.506	3.496	0.010	45	2373	2.68	
26 Carbon disulfide	76		3.691				ND	
31 Methylene Chloride	84	4.144	4.147	-0.003	0	2032	0.7043	
33 Acrylonitrile	53		4.548				ND	
34 trans-1,2-Dichloroethene	96		4.561				ND	
35 Methyl tert-butyl ether	73		4.591				ND	
37 1,1-Dichloroethane	63		5.175				ND	
45 cis-1,2-Dichloroethene	96		5.935				ND	
46 2-Butanone (MEK)	43		5.990				ND	
49 Chlorobromomethane	128		6.227				ND	
52 Chloroform	83		6.337				ND	
53 1,1,1-Trichloroethane	97		6.525				ND	
56 Carbon tetrachloride	117		6.720				ND	
58 Benzene	78		6.951				ND	
59 1,2-Dichloroethane	62		6.982				ND	
64 Trichloroethene	130		7.669				ND	
67 1,2-Dichloropropane	63		7.900				ND	
70 1,4-Dioxane	88		8.052				ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ng	Flags
71 Dichlorobromomethane	83		8.192				ND	
74 cis-1,3-Dichloropropene	75		8.655				ND	
75 4-Methyl-2-pentanone (MIBK)	43		8.819				ND	
76 Toluene	91		8.989				ND	
77 trans-1,3-Dichloropropene	75		9.220				ND	
79 1,1,2-Trichloroethane	97		9.397				ND	
80 Tetrachloroethene	164		9.531				ND	
82 2-Hexanone	43		9.652				ND	
84 Chlorodibromomethane	129		9.786				ND	
85 Ethylene Dibromide	107		9.902				ND	
87 Chlorobenzene	112		10.388				ND	
89 1,1,1,2-Tetrachloroethane	131		10.474				ND	
90 Ethylbenzene	106		10.498				ND	
91 m-Xylene & p-Xylene	106		10.614				ND	
92 o-Xylene	106		11.009				ND	
93 Styrene	104		11.021				ND	
94 Bromoform	173		11.204				ND	
99 1,1,2,2-Tetrachloroethane	83		11.672				ND	
S 133 Xylenes, Total	106		1.000				ND	

**Reagents:**

VOA8260INT\_00031  
 VOA8260SURR\_00033

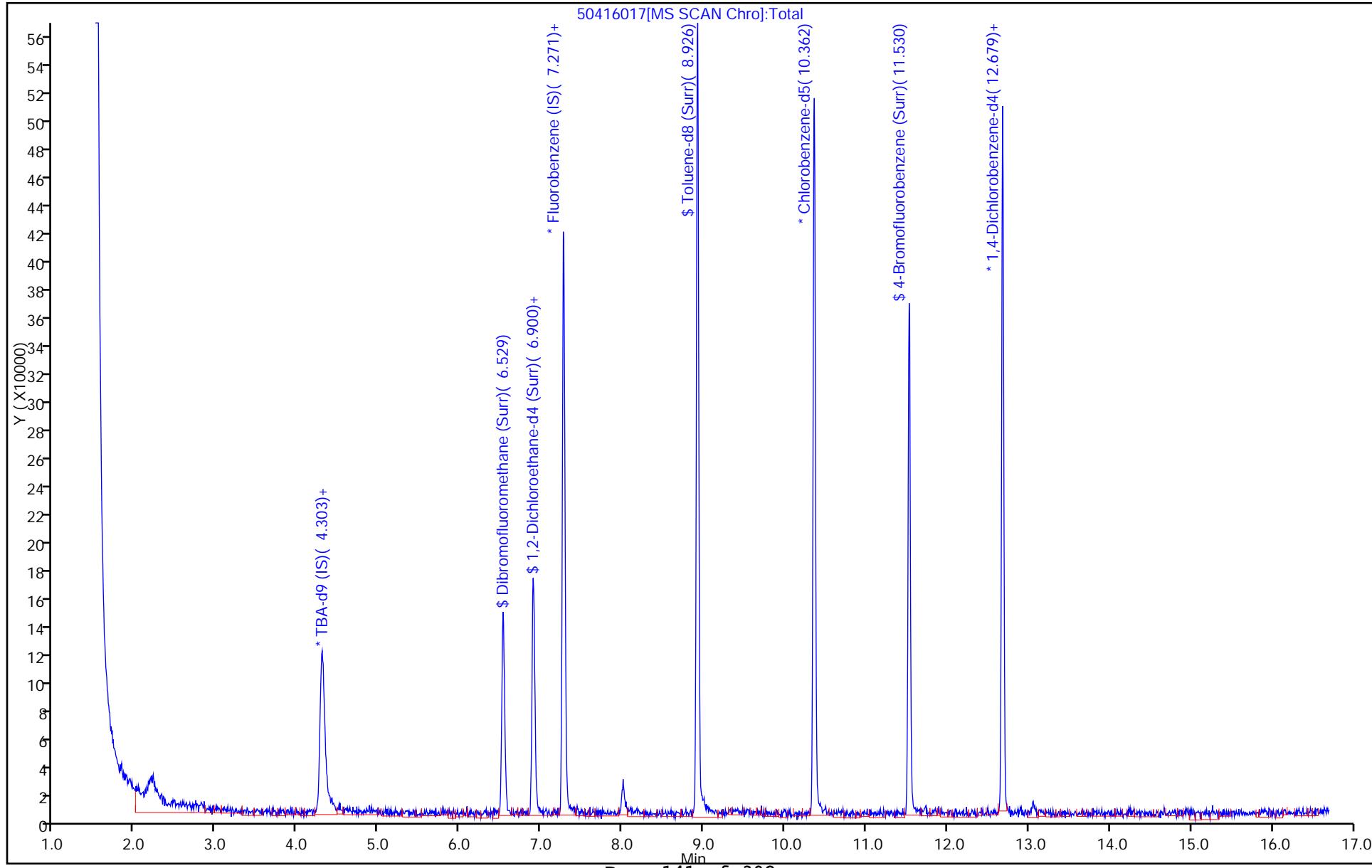
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 Amount Added: 2.00 Units: uL Run Reagent

Report Date: 17-Apr-2015 07:48:32

Chrom Revision: 2.2 13-Mar-2015 11:20:44

## TestAmerica Pittsburgh

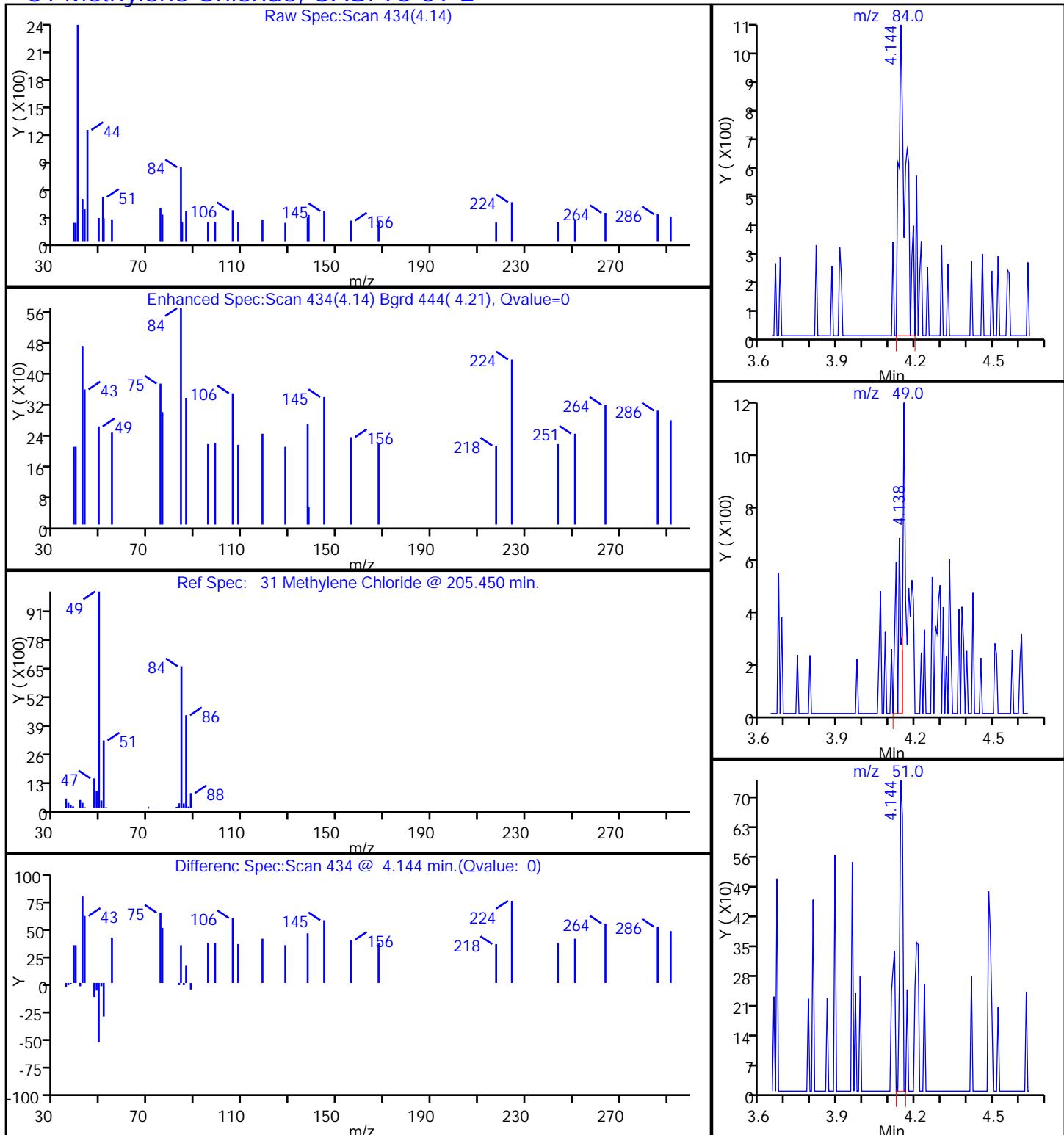
Data File: \PITCHROM\ChromData\CHHP5\20150416-6494.b\50416017.D  
Injection Date: 16-Apr-2015 16:53:30 Instrument ID: CHHP5 Operator ID: 001562  
Lims ID: 180-42975-C-3 Lab Sample ID: 180-42975-3 Worklist Smp#: 17  
Client ID: HD-MW-169-0/1-0  
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 17  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)



## TestAmerica Pittsburgh

Data File: \PITCHROM\ChromData\CHHP5\20150416-6494.b\50416017.D  
 Injection Date: 16-Apr-2015 16:53:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-C-3 Lab Sample ID: 180-42975-3  
 Client ID: HD-MW-169-0/1-0  
 Operator ID: 001562 ALS Bottle#: 17 Worklist Smp#: 17  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 ( 0.18 mm) Detector: MS SCAN

## 31 Methylene Chloride, CAS: 75-09-2



FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Client Sample ID: HD-QC2-0/1-2 Lab Sample ID: 180-42975-4

Matrix: Water Lab File ID: 50415007.D

Analysis Method: 8260C Date Collected: 04/10/2015 12:00

Sample wt/vol: 5 (mL) Date Analyzed: 04/15/2015 15:33

Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1

Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)

% Moisture: \_\_\_\_\_ Level: (low/med) Low

Analysis Batch No.: 138583 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
74-87-3	Chloromethane	1.0	U	1.0	0.28
75-01-4	Vinyl chloride	1.0	U	1.0	0.23
74-83-9	Bromomethane	1.0	U	1.0	0.31
75-00-3	Chloroethane	1.0	U	1.0	0.21
75-35-4	1,1-Dichloroethene	1.0	U	1.0	0.30
67-64-1	Acetone	5.0	U	5.0	2.5
75-15-0	Carbon disulfide	1.0	U *	1.0	0.21
75-09-2	Methylene Chloride	0.22	J	1.0	0.13
156-60-5	trans-1,2-Dichloroethene	1.0	U	1.0	0.17
1634-04-4	Methyl tert-butyl ether	1.0	U	1.0	0.18
75-34-3	1,1-Dichloroethane	1.0	U	1.0	0.12
156-59-2	cis-1,2-Dichloroethene	1.0	U	1.0	0.24
74-97-5	Bromochloromethane	1.0	U	1.0	0.18
78-93-3	2-Butanone (MEK)	5.0	U	5.0	0.55
67-66-3	Chloroform	1.0	U	1.0	0.17
71-55-6	1,1,1-Trichloroethane	1.0	U	1.0	0.29
56-23-5	Carbon tetrachloride	1.0	U	1.0	0.14
71-43-2	Benzene	1.0	U	1.0	0.11
107-06-2	1,2-Dichloroethane	1.0	U	1.0	0.21
79-01-6	Trichloroethene	1.0	U	1.0	0.14
78-87-5	1,2-Dichloropropane	1.0	U	1.0	0.095
75-27-4	Bromodichloromethane	1.0	U	1.0	0.13
10061-01-5	cis-1,3-Dichloropropene	1.0	U	1.0	0.19
108-10-1	4-Methyl-2-pentanone (MIBK)	5.0	U	5.0	0.53
108-88-3	Toluene	1.0	U	1.0	0.15
10061-02-6	trans-1,3-Dichloropropene	1.0	U	1.0	0.15
79-00-5	1,1,2-Trichloroethane	1.0	U	1.0	0.20
127-18-4	Tetrachloroethene	1.0	U	1.0	0.15
591-78-6	2-Hexanone	5.0	U	5.0	0.16
124-48-1	Dibromochloromethane	1.0	U	1.0	0.14
106-93-4	1,2-Dibromoethane (EDB)	1.0	U	1.0	0.18
108-90-7	Chlorobenzene	1.0	U	1.0	0.14
630-20-6	1,1,1,2-Tetrachloroethane	1.0	U	1.0	0.28
100-41-4	Ethylbenzene	1.0	U	1.0	0.23
1330-20-7	Xylenes, Total	3.0	U	3.0	0.49
100-42-5	Styrene	1.0	U	1.0	0.097

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: HD-QC2-0/1-2 Lab Sample ID: 180-42975-4  
Matrix: Water Lab File ID: 50415007.D  
Analysis Method: 8260C Date Collected: 04/10/2015 12:00  
Sample wt/vol: 5 (mL) Date Analyzed: 04/15/2015 15:33  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138583 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-25-2	Bromoform	1.0	U	1.0	0.19
79-34-5	1,1,2,2-Tetrachloroethane	1.0	U	1.0	0.20
107-13-1	Acrylonitrile	20	U	20	0.55
123-91-1	1,4-Dioxane	200	U	200	34

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	114		64-135
2037-26-5	Toluene-d8 (Surr)	103		71-118
460-00-4	4-Bromofluorobenzene (Surr)	97		70-118
1868-53-7	Dibromofluoromethane (Surr)	112		70-128

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415007.D  
 Lims ID: 180-42975-A-4 Lab Sample ID: 180-42975-4  
 Client ID: HD-QC2-01/2  
 Sample Type: Client  
 Inject. Date: 15-Apr-2015 15:33:30 ALS Bottle#: 7 Worklist Smp#: 7  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: 180-42975-A-4  
 Misc. Info.: 180-0006480-007  
 Operator ID: 001562 Instrument ID: CHHP5  
 Method: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 07:33:58 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICAL File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 16-Apr-2015 07:33:58

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.296	4.302	-0.006	0	167449	1000.0	
* 2 Fluorobenzene (IS)	96	7.271	7.271	0.000	99	422314	50.0	
* 3 Chlorobenzene-d5	119	10.361	10.361	0.000	89	101006	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.679	12.679	0.000	96	149655	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.529	6.528	0.001	93	107812	56.1	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.906	6.899	0.007	0	144790	57.2	
\$ 7 Toluene-d8 (Surr)	98	8.926	8.919	0.007	94	414777	51.5	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.529	11.529	0.000	88	140487	48.4	
12 Chloromethane	50		1.789				ND	
13 Vinyl chloride	62		1.917				ND	
15 Bromomethane	94		2.264				ND	
16 Chloroethane	64		2.416				ND	
22 1,1-Dichloroethene	96		3.395				ND	
24 Acetone	43		3.499				ND	
26 Carbon disulfide	76		3.675				ND	
31 Methylene Chloride	84	4.150	4.143	0.007	50	3097	1.10	
33 Acrylonitrile	53		4.551				ND	
34 trans-1,2-Dichloroethene	96		4.563				ND	
35 Methyl tert-butyl ether	73		4.600				ND	
37 1,1-Dichloroethane	63		5.165				ND	
45 cis-1,2-Dichloroethene	96		5.938				ND	
46 2-Butanone (MEK)	43		5.987				ND	
49 Chlorobromomethane	128		6.224				ND	
52 Chloroform	83		6.339				ND	
53 1,1,1-Trichloroethane	97		6.528				ND	
56 Carbon tetrachloride	117		6.723				ND	
58 Benzene	78		6.954				ND	
59 1,2-Dichloroethane	62		6.984				ND	
64 Trichloroethene	130		7.666				ND	
67 1,2-Dichloropropane	63		7.897				ND	
70 1,4-Dioxane	88		8.049				ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ng	Flags
71 Dichlorobromomethane	83		8.195				ND	
74 cis-1,3-Dichloropropene	75		8.651				ND	
75 4-Methyl-2-pentanone (MIBK)	43		8.822				ND	
76 Toluene	91		8.986				ND	
77 trans-1,3-Dichloropropene	75		9.217				ND	
79 1,1,2-Trichloroethane	97		9.399				ND	
80 Tetrachloroethene	164		9.533				ND	
82 2-Hexanone	43		9.655				ND	
84 Chlorodibromomethane	129		9.789				ND	
85 Ethylene Dibromide	107		9.898				ND	
87 Chlorobenzene	112		10.391				ND	
89 1,1,1,2-Tetrachloroethane	131		10.470				ND	
90 Ethylbenzene	106		10.501				ND	
91 m-Xylene & p-Xylene	106		10.616				ND	
92 o-Xylene	106		11.012				ND	
93 Styrene	104		11.024				ND	
94 Bromoform	173		11.212				ND	
99 1,1,2,2-Tetrachloroethane	83		11.675				ND	
S 133 Xylenes, Total	106		1.000				ND	

**Reagents:**

VOA8260INT\_00031  
 VOA8260SURR\_00033

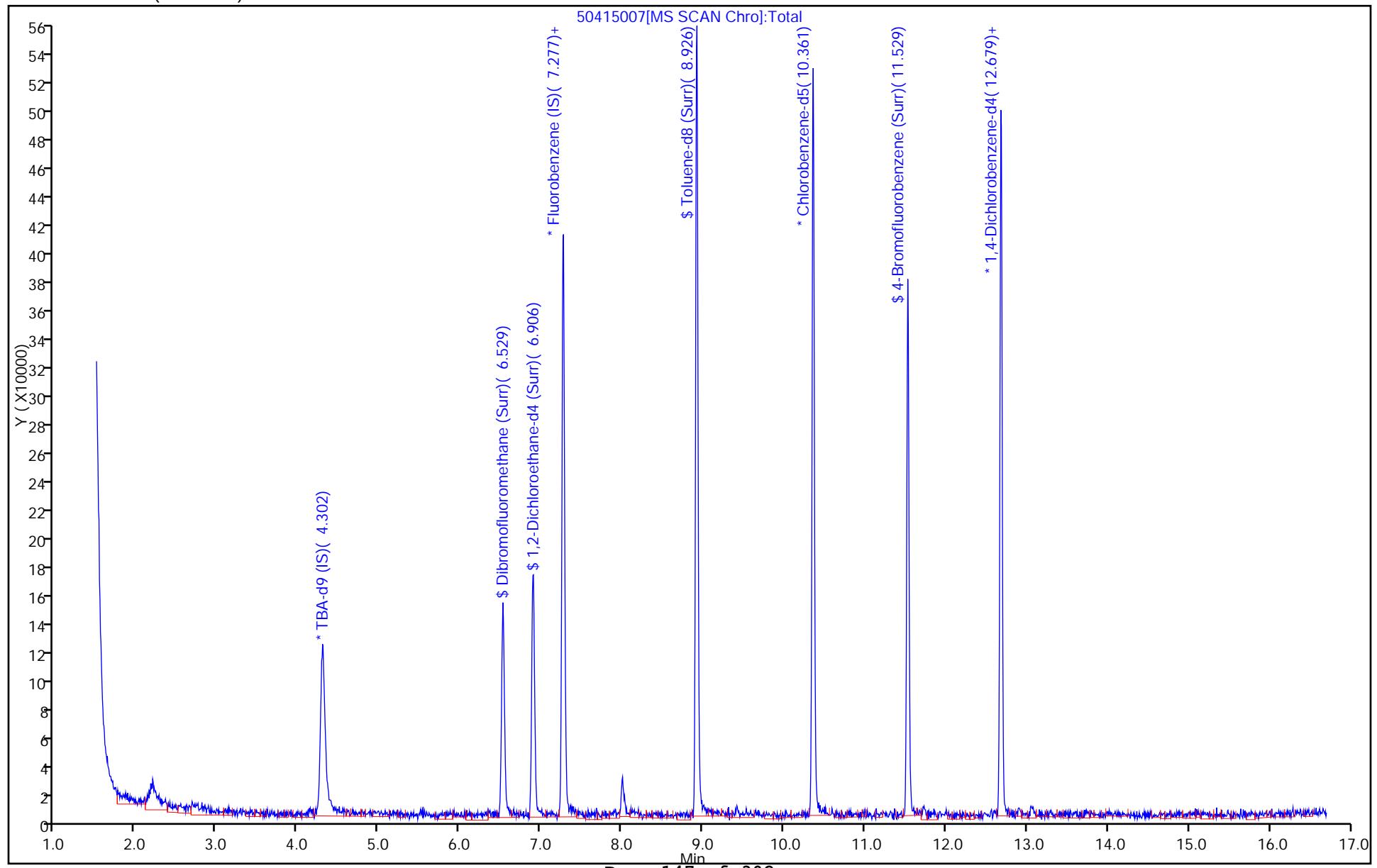
Amount Added: 2.00 Units: uL Run Reagent  
 Amount Added: 2.00 Units: uL Run Reagent

Report Date: 16-Apr-2015 07:33:58

Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Pittsburgh

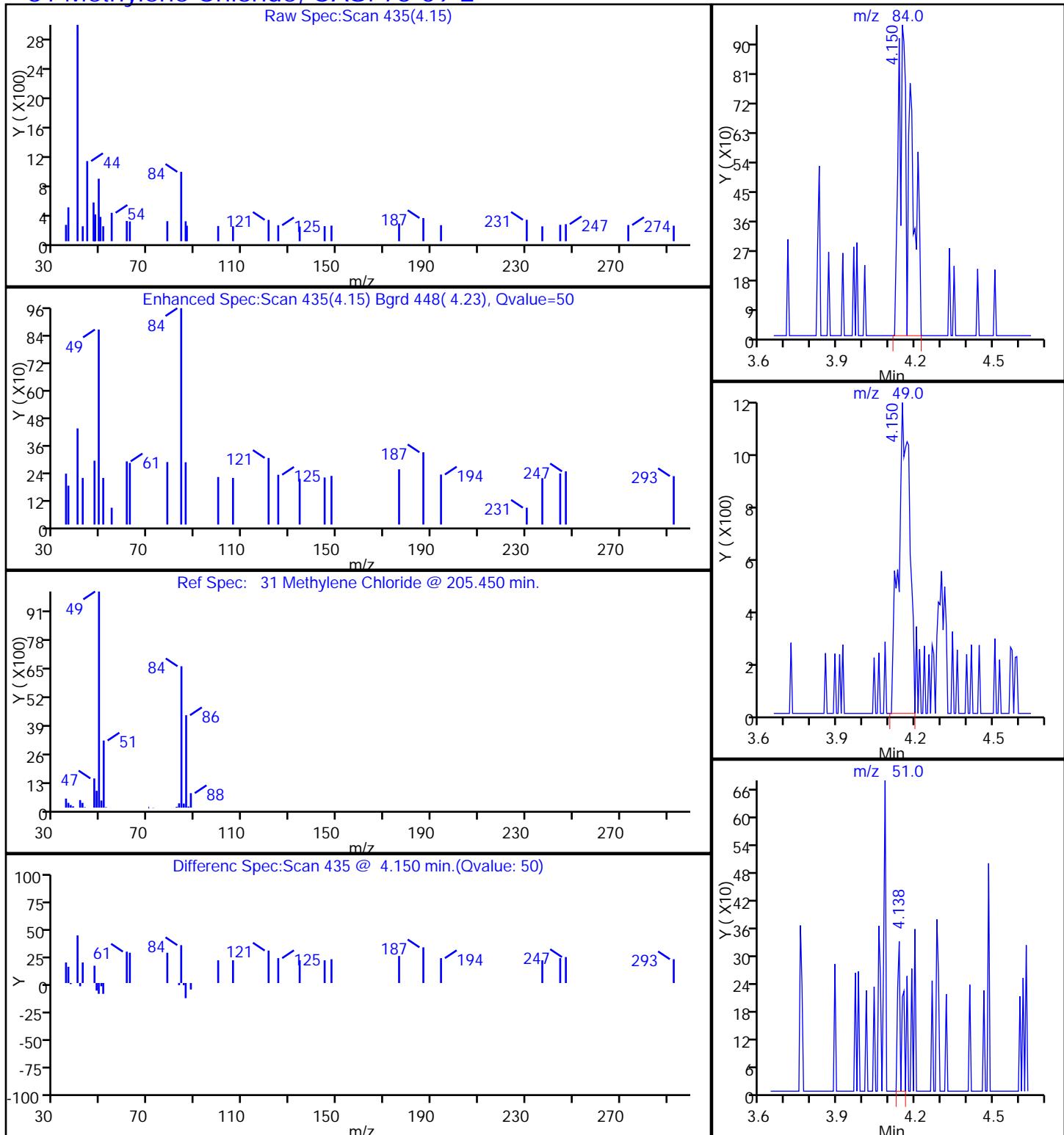
Data File:	\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415007.D	Instrument ID:	CHHP5	Operator ID:	001562
Injection Date:	15-Apr-2015 15:33:30	Lab Sample ID:	180-42975-4	Worklist Smp#:	7
Lims ID:	180-42975-A-4	Dil. Factor:	1.0000	ALS Bottle#:	7
Client ID:	HD-QC2-01-2	Limit Group:	VOA 8260C ICAL		
Purge Vol:	5.000 mL				
Method:	MSVOA_LL_CHHP5				
Column:	DB-624 ( 0.18 mm)				



## TestAmerica Pittsburgh

Data File: \PITCHROM\ChromData\CHHP5\20150415-6480.b\50415007.D  
 Injection Date: 15-Apr-2015 15:33:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-A-4 Lab Sample ID: 180-42975-4  
 Client ID: HD-QC2-01-2  
 Operator ID: 001562 ALS Bottle#: 7 Worklist Smp#: 7  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 ( 0.18 mm) Detector: MS SCAN

## 31 Methylene Chloride, CAS: 75-09-2



FORM VI  
GC/MS VOA INITIAL CALIBRATION DATA  
INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

Analy Batch No.: 135593

SDG No.: \_\_\_\_\_

Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 03/16/2015 12:41 Calibration End Date: 03/16/2015 16:17 Calibration ID: 22457

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	IC 180-135593/13	50316013.D
Level 2	IC 180-135593/4	50316004.D
Level 3	ICIS 180-135593/5	50316005.D
Level 4	IC 180-135593/6	50316006.D
Level 5	IC 180-135593/7	50316007.D
Level 6	IC 180-135593/8	50316008.D
Level 7	IC 180-135593/9	50316009.D
Level 8	IC 180-135593/10	50316010.D

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5		B	M1	M2								
Dichlorodifluoromethane	0.1981 0.2196	0.2184 0.2203	0.2158 0.2064	0.2176	0.2184	Ave		0.2143			0.1000	3.7		20.0			
Chloromethane	0.3161 0.2913	0.3036 0.2846	0.2971 0.2760	0.3139	0.2839	Ave		0.2958			0.1000	4.9		20.0			
Vinyl chloride	0.3339 0.3170	0.3476 0.3238	0.3406 0.2981	0.3521	0.3317	Ave		0.3306			0.1000	5.3		20.0			
1,3-Butadiene	0.4238 0.3606	0.3989 0.3546	0.3880 0.3243	0.3988	0.3720	Ave		0.3776			0.0100	8.3		20.0			
Bromomethane	0.3177 0.1565	0.2026 0.1546	0.1872 0.1489	0.2009	0.1727	Lin2	0.7885	0.1633			0.0500				0.9910		0.9900
Chloroethane	0.2320 0.2316	0.2215 0.2239	0.2348 0.2259	0.2403	0.2201	Ave		0.2287			0.0500	3.1		20.0			
Dichlorofluoromethane	0.6033 0.4953	0.5246 0.5015	0.5246 0.4874	0.5502	0.4911	Ave		0.5222			0.0100	7.5		20.0			
Trichlorofluoromethane	0.3610 0.3924	0.3936 0.3991	0.4043 0.3800	0.4504	0.3921	Ave		0.3966			0.1000	6.4		20.0			
Ethyl ether	0.2888 0.2638	0.2444 0.2500	0.2576 0.2556	0.2691	0.2633	Ave		0.2615			0.0100	5.2		20.0			
Acrolein	0.0310 0.0323	0.0302 0.0321	0.0313 0.0320	0.0335	0.0318	Ave		0.0318			0.0100	3.1		20.0			
1,1-Dichloroethene	0.3207 0.2859	0.2901 0.2792	0.2822 0.2667	0.2965	0.2853	Ave		0.2883			0.1000	5.4		20.0			
1,1,2-Trichloro-1,2,2-trifluoroethane	0.2914 0.2935	0.2973 0.2885	0.2973 0.2692	0.3100	0.2859	Ave		0.2916			0.1000	4.0		20.0			
Acetone	0.1044 0.1092	0.0964 0.1031	0.0956 0.1001	0.1134	0.0972	Ave		0.1024			0.0500	6.2		20.0			
Iodomethane	0.4015 0.3985	0.4019 0.3989	0.4026 0.3873	0.4200	0.3937	Ave		0.4005			0.0100	2.3		20.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI  
GC/MS VOA INITIAL CALIBRATION DATA  
INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

Analy Batch No.: 135593

SDG No.: \_\_\_\_\_

Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 03/16/2015 12:41 Calibration End Date: 03/16/2015 16:17 Calibration ID: 22457

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5		B	M1	M2								
Carbon disulfide	0.7271 0.7017	0.7065 0.6934	0.7209 0.6551	0.7444	0.6916	Ave		0.7051			0.1000	3.8		20.0			
Allyl chloride	0.1408 0.1596	0.1433 0.1659	0.1497 0.1554	0.1576	0.1468	Ave		0.1524			0.0100	5.7		20.0			
Methyl acetate	0.2499 0.2446	0.2206 0.2371	0.2383 0.2398	0.2500	0.2368	Ave		0.2396			0.1000	3.9		20.0			
Methylene Chloride	0.4921 0.3038	0.3340 0.2965	0.3132 0.2964	0.3223	0.3092	Ave		0.3335			0.1000	20.0		20.0			
tert-Butyl alcohol	1.4634 1.1634	1.1166 1.0879	1.2271 1.0609	1.1679	1.1362	Ave		1.1779			0.0100	11.0		20.0			
Acrylonitrile	0.1262 0.1243	0.1185 0.1210	0.1238 0.1200	0.1302	0.1222	Ave		0.1233			0.0100	3.0		20.0			
trans-1,2-Dichloroethene	0.3010 0.2955	0.3039 0.2920	0.2999 0.2846	0.3158	0.2932	Ave		0.2982			0.1000	3.1		20.0			
Methyl tert-butyl ether	0.7046 0.6848	0.5895 0.6670	0.6262 0.6870	0.6643	0.6513	Ave		0.6593			0.1000	5.6		20.0			
Hexane	0.5105 0.4724	0.4808 0.4625	0.4867 0.4447	0.4928	0.4612	Ave		0.4764			0.0100	4.3		20.0			
1,1-Dichloroethane	0.5210 0.5346	0.5355 0.5274	0.5415 0.5173	0.5479	0.5333	Ave		0.5323			0.2000	1.9		20.0			
Vinyl acetate	0.3354 0.4226	0.3143 0.4225	0.3492 0.4312	0.3701	0.3754	Ave		0.3776			0.0100	12.0		20.0			
2,2-Dichloropropane	0.1102 0.1425	0.1245 0.1427	0.1303 0.1457	0.1368	0.1319	Ave		0.1331			0.0100	8.8		20.0			
cis-1,2-Dichloroethene	0.3333 0.3114	0.3188 0.3041	0.3064 0.2999	0.3262	0.3133	Ave		0.3142			0.1000	3.6		20.0			
2-Butanone (MEK)	0.1479 0.1689	0.1544 0.1707	0.1682 0.1707	0.1629	0.1664	Ave		0.1638			0.0500	5.1		20.0			
Bromochloromethane	0.1516 0.1369	0.1328 0.1312	0.1322 0.1303	0.1382	0.1345	Ave		0.1360			0.0100	5.1		20.0			
Tetrahydrofuran	0.1048 0.1057	0.0960 0.1019	0.1025 0.1042	0.1047	0.1007	Ave		0.1026			0.0100	3.0		20.0			
Chloroform	0.5131 0.4845	0.4800 0.4679	0.4876 0.4593	0.4976	0.4787	Ave		0.4836			0.2000	3.5		20.0			
1,1,1-Trichloroethane	0.2755 0.3251	0.2860 0.3242	0.3106 0.3133	0.3267	0.3088	Ave		0.3088			0.1000	6.1		20.0			
Cyclohexane	0.6382 0.5901	0.5930 0.5765	0.5992 0.5384	0.6258	0.5817	Ave		0.5929			0.1000	5.2		20.0			
Carbon tetrachloride	0.2289 0.2566	0.2357 0.2582	0.2463 0.2549	0.2561	0.2457	Ave		0.2478			0.1000	4.4		20.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI  
GC/MS VOA INITIAL CALIBRATION DATA  
INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

Analy Batch No.: 135593

SDG No.: \_\_\_\_\_

Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 03/16/2015 12:41 Calibration End Date: 03/16/2015 16:17 Calibration ID: 22457

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5		B	M1	M2								
1,1-Dichloropropene	0.4232 0.3982	0.4094 0.3936	0.4088 0.3740	0.4106	0.3910	Ave		0.4011			0.0100	3.8		20.0			
Isobutyl alcohol	0.0062 0.0078	0.0044 0.0073	0.0062 0.0083	0.0069	0.0062	Ave		0.0067		*	0.0100	18.0		20.0			
Benzene	1.2964 1.1615	1.1929 1.1190	1.2156 1.0805	1.2375	1.1757	Ave		1.1849			0.5000	5.7		20.0			
1,2-Dichloroethane	0.3889 0.3972	0.3853 0.3828	0.3878 0.3740	0.4029	0.3849	Ave		0.3880			0.1000	2.3		20.0			
n-Heptane	0.4132 0.4165	0.4157 0.3968	0.4135 0.3813	0.4223	0.3971	Ave		0.4071			0.0100	3.4		20.0			
Trichloroethene	0.3236 0.2977	0.2885 0.2885	0.3022 0.2774	0.3045	0.2926	Ave		0.2969			0.2000	4.7		20.0			
Methylcyclohexane	0.5265 0.5361	0.5390 0.5114	0.5500 0.4900	0.5614	0.5230	Ave		0.5297			0.1000	4.2		20.0			
1,2-Dichloropropane	0.2976 0.3036	0.2675 0.2956	0.2870 0.2962	0.2996	0.2976	Ave		0.2931			0.1000	3.9		20.0			
Dibromomethane	0.1682 0.1567	0.1491 0.1563	0.1532 0.1546	0.1640	0.1603	Ave		0.1578			0.0100	3.9		20.0			
1,4-Dioxane	0.0033 0.0034	0.0029 0.0031	0.0029 0.0030	0.0032	0.0030	Ave		0.0031		*	0.0100	5.9		20.0			
Bromodichloromethane	0.2966 0.3370	0.3114 0.3262	0.3286 0.3235	0.3266	0.3259	Ave		0.3220			0.2000	3.9		20.0			
cis-1,3-Dichloropropene	0.2720 0.3463	0.2598 0.3498	0.2835 0.3541	0.3106	0.3095	Ave		0.3107			0.2000	12.0		20.0			
4-Methyl-2-pentanone (MIBK)	1.2503 1.3434	1.2818 1.3687	1.4091 1.3065	1.4145	1.4492	Ave		1.3529			0.1000	5.2		20.0			
Toluene	5.9882 4.5343	5.4946 4.5939	5.5890 4.1718	5.4186	5.2011	Ave		5.1239			0.4000	12.0		20.0			
trans-1,3-Dichloropropene	0.8645 0.9716	0.7455 1.0385	0.8963 1.0484	0.8911	0.9475	Ave		0.9254			0.1000	11.0		20.0			
Ethyl methacrylate	1.1000 1.2637	0.9953 1.3239	1.1753 1.3175	1.1818	1.2989	Ave		1.2070			0.0100	9.7		20.0			
1,1,2-Trichloroethane	1.0794 0.8993	0.9278 0.9152	1.0316 0.8752	0.9797	0.9793	Ave		0.9609			0.1000	7.3		20.0			
Tetrachloroethene	1.1314 0.9214	1.0730 0.9231	1.0654 0.8552	1.0357	1.0130	Ave		1.0023			0.2000	9.3		20.0			
1,3-Dichloropropane	1.9127 1.6507	1.8290 1.6948	1.9187 1.6444	1.8257	1.8122	Ave		1.7860			0.0100	6.1		20.0			
2-Hexanone	0.8865 1.0653	0.9324 1.1043	1.1169 1.0437	1.0718	1.0506	Ave		1.0339			0.1000	7.9		20.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI  
GC/MS VOA INITIAL CALIBRATION DATA  
INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

Analy Batch No.: 135593

SDG No.:

Instrument ID: CHHP5

GC Column: DB-624

ID: 0.18 (mm)

Heated Purge: (Y/N) N

Calibration Start Date: 03/16/2015 12:41

Calibration End Date: 03/16/2015 16:17

Calibration ID: 22457

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5		B	M1	M2								
Dibromochloromethane	0.6589 0.7861	0.7302 0.8019	0.7961 0.7741	0.7741	0.8146	Ave		0.7670			0.1000	6.6		20.0			
1,2-Dibromoethane (EDB)	0.9462 0.8909	0.8507 0.9041	0.9478 0.8836	0.9579	0.9540	Ave		0.9169			0.1000	4.4		20.0			
3-Chlorobenzotrifluoride	2.1568 1.7885	2.0616 1.8999	2.0657 1.6136	2.0676	1.9855	Ave		1.9549			0.0100	9.2		20.0			
Chlorobenzene	3.9165 2.9120	3.3811 2.9538	3.4265 2.7856	3.3185	3.2780	Ave		3.2465			0.5000	11.0		20.0			
4-Chlorobenzotrifluoride	2.1386 1.7554	1.9292 1.8762	1.9271 1.5481	1.9634	1.9831	Ave		1.8901			0.0100	9.2		20.0			
1,1,1,2-Tetrachloroethane	0.7551 0.8493	0.8012 0.8680	0.8363 0.8428	0.8482	0.9047	Ave		0.8382			0.0100	5.3		20.0			
Ethylbenzene	1.9914 1.7179	1.9333 1.7672	1.9980 1.6464	1.9518	1.8953	Ave		1.8627			0.1000	7.2		20.0			
m-Xylene & p-Xylene	2.4849 2.1093	2.3674 2.1267	2.4171 1.9994	2.4234	2.2969	Ave		2.2781			0.1000	7.8		20.0			
o-Xylene	2.6403 2.0475	2.2064 2.0545	2.3516 1.9292	2.3257	2.2716	Ave		2.2283			0.3000	10.0		20.0			
Styrene	3.8818 3.3296	3.6611 3.3147	3.8658 3.1277	3.7940	3.7504	Ave		3.5907			0.3000	8.1		20.0			
Bromoform	0.4254 0.4898	0.4398 0.4974	0.4744 0.4894	0.4822	0.4911	Ave		0.4737			0.1000	5.6		20.0			
2-Chlorobenzotrifluoride	2.0985 1.7811	2.0764 1.8958	2.0751 1.6078	2.0615	2.0224	Ave		1.9523			0.0100	9.1		20.0			
Isopropylbenzene	6.2252 4.9838	6.1153 4.8827	6.0965 4.4013	6.0579	5.7184	Ave		5.5601			0.1000	13.0		20.0			
1,1,2,2-Tetrachloroethane	1.5778 1.3165	1.3921 1.3063	1.4139 1.2430	1.4088	1.3646	Ave		1.3779			0.3000	7.2		20.0			
Bromobenzene	0.9601 0.9043	0.9163 0.9102	0.9670 0.9012	0.9241	0.9202	Ave		0.9254			0.0100	2.7		20.0			
1,2,3-Trichloropropane	0.3380 0.3040	0.2838 0.2874	0.3205 0.3069	0.2961	0.2961	Ave		0.3041			0.0100	5.9		20.0			
trans-1,4-Dichloro-2-butene	0.2572 0.2562	0.2443 0.2601	0.2456 0.2696	0.2438	0.2460	Ave		0.2528			0.0100	3.7		20.0			
N-Propylbenzene	1.2305 1.1066	1.1620 1.0908	1.2081 1.0656	1.1555	1.1135	Ave		1.1416			0.0100	5.1		20.0			
2-Chlorotoluene	1.0248 0.9458	0.9575 0.9297	1.0195 0.9076	0.9558	0.9319	Ave		0.9591			0.0100	4.4		20.0			
3-Chlorotoluene	1.1523 1.0737	1.0357 1.0942	1.0635 0.9927	1.0618	1.1018	Ave		1.0720			0.0100	4.4		20.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI  
GC/MS VOA INITIAL CALIBRATION DATA  
INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1 Analy Batch No.: 135593

SDG No.: \_\_\_\_\_

Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 03/16/2015 12:41 Calibration End Date: 03/16/2015 16:17 Calibration ID: 22457

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5		B	M1	M2								
1,3,5-Trimethylbenzene	3.5091 3.0301	3.2905 2.9338	3.3765 2.8104	3.3601	3.1525	Ave		3.1829			0.0100	7.6		20.0			
4-Chlorotoluene	1.1316 1.0097	1.0151 0.9774	1.0863 1.0009	1.0825	1.0021	Ave		1.0382			0.0100	5.2		20.0			
tert-Butylbenzene	3.1830 2.5928	2.8173 2.5318	2.9656 2.3701	2.8959	2.7052	Ave		2.7577			0.0100	9.5		20.0			
1,2,4-Trimethylbenzene	3.6039 3.1029	3.3270 3.0238	3.4986 2.8908	3.4674	3.2206	Ave		3.2669			0.0100	7.7		20.0			
3,4-Dichlorobenzotrifluoride	1.1042 1.0202	0.9953 1.0227	1.0269 0.9335	1.1049	1.0507	Ave		1.0323			0.0100	5.5		20.0			
sec-Butylbenzene	4.3054 3.6389	4.1108 3.5066	4.1991 3.2620	4.1487	3.8794	Ave		3.8814			0.0100	9.7		20.0			
1,3-Dichlorobenzene	1.9132 1.6438	1.7258 1.6071	1.7369 1.5897	1.7497	1.6725	Ave		1.7048			0.6000	6.1		20.0			
4-Isopropyltoluene	3.4872 3.0606	3.2348 2.9586	3.4694 2.7984	3.4562	3.1691	Ave		3.2043			0.0100	8.0		20.0			
1,4-Dichlorobenzene	1.9760 1.6976	1.7145 1.6569	1.7807 1.6355	1.7648	1.7035	Ave		1.7412			0.5000	6.1		20.0			
2,4-Dichlorobenzotrifluoride	1.0162 0.9585	0.9307 0.9665	1.0004 0.8567	1.0551	0.9508	Ave		0.9669			0.0100	6.2		20.0			
2,5-Dichlorobenzotrifluoride	1.1811 1.0613	1.0765 1.0776	1.0685 0.9818	1.1269	1.0793	Ave		1.0816			0.0100	5.2		20.0			
n-Butylbenzene	3.1276 2.8128	2.9811 2.7148	3.1079 2.5582	3.1414	2.9001	Ave		2.9180			0.0100	7.3		20.0			
1,2-Dichlorobenzene	1.7371 1.5488	1.5543 1.5042	1.6235 1.4749	1.6066	1.5803	Ave		1.5787			0.4000	5.1		20.0			
1,2-Dibromo-3-Chloropropane	0.1313 0.1386	0.1067 0.1383	0.1229 0.1385	0.1324	0.1248	Ave		0.1292			0.0500	8.5		20.0			
1,2,4-Trichlorobenzene	0.9720 0.8625	0.7083 0.8349	0.7579 0.7778	0.8780	0.7835	Ave		0.8219			0.2000	10.0		20.0			
Hexachlorobutadiene	0.4883 0.3899	0.3825 0.3778	0.3866 0.3464	0.4091	0.3724	Ave		0.3941			0.0100	11.0		20.0			
Naphthalene	2.3899 2.2683	1.8332 2.1948	1.9931 2.0920	2.3983	2.0941	Ave		2.1580			0.0100	9.0		20.0			
1,2,3-Trichlorobenzene	0.7895 0.7155	0.5376 0.7162	0.6024 0.6573	0.7303	0.6432	Ave		0.6740			0.0100	12.0		20.0			
2,4,5-Trichlorotoluene	0.4907 0.3881	0.2750 0.3876	0.2929 0.3431	0.3938	0.3283	Ave		0.3624			0.0100	19.0		20.0			
2,3,6-Trichlorotoluene	0.4374 0.3491	0.2501 0.3509	0.2713 0.3051	0.3608	0.2936	Ave		0.3273			0.0100	18.0		20.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI  
GC/MS VOA INITIAL CALIBRATION DATA  
INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1 Analy Batch No.: 135593

SDG No.: \_\_\_\_\_

Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 03/16/2015 12:41 Calibration End Date: 03/16/2015 16:17 Calibration ID: 22457

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5		B	M1	M2								
Dibromofluoromethane (Surr)	0.2497 0.2212	0.2276 0.2219	0.2284 0.2143	0.2334	0.2228	Ave		0.2274				4.7		20.0			
1,2-Dichloroethane-d4 (Surr)	0.3017 0.2995	0.3055 0.2914	0.3015 0.2867	0.3115	0.3008	Ave		0.2998				2.6		20.0			
Toluene-d8 (Surr)	4.5313 3.5890	4.2126 3.6439	4.3365 3.2599	4.2301	4.0882	Ave		3.9864				11.0		20.0			
4-Bromofluorobenzene (Surr)	1.5722 1.3558	1.4371 1.3519	1.5107 1.2944	1.4891	1.4730	Ave		1.4356				6.6		20.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI  
GC/MS VOA INITIAL CALIBRATION DATA  
INTERNAL STANDARD RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

Analy Batch No.: 135593

SDG No.: \_\_\_\_\_

Instrument ID: CHHP5      GC Column: DB-624      ID: 0.18 (mm)      Heated Purge: (Y/N) N

Calibration Start Date: 03/16/2015 12:41      Calibration End Date: 03/16/2015 16:17      Calibration ID: 22457

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	IC 180-135593/13	50316013.D
Level 2	IC 180-135593/4	50316004.D
Level 3	ICIS 180-135593/5	50316005.D
Level 4	IC 180-135593/6	50316006.D
Level 5	IC 180-135593/7	50316007.D
Level 6	IC 180-135593/8	50316008.D
Level 7	IC 180-135593/9	50316009.D
Level 8	IC 180-135593/10	50316010.D

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (NG)				
			LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5
Dichlorodifluoromethane	FB	Ave	11265 432190	59394 522240	116111 640090	173113	243823	5.00 175	25.0 200	50.0 250	75.0	100
Chloromethane	FB	Ave	17972 573343	82552 674845	159885 855933	249772	316915	5.00 175	25.0 200	50.0 250	75.0	100
Vinyl chloride	FB	Ave	18981 624000	94520 767804	183317 924535	280135	370271	5.00 175	25.0 200	50.0 250	75.0	100
1,3-Butadiene	FB	Ave	24095 709784	108469 840803	208815 1005925	317272	415323	5.00 175	25.0 200	50.0 250	75.0	100
Bromomethane	FB	Lin2	18060 307964	55097 366671	100717 461680	159846	192846	5.00 175	25.0 200	50.0 250	75.0	100
Chloroethane	FB	Ave	13187 455903	60248 530813	126349 700467	191164	245673	5.00 175	25.0 200	50.0 250	75.0	100
Dichlorofluoromethane	FB	Ave	34297 974888	142662 1188936	282324 1511714	437737	548270	5.00 175	25.0 200	50.0 250	75.0	100
Trichlorofluoromethane	FB	Ave	20521 772293	107038 946313	217544 1178605	358375	437688	5.00 175	25.0 200	50.0 250	75.0	100
Ethyl ether	FB	Ave	16416 519119	66452 592652	138609 792637	214135	293889	5.00 175	25.0 200	50.0 250	75.0	100
Acrolein	FB	Ave	35289 81646	41017 95028	50582 109180	62132	71073	100 225	125 250	150 275	175	200
1,1-Dichloroethene	FB	Ave	18234 562804	78897 662050	151843 827120	235889	318457	5.00 175	25.0 200	50.0 250	75.0	100
1,1,2-Trichloro-1,2,2-trifluoroethane	FB	Ave	16567 577719	80854 684103	159979 834802	246660	319162	5.00 175	25.0 200	50.0 250	75.0	100
Acetone	FB	Ave	29674 429781	52410 489133	102899 621064	180387	217095	25.0 350	50.0 400	100 500	150	200
Iodomethane	FB	Ave	22824 784350	109309 945860	216640 1201056	334141	439512	5.00 175	25.0 200	50.0 250	75.0	100
Carbon disulfide	FB	Ave	41336 1381152	192118 1643948	387934 2031733	592248	772081	5.00 175	25.0 200	50.0 250	75.0	100

FORM VI  
GC/MS VOA INITIAL CALIBRATION DATA  
INTERNAL STANDARD RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

Analy Batch No.: 135593

SDG No.:

Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 03/16/2015 12:41 Calibration End Date: 03/16/2015 16:17 Calibration ID: 22457

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (NG)				
			LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5
Allyl chloride	FB	Ave	8006 314052	38957 393345	80577 482122	125423	163875	5.00 175	25.0 200	50.0 250	75.0	100
Methyl acetate	FB	Ave	71022 2407305	299965 2810332	641136 3718382	994505	1321970	25.0 875	125 1000	250 1250	375	500
Methylene Chloride	FB	Ave	27978 597904	90836 703059	168570 919183	256424	345226	5.00 175	25.0 200	50.0 250	75.0	100
tert-Butyl alcohol	TBA	Ave	10830 351016	39251 399281	83098 537174	133756	175500	50.0 1750	250 2000	500 2500	750	1000
Acrylonitrile	FB	Ave	71728 2446379	322268 2868164	666088 3721902	1035956	1363975	50.0 1750	250 2000	500 2500	750	1000
trans-1,2-Dichloroethene	FB	Ave	17111 581552	82640 692220	161381 882651	251288	327278	5.00 175	25.0 200	50.0 250	75.0	100
Methyl tert-butyl ether	FB	Ave	40058 1347848	160325 1581345	336961 2130684	528520	727030	5.00 175	25.0 200	50.0 250	75.0	100
Hexane	FB	Ave	29021 929791	130741 1096478	261916 1379168	392065	514868	5.00 175	25.0 200	50.0 250	75.0	100
1,1-Dichloroethane	FB	Ave	29622 1052201	145639 1250453	291408 1604398	435915	595324	5.00 175	25.0 200	50.0 250	75.0	100
Vinyl acetate	FB	Ave	19067 831670	85462 1001771	187915 1337263	294456	419086	5.00 175	25.0 200	50.0 250	75.0	100
2,2-Dichloropropane	FB	Ave	6267 280515	33850 338302	70106 452022	108858	147216	5.00 175	25.0 200	50.0 250	75.0	100
cis-1,2-Dichloroethene	FB	Ave	18951 612812	86701 721075	164893 930230	259517	349805	5.00 175	25.0 200	50.0 250	75.0	100
2-Butanone (MEK)	FB	Ave	42054 665013	83987 809232	180996 1059138	259227	371447	25.0 350	50.0 400	100 500	150	200
Bromochloromethane	FB	Ave	8619 269375	36107 311076	71124 404105	109930	150204	5.00 175	25.0 200	50.0 250	75.0	100
Tetrahydrofuran	FB	Ave	11913 415944	52231 483324	110274 646482	166594	224920	10.0 350	50.0 400	100 500	150	200
Chloroform	FB	Ave	29168 953676	130523 1109416	262371 1424461	395935	534362	5.00 175	25.0 200	50.0 250	75.0	100
1,1,1-Trichloroethane	FB	Ave	15663 639960	77770 768585	167130 971626	259963	344772	5.00 175	25.0 200	50.0 250	75.0	100
Cyclohexane	FB	Ave	36280 1161488	161271 1366913	322468 1669676	497889	649387	5.00 175	25.0 200	50.0 250	75.0	100
Carbon tetrachloride	FB	Ave	13013 504991	64089 612080	132517 790495	203736	274328	5.00 175	25.0 200	50.0 250	75.0	100
1,1-Dichloropropene	FB	Ave	24060 783682	111342 933326	219974 1159811	326699	436454	5.00 175	25.0 200	50.0 250	75.0	100
Isobutyl alcohol	FB	Ave	8820 386141	29897 433313	83109 644697	137203	174166	125 4375	625 5000	1250 6250	1875	2500

FORM VI  
GC/MS VOA INITIAL CALIBRATION DATA  
INTERNAL STANDARD RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

Analy Batch No.: 135593

SDG No.: \_\_\_\_\_

Instrument ID: CHHP5      GC Column: DB-624      ID: 0.18 (mm)      Heated Purge: (Y/N) N

Calibration Start Date: 03/16/2015 12:41      Calibration End Date: 03/16/2015 16:17      Calibration ID: 22457

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (NG)				
			LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5
Benzene	FB	Ave	73700 2286079	324419 2653105	654151 3351151	984614	1312435	5.00 175	25.0 200	50.0 250	75.0	100
1,2-Dichloroethane	FB	Ave	22108 781760	104777 907622	208683 1159879	320594	429724	5.00 175	25.0 200	50.0 250	75.0	100
n-Heptane	FB	Ave	23490 819785	113041 940924	222515 1182643	335961	443357	5.00 175	25.0 200	50.0 250	75.0	100
Trichloroethene	FB	Ave	18397 586010	78459 684010	162608 860273	242252	326599	5.00 175	25.0 200	50.0 250	75.0	100
Methylcyclohexane	FB	Ave	29934 1055175	146574 1212427	295972 1519674	446628	583894	5.00 175	25.0 200	50.0 250	75.0	100
1,2-Dichloropropane	FB	Ave	16916 597514	72742 700921	154467 918714	238331	332279	5.00 175	25.0 200	50.0 250	75.0	100
Dibromomethane	FB	Ave	9562 308441	40542 370624	82469 479407	130496	178905	5.00 175	25.0 200	50.0 250	75.0	100
1,4-Dioxane	FB	Ave	3746 132396	15563 146272	31354 185631	50907	66490	100 3500	500 4000	1000 5000	1500	2000
Bromodichloromethane	FB	Ave	16863 663337	84673 773432	176851 1003399	259871	363842	5.00 175	25.0 200	50.0 250	75.0	100
cis-1,3-Dichloropropene	FB	Ave	15462 681682	70642 829306	152581 1098242	247138	345528	5.00 175	25.0 200	50.0 250	75.0	100
4-Methyl-2-pentanone (MIBK)	CBZ	Ave	75787 1390980	154453 1617802	342539 2109966	531084	747218	25.0 350	50.0 400	100 500	150	200
Toluene	CBZ	Ave	72597 2347437	331041 2714932	679332 3368812	1017198	1340817	5.00 175	25.0 200	50.0 250	75.0	100
trans-1,3-Dichloropropene	CBZ	Ave	10481 502980	44917 613747	108942 846559	167274	244258	5.00 175	25.0 200	50.0 250	75.0	100
Ethyl methacrylate	CBZ	Ave	13336 654210	59964 782394	142858 1063861	221852	334858	5.00 175	25.0 200	50.0 250	75.0	100
1,1,2-Trichloroethane	CBZ	Ave	13086 465584	55897 540864	125390 706748	183907	252461	5.00 175	25.0 200	50.0 250	75.0	100
Tetrachloroethylene	CBZ	Ave	13716 477004	64647 545517	129494 690601	194422	261148	5.00 175	25.0 200	50.0 250	75.0	100
1,3-Dichloropropane	CBZ	Ave	23188 854593	110194 1001573	233217 1327847	342719	467174	5.00 175	25.0 200	50.0 250	75.0	100
2-Hexanone	CBZ	Ave	53734 1103034	112348 1305223	271508 1685534	402386	541680	25.0 350	50.0 400	100 500	150	200
Dibromochloromethane	CBZ	Ave	7988 406960	43996 473922	96762 625118	145315	210013	5.00 175	25.0 200	50.0 250	75.0	100
1,2-Dibromoethane (EDB)	CBZ	Ave	11471 461219	51254 534328	115204 713501	179814	245946	5.00 175	25.0 200	50.0 250	75.0	100
3-Chlorobenzotrifluoride	CBZ	Ave	26148 925933	124209 1122812	251080 1303041	388132	511845	5.00 175	25.0 200	50.0 250	75.0	100

FORM VI  
GC/MS VOA INITIAL CALIBRATION DATA  
INTERNAL STANDARD RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

Analy Batch No.: 135593

SDG No.:

Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 03/16/2015 12:41 Calibration End Date: 03/16/2015 16:17 Calibration ID: 22457

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (NG)				
			LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5
Chlorobenzene	CBZ	Ave	47481 1507544	203702 1745676	416488 2249414	622968	845046	5.00 175	25.0 200	50.0 250	75.0	100
4-Chlorobenzotrifluoride	CBZ	Ave	25927 908777	116232 1108797	234233 1250140	368570	511237	5.00 175	25.0 200	50.0 250	75.0	100
1,1,1,2-Tetrachloroethane	CBZ	Ave	9154 439701	48269 512980	101650 680608	159225	233228	5.00 175	25.0 200	50.0 250	75.0	100
Ethylbenzene	CBZ	Ave	24142 889389	116477 1044399	242856 1329470	366398	488611	5.00 175	25.0 200	50.0 250	75.0	100
m-Xylene & p-Xylene	CBZ	Ave	30126 1092005	142634 1256840	293796 1614511	454933	592135	5.00 175	25.0 200	50.0 250	75.0	100
o-Xylene	CBZ	Ave	32009 1059986	132929 1214164	285835 1557898	436586	585609	5.00 175	25.0 200	50.0 250	75.0	100
Styrene	CBZ	Ave	47061 1723778	220574 1958961	469890 2525667	712222	966850	5.00 175	25.0 200	50.0 250	75.0	100
Bromoform	CBZ	Ave	5157 253560	26498 293938	57667 395201	90522	126605	5.00 175	25.0 200	50.0 250	75.0	100
2-Chlorobenzotrifluoride	CBZ	Ave	25441 922108	125099 1120386	252226 1298335	386985	521379	5.00 175	25.0 200	50.0 250	75.0	100
Isopropylbenzene	CBZ	Ave	75470 2580136	368436 2885608	741027 3554151	1137215	1474178	5.00 175	25.0 200	50.0 250	75.0	100
1,1,2,2-Tetrachloroethane	CBZ	Ave	19128 681581	83874 772016	171864 1003707	264462	351798	5.00 175	25.0 200	50.0 250	75.0	100
Bromobenzene	DCB	Ave	16809 637569	80670 740842	168649 956763	253502	346996	5.00 175	25.0 200	50.0 250	75.0	100
1,2,3-Trichloropropane	DCB	Ave	5918 214358	24990 233938	55900 325768	81225	111668	5.00 175	25.0 200	50.0 250	75.0	100
trans-1,4-Dichloro-2-butene	DCB	Ave	4503 180624	21505 211691	42827 286166	66879	92761	5.00 175	25.0 200	50.0 250	75.0	100
N-Propylbenzene	DCB	Ave	21543 780243	102304 887838	210687 1131297	316980	419888	5.00 175	25.0 200	50.0 250	75.0	100
2-Chlorotoluene	DCB	Ave	17942 666866	84295 756732	177793 963573	262207	351403	5.00 175	25.0 200	50.0 250	75.0	100
3-Chlorotoluene	DCB	Ave	20174 757051	91182 890638	185477 1053875	291288	415463	5.00 175	25.0 200	50.0 250	75.0	100
1,3,5-Trimethylbenzene	DCB	Ave	61438 2136446	289696 2387945	588847 2983647	921783	1188743	5.00 175	25.0 200	50.0 250	75.0	100
4-Chlorotoluene	DCB	Ave	19812 711885	89370 795532	189449 1062581	296950	377870	5.00 175	25.0 200	50.0 250	75.0	100
tert-Butylbenzene	DCB	Ave	55729 1828125	248042 2060731	517188 2516209	794422	1020106	5.00 175	25.0 200	50.0 250	75.0	100
1,2,4-Trimethylbenzene	DCB	Ave	63098 2187785	292909 2461131	610150 3068942	951216	1214438	5.00 175	25.0 200	50.0 250	75.0	100

FORM VI  
GC/MS VOA INITIAL CALIBRATION DATA  
INTERNAL STANDARD RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Pittsburgh

Job No.: 180-42975-1

Analy Batch No.: 135593

SDG No.: \_\_\_\_\_

Instrument ID: CHHP5      GC Column: DB-624      ID: 0.18 (mm)      Heated Purge: (Y/N) N

Calibration Start Date: 03/16/2015 12:41      Calibration End Date: 03/16/2015 16:17      Calibration ID: 22457

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (NG)				
			LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5
3,4-Dichlorobenzotrifluoride	DCB	Ave	19333 719294	87627 832435	179092 991010	303120	396211	5.00 175	25.0 200	50.0 250	75.0	100
sec-Butylbenzene	DCB	Ave	75379 2565671	361915 2854173	732318 3463106	1138120	1462842	5.00 175	25.0 200	50.0 250	75.0	100
1,3-Dichlorobenzene	DCB	Ave	33497 1159025	151937 1308081	302903 1687649	480001	630675	5.00 175	25.0 200	50.0 250	75.0	100
4-Isopropyltoluene	DCB	Ave	61054 2157955	284792 2408127	605051 2970922	948139	1195021	5.00 175	25.0 200	50.0 250	75.0	100
1,4-Dichlorobenzene	DCB	Ave	34596 1196958	150942 1348596	310551 1736319	484138	642365	5.00 175	25.0 200	50.0 250	75.0	100
2,4-Dichlorobenzotrifluoride	DCB	Ave	17792 675783	81937 786683	174468 909481	289446	358539	5.00 175	25.0 200	50.0 250	75.0	100
2,5-Dichlorobenzotrifluoride	DCB	Ave	20678 748317	94772 877059	186350 1042359	309155	406971	5.00 175	25.0 200	50.0 250	75.0	100
n-Butylbenzene	DCB	Ave	54758 1983203	262455 2209671	542017 2715831	861784	1093564	5.00 175	25.0 200	50.0 250	75.0	100
1,2-Dichlorobenzene	DCB	Ave	30414 1092014	136843 1224311	283138 1565775	440732	595901	5.00 175	25.0 200	50.0 250	75.0	100
1,2-Dibromo-3-Chloropropane	DCB	Ave	2299 97714	9396 112547	21428 147059	36318	47067	5.00 175	25.0 200	50.0 250	75.0	100
1,2,4-Trichlorobenzene	DCB	Ave	17018 608110	62363 679520	132179 825772	240861	295444	5.00 175	25.0 200	50.0 250	75.0	100
Hexachlorobutadiene	DCB	Ave	8549 274932	33676 307470	67414 367792	112236	140410	5.00 175	25.0 200	50.0 250	75.0	100
Naphthalene	DCB	Ave	41842 1599300	161398 1786434	347596 2220927	657935	789643	5.00 175	25.0 200	50.0 250	75.0	100
1,2,3-Trichlorobenzene	DCB	Ave	13823 504504	47333 582911	105062 697862	200345	242534	5.00 175	25.0 200	50.0 250	75.0	100
2,4,5-Trichlorotoluene	DCB	Ave	8592 273662	24209 315499	51080 364223	108037	123791	5.00 175	25.0 200	50.0 250	75.0	100
2,3,6-Trichlorotoluene	DCB	Ave	7658 246163	22020 285573	47319 323920	98974	110702	5.00 175	25.0 200	50.0 250	75.0	100
Dibromofluoromethane (Surr)	FB	Ave	14193 435320	61901 526164	122918 664693	185698	248750	5.00 175	25.0 200	50.0 250	75.0	100
1,2-Dichloroethane-d4 (Surr)	FB	Ave	17152 589491	83077 691002	162227 889045	247858	335757	5.00 175	25.0 200	50.0 250	75.0	100
Toluene-d8 (Surr)	CBZ	Ave	54935 1858068	253798 2153477	527093 2632400	794092	1053927	5.00 175	25.0 200	50.0 250	75.0	100
4-Bromofluorobenzene (Surr)	CBZ	Ave	19061 701915	86585 798953	183629 1045249	279546	379740	5.00 175	25.0 200	50.0 250	75.0	100

FORM VI  
GC/MS VOA INITIAL CALIBRATION DATA  
INTERNAL STANDARD RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1 Analy Batch No.: 135593

SDG No.: \_\_\_\_\_

Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 03/16/2015 12:41 Calibration End Date: 03/16/2015 16:17 Calibration ID: 22457

Curve Type Legend:

Ave = Average ISTD

Lin2 = Linear 1/conc<sup>2</sup> ISTD

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316004.D  
 Lims ID: IC VSTD5  
 Client ID:  
 Sample Type: IC Calib Level: 2  
 Inject. Date: 16-Mar-2015 12:41:30 ALS Bottle#: 4 Worklist Smp#: 4  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: IC VSTD5  
 Misc. Info.: 180-0006031-004  
 Operator ID: 001562 Instrument ID: CHHP5  
 Sublist: chrom-MSVOA\_LL\_CHHP5\*sub4  
 Method: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 17-Mar-2015 10:59:20 Calib Date: 16-Mar-2015 16:17:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICal File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316013.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK012

First Level Reviewer: fergusond Date: 17-Mar-2015 09:28:19

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.302	4.305	-0.003	88	140612	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.277	7.273	0.004	97	543896	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.361	10.364	-0.003	99	120496	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.685	12.682	0.003	97	176082	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.535	6.525	0.010	94	61901	25.0	25.0	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.906	6.902	0.004	96	83077	25.0	25.5	
\$ 7 Toluene-d8 (Surr)	98	8.919	8.922	-0.003	100	253798	25.0	26.4	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.529	11.532	-0.003	95	86585	25.0	25.0	
11 Dichlorodifluoromethane	85	1.619	1.616	0.003	98	59394	25.0	25.5	
12 Chloromethane	50	1.771	1.774	-0.003	100	82552	25.0	25.7	
13 Vinyl chloride	62	1.905	1.902	0.003	99	94520	25.0	26.3	
14 Butadiene	39	1.948	1.944	0.004	98	108469	25.0	26.4	
15 Bromomethane	94	2.252	2.249	0.003	90	55097	25.0	26.2	M
16 Chloroethane	64	2.392	2.370	0.022	97	60248	25.0	24.2	
17 Dichlorofluoromethane	67	2.660	2.650	0.010	98	142662	25.0	25.1	
18 Trichlorofluoromethane	101	2.690	2.711	-0.021	96	107038	25.0	24.8	
20 Ethyl ether	59	3.085	3.088	-0.003	98	66452	25.0	23.4	
21 Acrolein	56	3.256	3.252	0.004	96	41017	125.0	118.7	
22 1,1-Dichloroethene	96	3.371	3.386	-0.015	97	78897	25.0	25.2	
23 1,1,2-Trichloro-1,2,2-trif	101	3.444	3.429	0.015	97	80854	25.0	25.5	
24 Acetone	43	3.493	3.496	-0.003	98	52410	50.0	47.0	
25 Iodomethane	142	3.572	3.587	-0.015	98	109309	25.0	25.1	
26 Carbon disulfide	76	3.651	3.654	-0.003	100	192118	25.0	25.0	
28 3-Chloro-1-propene	76	3.931	3.940	-0.009	92	38957	25.0	23.5	
30 Methyl acetate	43	4.022	4.019	0.003	100	299965	125.0	115.1	
31 Methylene Chloride	84	4.144	4.134	0.010	95	90836	25.0	25.0	
32 2-Methyl-2-propanol	59	4.430	4.445	-0.015	90	39251	250.0	237.0	
33 Acrylonitrile	53	4.552	4.554	-0.002	100	322268	250.0	240.3	
34 trans-1,2-Dichloroethene	96	4.564	4.560	0.004	61	82640	25.0	25.5	
35 Methyl tert-butyl ether	73	4.594	4.591	0.003	96	160325	25.0	22.4	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
36 Hexane	57	4.983	4.980	0.003	97	130741	25.0	25.2	
37 1,1-Dichloroethane	63	5.178	5.169	0.009	99	145639	25.0	25.2	
38 Vinyl acetate	43	5.300	5.290	0.010	100	85462	25.0	20.8	
44 2,2-Dichloropropane	77	5.926	5.923	0.003	85	33850	25.0	23.4	
45 cis-1,2-Dichloroethene	96	5.939	5.935	0.004	94	86701	25.0	25.4	
46 2-Butanone (MEK)	43	5.987	5.990	-0.003	99	83987	50.0	47.1	
49 Chlorobromomethane	128	6.224	6.233	-0.009	96	36107	25.0	24.4	
51 Tetrahydrofuran	42	6.285	6.288	-0.003	97	52231	50.0	46.8	
52 Chloroform	83	6.340	6.343	-0.003	96	130523	25.0	24.8	
53 1,1,1-Trichloroethane	97	6.529	6.531	-0.002	95	77770	25.0	23.2	
54 Cyclohexane	56	6.589	6.586	0.003	96	161271	25.0	25.0	
56 Carbon tetrachloride	117	6.723	6.720	0.003	69	64089	25.0	23.8	
55 1,1-Dichloropropene	75	6.729	6.726	0.003	96	111342	25.0	25.5	
57 Isobutyl alcohol	41	6.942	6.945	-0.003	33	29897	625.0	411.8	
58 Benzene	78	6.954	6.957	-0.003	98	324419	25.0	25.2	
59 1,2-Dichloroethane	62	6.985	6.981	0.004	98	104777	25.0	24.8	
62 n-Heptane	43	7.277	7.280	-0.003	65	113041	25.0	25.5	
64 Trichloroethene	130	7.666	7.669	-0.003	99	78459	25.0	24.3	
66 Methylcyclohexane	83	7.867	7.864	0.003	96	146574	25.0	25.4	
67 1,2-Dichloropropane	63	7.903	7.906	-0.003	95	72742	25.0	22.8	
68 Dibromomethane	93	8.031	8.028	0.003	94	40542	25.0	23.6	
70 1,4-Dioxane	88	8.068	8.058	0.010	87	15563	500.0	463.6	
71 Dichlorobromomethane	83	8.196	8.198	-0.002	97	84673	25.0	24.2	
74 cis-1,3-Dichloropropene	75	8.658	8.661	-0.002	98	70642	25.0	20.9	
75 4-Methyl-2-pentanone (MIBK)	43	8.822	8.825	-0.003	98	154453	50.0	47.4	
76 Toluene	91	8.992	8.989	0.003	99	331041	25.0	26.8	
77 trans-1,3-Dichloropropene	75	9.224	9.220	0.004	95	44917	25.0	20.1	
78 Ethyl methacrylate	69	9.321	9.318	0.003	97	59964	25.0	20.6	
79 1,1,2-Trichloroethane	97	9.400	9.397	0.003	98	55897	25.0	24.1	
80 Tetrachloroethene	164	9.534	9.537	-0.003	96	64647	25.0	26.8	
81 1,3-Dichloropropane	76	9.570	9.567	0.003	98	110194	25.0	25.6	
82 2-Hexanone	43	9.662	9.658	0.004	99	112348	50.0	45.1	
84 Chlorodibromomethane	129	9.795	9.786	0.009	98	43996	25.0	23.8	
85 Ethylene Dibromide	107	9.899	9.902	-0.003	99	51254	25.0	23.2	
86 3-Chlorobenzotrifluoride	180	10.373	10.370	0.003	89	124209	25.0	26.4	
87 Chlorobenzene	112	10.392	10.394	-0.002	99	203702	25.0	26.0	
88 4-Chlorobenzotrifluoride	180	10.434	10.431	0.003	99	116232	25.0	25.5	
89 1,1,2-Tetrachloroethane	131	10.477	10.473	0.004	96	48269	25.0	23.9	
90 Ethylbenzene	106	10.501	10.498	0.003	100	116477	25.0	25.9	
91 m-Xylene & p-Xylene	106	10.617	10.619	-0.002	99	142634	25.0	26.0	
92 o-Xylene	106	11.012	11.015	-0.003	97	132929	25.0	24.8	
93 Styrene	104	11.024	11.027	-0.003	99	220574	25.0	25.5	
94 Bromoform	173	11.213	11.209	0.004	97	26498	25.0	23.2	
96 2-Chlorobenzotrifluoride	180	11.274	11.276	-0.002	97	125099	25.0	26.6	
97 Isopropylbenzene	105	11.377	11.380	-0.003	99	368436	25.0	27.5	
99 1,1,2,2-Tetrachloroethane	83	11.675	11.672	0.003	70	83874	25.0	25.3	
100 Bromobenzene	156	11.681	11.684	-0.003	98	80670	25.0	24.8	
101 1,2,3-Trichloropropane	110	11.718	11.720	-0.002	95	24990	25.0	23.3	
102 trans-1,4-Dichloro-2-butene	53	11.730	11.727	0.003	85	21505	25.0	24.2	
103 N-Propylbenzene	120	11.791	11.787	0.004	100	102304	25.0	25.4	
104 2-Chlorotoluene	126	11.876	11.873	0.003	99	84295	25.0	25.0	
105 3-Chlorotoluene	126	11.937	11.933	0.004	98	91182	25.0	24.2	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
106 1,3,5-Trimethylbenzene	105	11.961	11.964	-0.003	100	289696	25.0	25.8	
107 4-Chlorotoluene	126	11.986	11.982	0.004	96	89370	25.0	24.4	
108 tert-Butylbenzene	119	12.290	12.286	0.004	98	248042	25.0	25.5	
110 1,2,4-Trimethylbenzene	105	12.338	12.335	0.003	99	292909	25.0	25.5	
111 1,2-dichloro-4-(trifluoromethyl)	214	12.399	12.402	-0.003	98	87627	25.0	24.1	
112 sec-Butylbenzene	105	12.509	12.505	0.004	100	361915	25.0	26.5	
113 1,3-Dichlorobenzene	146	12.618	12.615	0.003	99	151937	25.0	25.3	
114 4-Isopropyltoluene	119	12.655	12.651	0.004	99	284792	25.0	25.2	
115 1,4-Dichlorobenzene	146	12.709	12.706	0.003	98	150942	25.0	24.6	
116 2,4-Dichloro-1-(trifluoromethyl)	214	12.758	12.761	-0.003	94	81937	25.0	24.1	
118 2,5-Dichlorobenzotrifluoride	214	12.807	12.809	-0.002	97	94772	25.0	24.9	
120 n-Butylbenzene	91	13.062	13.059	0.003	100	262455	25.0	25.5	
121 1,2-Dichlorobenzene	146	13.081	13.083	-0.002	99	136843	25.0	24.6	
122 1,2-Dibromo-3-Chloropropan	75	13.859	13.862	-0.003	92	9396	25.0	20.7	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125	14.011	14.008	0.003	99	300911	75.0	71.6	
125 2,3- & 3,4- Dichlorotoluene	125	14.425	14.428	-0.003	99	191256	50.0	46.8	
126 1,2,4-Trichlorobenzene	180	14.693	14.689	0.004	98	62363	25.0	21.5	
127 Hexachlorobutadiene	225	14.863	14.860	0.003	95	33676	25.0	24.3	
128 Naphthalene	128	14.942	14.939	0.003	99	161398	25.0	21.2	
129 1,2,3-Trichlorobenzene	180	15.185	15.188	-0.003	97	47333	25.0	19.9	
131 2,4,5-Trichlorotoluene	159	15.964	15.961	0.003	95	24209	25.0	19.0	
130 2,3,6-Trichlorotoluene	159	16.061	16.064	-0.003	95	22020	25.0	19.1	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
S 133 Xylenes, Total	106				0		50.0	50.7	
S 134 1,2-Dichloroethene, Total	96				0		50.0	50.8	
S 135 1,3-Dichloropropene, Total	1				0		50.0	41.0	

### QC Flag Legend

#### Processing Flags

ND - Not Detected or Marked ND

#### Review Flags

M - Manually Integrated

### Reagents:

VOAACRPRI_00003	Amount Added: 5.00	Units: uL	
voaWEEpri Res_00003	Amount Added: 1.00	Units: uL	
voaWKetpri Re_00003	Amount Added: 1.00	Units: uL	
VOA8260SURR_00032	Amount Added: 1.00	Units: uL	
VOA8260VOAPRI_00105	Amount Added: 1.00	Units: uL	
VOAVAPRI_00005	Amount Added: 1.00	Units: uL	
VOA8260INT_00030	Amount Added: 2.00	Units: uL	Run Reagent

Report Date: 17-Mar-2015 10:59:21

Chrom Revision: 2.2 15-Jan-2015 13:05:58

TestAmerica Pittsburgh

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150316-6031.b\\50316004.D

Injection Date: 16-Mar-2015 12:41:30

Instrument ID: CHHP5

Lims ID: IC VSTD5

Operator ID: 001562

Client ID:

Worklist Smp#: 4

Purge Vol: 5.000 mL

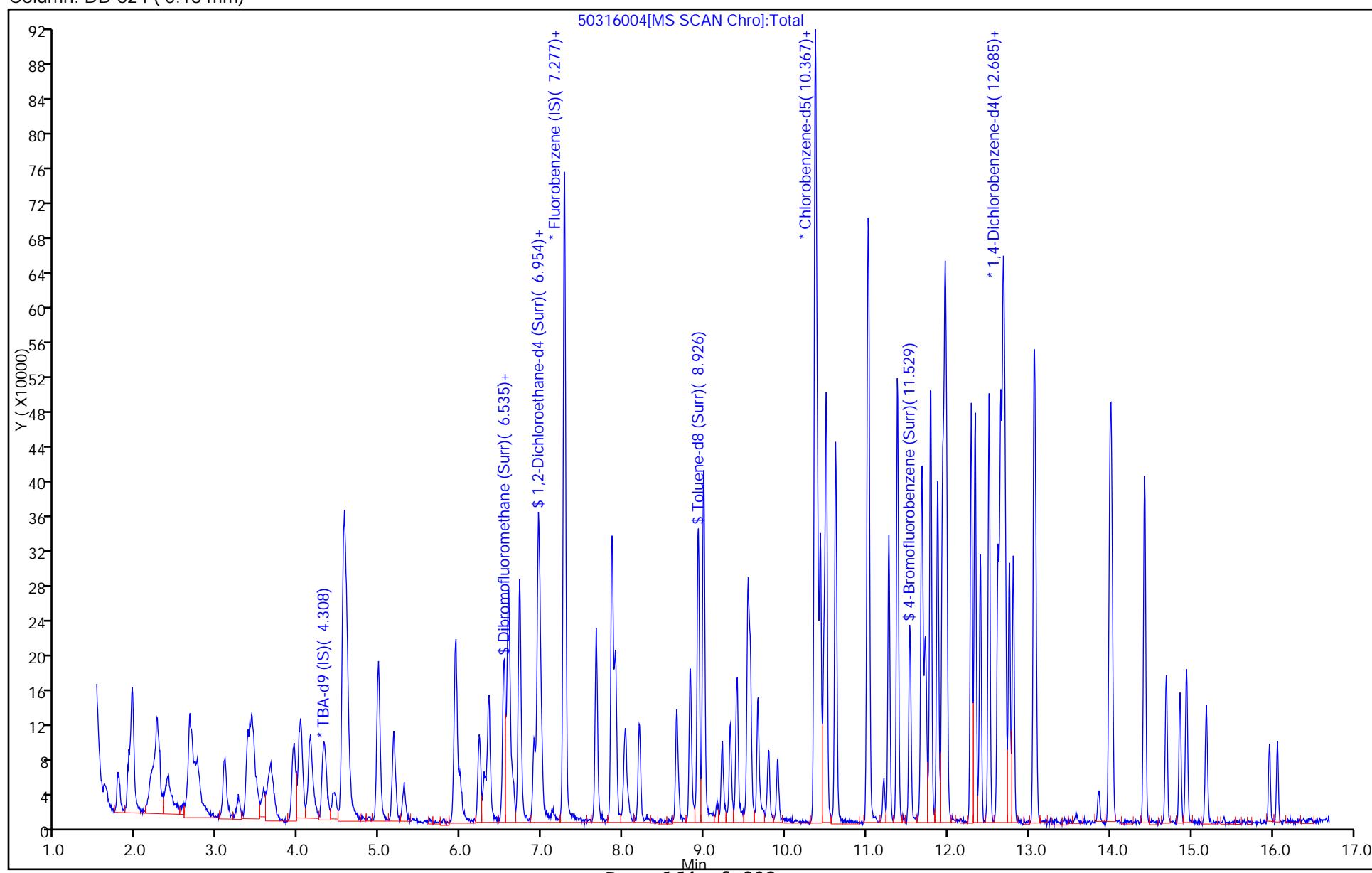
Dil. Factor: 1.0000

ALS Bottle#: 4

Method: MSVOA\_LL\_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 ( 0.18 mm)



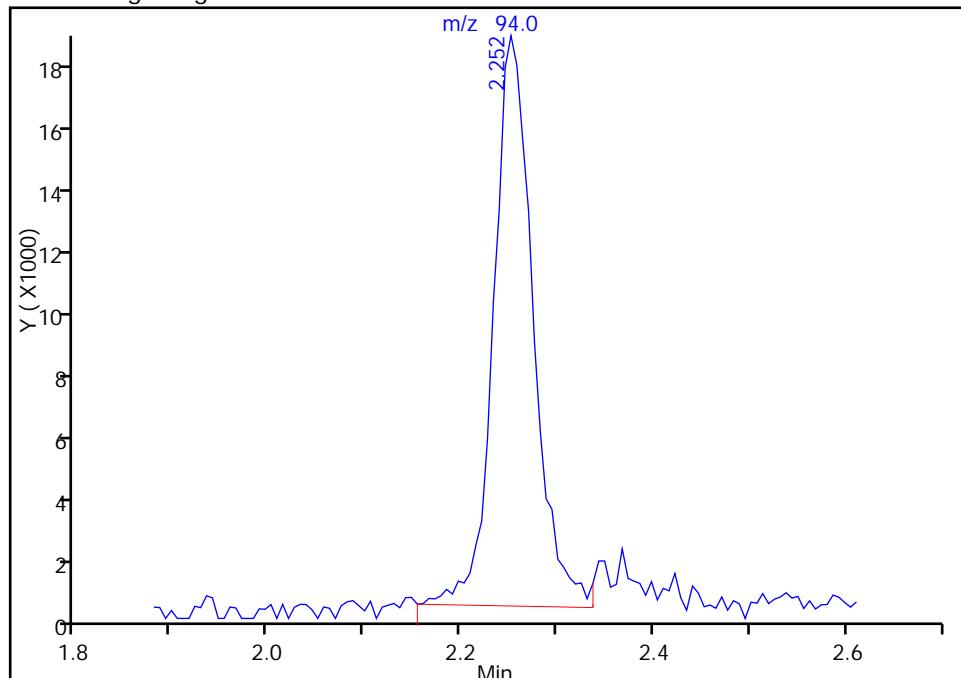
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316004.D  
 Injection Date: 16-Mar-2015 12:41:30 Instrument ID: CHHP5  
 Lims ID: IC VSTD5  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 4 Worklist Smp#: 4  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 15 Bromomethane, CAS: 74-83-9

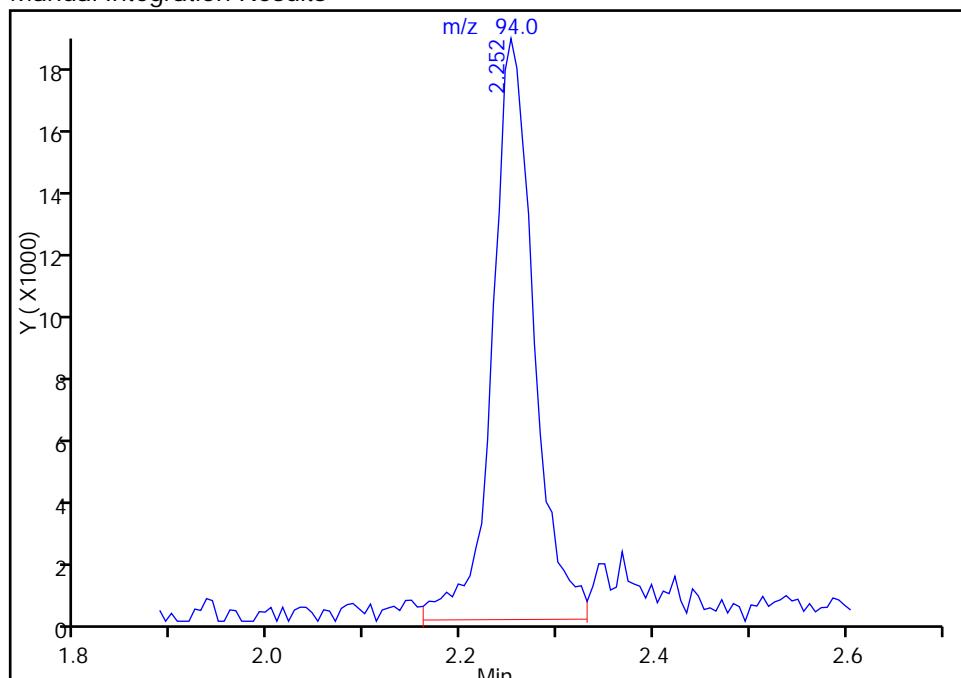
RT: 2.25  
 Area: 51742  
 Amount: 22.147125  
 Amount Units: ng

## Processing Integration Results



RT: 2.25  
 Area: 55097  
 Amount: 26.195176  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 17-Mar-2015 09:42:10

Audit Action: Manually Integrated

Audit Reason: Baseline

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316005.D  
 Lims ID: ICIS VSTD10  
 Client ID:  
 Sample Type: ICIS Calib Level: 3  
 Inject. Date: 16-Mar-2015 13:05:30 ALS Bottle#: 5 Worklist Smp#: 5  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: ICIS VSTD10  
 Misc. Info.: 180-0006031-005  
 Operator ID: 001562 Instrument ID: CHHP5  
 Sublist: chrom-MSVOA\_LL\_CHHP5\*sub4  
 Method: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 17-Mar-2015 10:59:21 Calib Date: 16-Mar-2015 16:17:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICal File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316013.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK012

First Level Reviewer: fergusond Date: 16-Mar-2015 15:03:52

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.305	4.305	0.000	86	135440	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.273	7.273	0.000	99	538139	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.364	10.364	0.000	97	121549	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.682	12.682	0.000	98	174397	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.525	6.525	0.000	95	122918	50.0	50.2	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.902	6.902	0.000	99	162227	50.0	50.3	
\$ 7 Toluene-d8 (Surr)	98	8.922	8.922	0.000	100	527093	50.0	54.4	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.532	11.532	0.000	95	183629	50.0	52.6	
11 Dichlorodifluoromethane	85	1.616	1.616	0.000	100	116111	50.0	50.3	
12 Chloromethane	50	1.774	1.774	0.000	100	159885	50.0	50.2	
13 Vinyl chloride	62	1.902	1.902	0.000	100	183317	50.0	51.5	
14 Butadiene	39	1.944	1.944	0.000	99	208815	50.0	51.4	
15 Bromomethane	94	2.249	2.249	0.000	93	100717	50.0	52.5	
16 Chloroethane	64	2.370	2.370	0.000	98	126349	50.0	51.3	
17 Dichlorofluoromethane	67	2.650	2.650	0.000	100	282324	50.0	50.2	
18 Trichlorofluoromethane	101	2.711	2.711	0.000	98	217544	50.0	51.0	
20 Ethyl ether	59	3.088	3.088	0.000	98	138609	50.0	49.2	
21 Acrolein	56	3.252	3.252	0.000	98	50582	150.0	147.9	
22 1,1-Dichloroethene	96	3.386	3.386	0.000	99	151843	50.0	48.9	
23 1,1,2-Trichloro-1,2,2-trif	101	3.429	3.429	0.000	97	159979	50.0	51.0	
24 Acetone	43	3.496	3.496	0.000	99	102899	100.0	93.3	
25 Iodomethane	142	3.587	3.587	0.000	96	216640	50.0	50.3	
26 Carbon disulfide	76	3.654	3.654	0.000	100	387934	50.0	51.1	
28 3-Chloro-1-propene	76	3.940	3.940	0.000	96	80577	50.0	49.1	
30 Methyl acetate	43	4.019	4.019	0.000	100	641136	250.0	248.6	
31 Methylene Chloride	84	4.134	4.134	0.000	86	168570	50.0	47.0	
32 2-Methyl-2-propanol	59	4.445	4.445	0.000	86	83098	500.0	520.9	
33 Acrylonitrile	53	4.554	4.554	0.000	99	666088	500.0	502.1	
34 trans-1,2-Dichloroethene	96	4.560	4.560	0.000	59	161381	50.0	50.3	
35 Methyl tert-butyl ether	73	4.591	4.591	0.000	96	336961	50.0	47.5	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
36 Hexane	57	4.980	4.980	0.000	98	261916	50.0	51.1	
37 1,1-Dichloroethane	63	5.169	5.169	0.000	100	291408	50.0	50.9	
38 Vinyl acetate	43	5.290	5.290	0.000	100	187915	50.0	46.2	
44 2,2-Dichloropropane	77	5.923	5.923	0.000	67	70106	50.0	48.9	
45 cis-1,2-Dichloroethene	96	5.935	5.935	0.000	92	164893	50.0	48.8	
46 2-Butanone (MEK)	43	5.990	5.990	0.000	100	180996	100.0	102.7	
49 Chlorobromomethane	128	6.233	6.233	0.000	95	71124	50.0	48.6	
51 Tetrahydrofuran	42	6.288	6.288	0.000	98	110274	100.0	99.9	
52 Chloroform	83	6.343	6.343	0.000	96	262371	50.0	50.4	
53 1,1,1-Trichloroethane	97	6.531	6.531	0.000	95	167130	50.0	50.3	
54 Cyclohexane	56	6.586	6.586	0.000	95	322468	50.0	50.5	
56 Carbon tetrachloride	117	6.720	6.720	0.000	69	132517	50.0	49.7	
55 1,1-Dichloropropene	75	6.726	6.726	0.000	97	219974	50.0	51.0	
57 Isobutyl alcohol	41	6.945	6.945	0.000	37	83109	1250.0	1157.0	
58 Benzene	78	6.957	6.957	0.000	99	654151	50.0	51.3	
59 1,2-Dichloroethane	62	6.981	6.981	0.000	97	208683	50.0	50.0	
62 n-Heptane	43	7.280	7.280	0.000	81	222515	50.0	50.8	
64 Trichloroethene	130	7.669	7.669	0.000	98	162608	50.0	50.9	
66 Methylcyclohexane	83	7.864	7.864	0.000	96	295972	50.0	51.9	
67 1,2-Dichloropropane	63	7.906	7.906	0.000	95	154467	50.0	49.0	
68 Dibromomethane	93	8.028	8.028	0.000	95	82469	50.0	48.6	
70 1,4-Dioxane	88	8.058	8.058	0.000	96	31354	1000.0	944.0	M
71 Dichlorobromomethane	83	8.198	8.198	0.000	99	176851	50.0	51.0	
74 cis-1,3-Dichloropropene	75	8.661	8.661	0.000	99	152581	50.0	45.6	
75 4-Methyl-2-pentanone (MIBK)	43	8.825	8.825	0.000	99	342539	100.0	104.1	
76 Toluene	91	8.989	8.989	0.000	100	679332	50.0	54.5	
77 trans-1,3-Dichloropropene	75	9.220	9.220	0.000	94	108942	50.0	48.4	
78 Ethyl methacrylate	69	9.318	9.318	0.000	96	142858	50.0	48.7	
79 1,1,2-Trichloroethane	97	9.397	9.397	0.000	99	125390	50.0	53.7	
80 Tetrachloroethene	164	9.537	9.537	0.000	95	129494	50.0	53.1	
81 1,3-Dichloropropane	76	9.567	9.567	0.000	98	233217	50.0	53.7	
82 2-Hexanone	43	9.658	9.658	0.000	99	271508	100.0	108.0	
84 Chlorodibromomethane	129	9.786	9.786	0.000	99	96762	50.0	51.9	
85 Ethylene Dibromide	107	9.902	9.902	0.000	98	115204	50.0	51.7	
86 3-Chlorobenzotrifluoride	180	10.370	10.370	0.000	97	251080	50.0	52.8	
87 Chlorobenzene	112	10.394	10.394	0.000	99	416488	50.0	52.8	
88 4-Chlorobenzotrifluoride	180	10.431	10.431	0.000	99	234233	50.0	51.0	
89 1,1,1,2-Tetrachloroethane	131	10.473	10.473	0.000	97	101650	50.0	49.9	
90 Ethylbenzene	106	10.498	10.498	0.000	100	242856	50.0	53.6	
91 m-Xylene & p-Xylene	106	10.619	10.619	0.000	99	293796	50.0	53.0	
92 o-Xylene	106	11.015	11.015	0.000	97	285835	50.0	52.8	
93 Styrene	104	11.027	11.027	0.000	99	469890	50.0	53.8	
94 Bromoform	173	11.209	11.209	0.000	96	57667	50.0	50.1	
96 2-Chlorobenzotrifluoride	180	11.276	11.276	0.000	99	252226	50.0	53.1	
97 Isopropylbenzene	105	11.380	11.380	0.000	100	741027	50.0	54.8	
99 1,1,2,2-Tetrachloroethane	83	11.672	11.672	0.000	97	171864	50.0	51.3	
100 Bromobenzene	156	11.684	11.684	0.000	98	168649	50.0	52.2	
101 1,2,3-Trichloropropane	110	11.720	11.720	0.000	97	55900	50.0	52.7	
102 trans-1,4-Dichloro-2-butene	53	11.727	11.727	0.000	88	42827	50.0	48.6	
103 N-Propylbenzene	120	11.787	11.787	0.000	100	210687	50.0	52.9	
104 2-Chlorotoluene	126	11.873	11.873	0.000	100	177793	50.0	53.1	
105 3-Chlorotoluene	126	11.933	11.933	0.000	99	185477	50.0	49.6	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
106 1,3,5-Trimethylbenzene	105	11.964	11.964	0.000	99	588847	50.0	53.0	
107 4-Chlorotoluene	126	11.982	11.982	0.000	99	189449	50.0	52.3	
108 tert-Butylbenzene	119	12.286	12.286	0.000	100	517188	50.0	53.8	
110 1,2,4-Trimethylbenzene	105	12.335	12.335	0.000	100	610150	50.0	53.5	
111 1,2-dichloro-4-(trifluoromethyl)	214	12.402	12.402	0.000	98	179092	50.0	49.7	
112 sec-Butylbenzene	105	12.505	12.505	0.000	100	732318	50.0	54.1	
113 1,3-Dichlorobenzene	146	12.615	12.615	0.000	98	302903	50.0	50.9	
114 4-Isopropyltoluene	119	12.651	12.651	0.000	99	605051	50.0	54.1	
115 1,4-Dichlorobenzene	146	12.706	12.706	0.000	98	310551	50.0	51.1	
116 2,4-Dichloro-1-(trifluoromethyl)	214	12.761	12.761	0.000	94	174468	50.0	51.7	
118 2,5-Dichlorobenzotrifluoride	214	12.809	12.809	0.000	98	186350	50.0	49.4	
120 n-Butylbenzene	91	13.059	13.059	0.000	100	542017	50.0	53.3	
121 1,2-Dichlorobenzene	146	13.083	13.083	0.000	100	283138	50.0	51.4	
122 1,2-Dibromo-3-Chloropropan	75	13.862	13.862	0.000	86	21428	50.0	47.6	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125	14.008	14.008	0.000	99	613057	150.0	147.3	
125 2,3- & 3,4- Dichlorotoluene	125	14.428	14.428	0.000	100	386758	100.0	95.5	
126 1,2,4-Trichlorobenzene	180	14.689	14.689	0.000	96	132179	50.0	46.1	
127 Hexachlorobutadiene	225	14.860	14.860	0.000	96	67414	50.0	49.0	
128 Naphthalene	128	14.939	14.939	0.000	100	347596	50.0	46.2	
129 1,2,3-Trichlorobenzene	180	15.188	15.188	0.000	98	105062	50.0	44.7	
131 2,4,5-Trichlorotoluene	159	15.961	15.961	0.000	96	51080	50.0	40.4	
130 2,3,6-Trichlorotoluene	159	16.064	16.064	0.000	97	47319	50.0	41.5	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
S 133 Xylenes, Total	106				0		100.0	105.8	
S 134 1,2-Dichloroethene, Total	96				0		100.0	99.0	
S 135 1,3-Dichloropropene, Total	1				0		100.0	94.1	

### QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

### Reagents:

VOAVAPRI_00005	Amount Added: 2.00	Units: uL	
VOA8260VOAPRI_00105	Amount Added: 2.00	Units: uL	
VOA8260SURR_00032	Amount Added: 2.00	Units: uL	
voaWEEpri Res_00003	Amount Added: 2.00	Units: uL	
voaWKetpri Re_00003	Amount Added: 2.00	Units: uL	
VOAACRPRI_00003	Amount Added: 6.00	Units: uL	
VOA8260INT_00030	Amount Added: 2.00	Units: uL	Run Reagent

Report Date: 17-Mar-2015 10:59:23

Chrom Revision: 2.2 15-Jan-2015 13:05:58

TestAmerica Pittsburgh

Data File: \PITCHROM\ChromData\CHHP5\20150316-6031.b\50316005.D

Injection Date: 16-Mar-2015 13:05:30

Instrument ID: CHHP5

Operator ID: 001562

Lims ID: ICIS VSTD10

Worklist Smp#: 5

Client ID:

Purge Vol: 5.000 mL

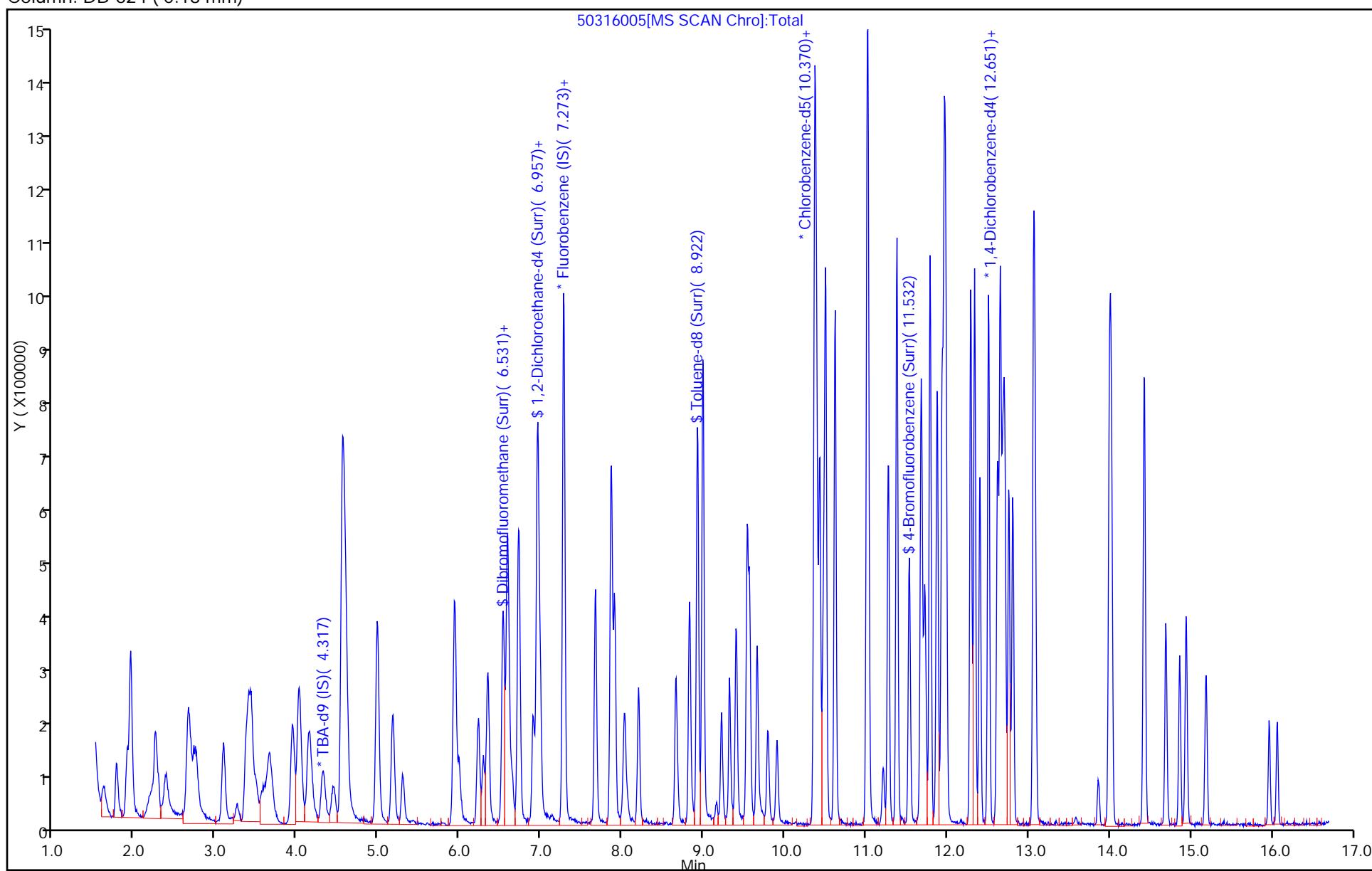
Dil. Factor: 1.0000

ALS Bottle#: 5

Method: MSVOA\_LL\_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 ( 0.18 mm)



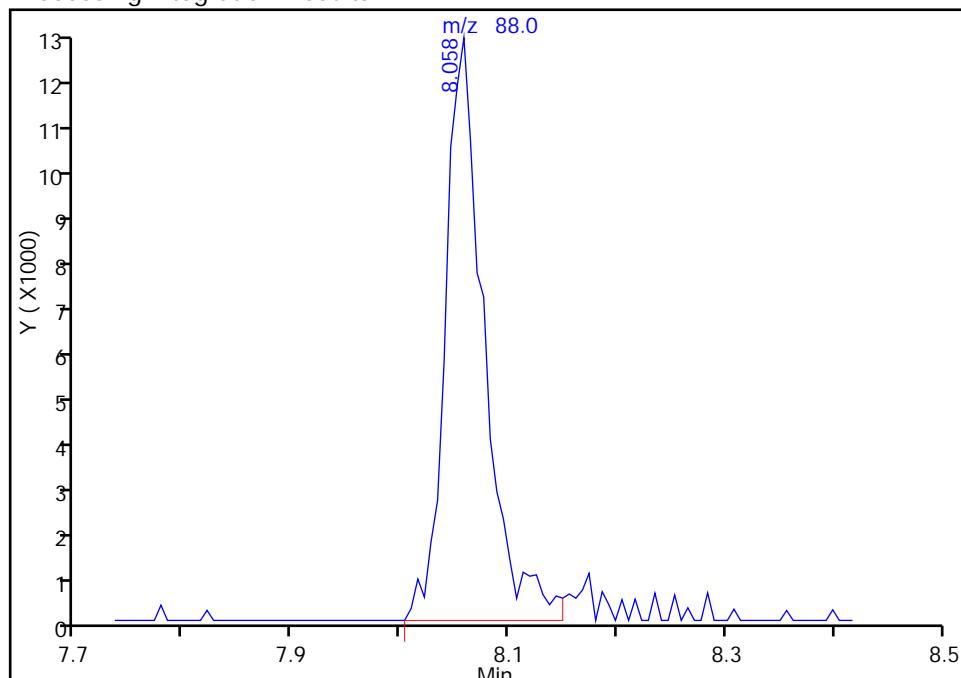
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316005.D  
 Injection Date: 16-Mar-2015 13:05:30 Instrument ID: CHHP5  
 Lims ID: ICIS VSTD10  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 5 Worklist Smp#: 5  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 70 1,4-Dioxane, CAS: 123-91-1

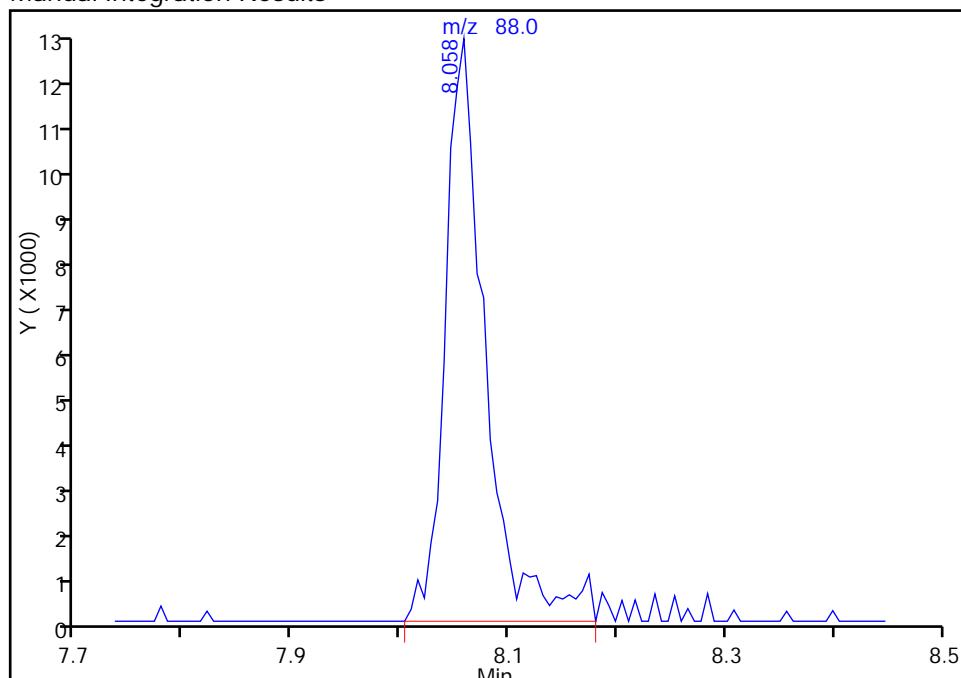
RT: 8.06  
 Area: 30397  
 Amount: 939.9751  
 Amount Units: ng

## Processing Integration Results



RT: 8.06  
 Area: 31354  
 Amount: 944.0403  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 17-Mar-2015 09:27:38

Audit Action: Manually Integrated

Audit Reason: Peak Tail

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316006.D  
 Lims ID: IC VSTD15  
 Client ID:  
 Sample Type: IC Calib Level: 4  
 Inject. Date: 16-Mar-2015 13:29:30 ALS Bottle#: 6 Worklist Smp#: 6  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: IC VSTD15  
 Misc. Info.: 180-0006031-006  
 Operator ID: 001562 Instrument ID: CHHP5  
 Sublist: chrom-MSVOA\_LL\_CHHP5\*sub4  
 Method: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 17-Mar-2015 10:59:26 Calib Date: 16-Mar-2015 16:17:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICal File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316013.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK012

First Level Reviewer: fergusond Date: 17-Mar-2015 09:45:06

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.305	4.305	0.000	89	152705	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.273	7.273	0.000	99	530419	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.364	10.364	0.000	99	125149	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.682	12.682	0.000	95	182887	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.525	6.525	0.000	97	185698	75.0	77.0	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.902	6.902	0.000	96	247858	75.0	77.9	
\$ 7 Toluene-d8 (Surr)	98	8.922	8.922	0.000	100	794092	75.0	79.6	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.532	11.532	0.000	97	279546	75.0	77.8	
11 Dichlorodifluoromethane	85	1.622	1.622	0.000	99	173113	75.0	76.1	
12 Chloromethane	50	1.768	1.768	0.000	99	249772	75.0	79.6	
13 Vinyl chloride	62	1.896	1.896	0.000	100	280135	75.0	79.9	
14 Butadiene	39	1.944	1.944	0.000	99	317272	75.0	79.2	
15 Bromomethane	94	2.249	2.249	0.000	92	159846	75.0	87.5	
16 Chloroethane	64	2.376	2.376	0.000	96	191164	75.0	78.8	
17 Dichlorofluoromethane	67	2.644	2.644	0.000	99	437737	75.0	79.0	
18 Trichlorofluoromethane	101	2.723	2.723	0.000	96	358375	75.0	85.2	
20 Ethyl ether	59	3.082	3.082	0.000	100	214135	75.0	77.2	
21 Acrolein	56	3.258	3.258	0.000	100	62132	175.0	184.3	
22 1,1-Dichloroethene	96	3.374	3.374	0.000	100	235889	75.0	77.1	
23 1,1,2-Trichloro-1,2,2-trif	101	3.423	3.423	0.000	100	246660	75.0	79.7	
24 Acetone	43	3.496	3.496	0.000	100	180387	150.0	166.0	
25 Iodomethane	142	3.581	3.581	0.000	100	334141	75.0	78.6	
26 Carbon disulfide	76	3.660	3.660	0.000	100	592248	75.0	79.2	
28 3-Chloro-1-propene	76	3.934	3.934	0.000	100	125423	75.0	77.6	
30 Methyl acetate	43	4.019	4.019	0.000	100	994505	375.0	391.2	
31 Methylene Chloride	84	4.147	4.147	0.000	100	256424	75.0	72.5	
32 2-Methyl-2-propanol	59	4.439	4.439	0.000	100	133756	750.0	743.6	
33 Acrylonitrile	53	4.554	4.554	0.000	100	1035956	750.0	792.2	
34 trans-1,2-Dichloroethene	96	4.560	4.560	0.000	100	251288	75.0	79.4	
35 Methyl tert-butyl ether	73	4.597	4.597	0.000	100	528520	75.0	75.6	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
36 Hexane	57	4.980	4.980	0.000	100	392065	75.0	77.6	
37 1,1-Dichloroethane	63	5.175	5.175	0.000	100	435915	75.0	77.2	
38 Vinyl acetate	43	5.296	5.296	0.000	100	294456	75.0	73.5	
44 2,2-Dichloropropane	77	5.929	5.929	0.000	100	108858	75.0	77.1	
45 cis-1,2-Dichloroethene	96	5.941	5.941	0.000	100	259517	75.0	77.9	
46 2-Butanone (MEK)	43	5.990	5.990	0.000	100	259227	150.0	149.2	
49 Chlorobromomethane	128	6.227	6.227	0.000	100	109930	75.0	76.2	
51 Tetrahydrofuran	42	6.282	6.282	0.000	100	166594	150.0	153.1	
52 Chloroform	83	6.343	6.343	0.000	100	395935	75.0	77.2	
53 1,1,1-Trichloroethane	97	6.531	6.531	0.000	100	259963	75.0	79.4	
54 Cyclohexane	56	6.586	6.586	0.000	100	497889	75.0	79.2	
56 Carbon tetrachloride	117	6.720	6.720	0.000	100	203736	75.0	77.5	
55 1,1-Dichloropropene	75	6.726	6.726	0.000	100	326699	75.0	76.8	
57 Isobutyl alcohol	41	6.945	6.945	0.000	100	137203	1875.0	1937.9	M
58 Benzene	78	6.957	6.957	0.000	100	984614	75.0	78.3	
59 1,2-Dichloroethane	62	6.988	6.988	0.000	100	320594	75.0	77.9	
62 n-Heptane	43	7.280	7.280	0.000	100	335961	75.0	77.8	
64 Trichloroethene	130	7.669	7.669	0.000	100	242252	75.0	76.9	
66 Methylcyclohexane	83	7.864	7.864	0.000	100	446628	75.0	79.5	
67 1,2-Dichloropropane	63	7.906	7.906	0.000	100	238331	75.0	76.7	
68 Dibromomethane	93	8.022	8.022	0.000	100	130496	75.0	78.0	
70 1,4-Dioxane	88	8.058	8.058	0.000	100	50907	1500.0	1555.1	
71 Dichlorobromomethane	83	8.198	8.198	0.000	100	259871	75.0	76.1	
74 cis-1,3-Dichloropropene	75	8.661	8.661	0.000	100	247138	75.0	75.0	
75 4-Methyl-2-pentanone (MIBK)	43	8.825	8.825	0.000	100	531084	150.0	156.8	
76 Toluene	91	8.989	8.989	0.000	100	1017198	75.0	79.3	
77 trans-1,3-Dichloropropene	75	9.220	9.220	0.000	100	167274	75.0	72.2	
78 Ethyl methacrylate	69	9.318	9.318	0.000	100	221852	75.0	73.4	
79 1,1,2-Trichloroethane	97	9.403	9.403	0.000	100	183907	75.0	76.5	
80 Tetrachloroethene	164	9.537	9.537	0.000	100	194422	75.0	77.5	
81 1,3-Dichloropropane	76	9.561	9.561	0.000	100	342719	75.0	76.7	
82 2-Hexanone	43	9.658	9.658	0.000	100	402386	150.0	155.5	
84 Chlorodibromomethane	129	9.792	9.792	0.000	100	145315	75.0	75.7	
85 Ethylene Dibromide	107	9.902	9.902	0.000	100	179814	75.0	78.4	
86 3-Chlorobenzotrifluoride	180	10.370	10.370	0.000	100	388132	75.0	79.3	
87 Chlorobenzene	112	10.388	10.388	0.000	100	622968	75.0	76.7	
88 4-Chlorobenzotrifluoride	180	10.431	10.431	0.000	100	368570	75.0	77.9	
89 1,1,1,2-Tetrachloroethane	131	10.473	10.473	0.000	100	159225	75.0	75.9	
90 Ethylbenzene	106	10.504	10.504	0.000	100	366398	75.0	78.6	
91 m-Xylene & p-Xylene	106	10.619	10.619	0.000	100	454933	75.0	79.8	
92 o-Xylene	106	11.009	11.009	0.000	100	436586	75.0	78.3	
93 Styrene	104	11.027	11.027	0.000	100	712222	75.0	79.2	
94 Bromoform	173	11.209	11.209	0.000	100	90522	75.0	76.3	
96 2-Chlorobenzotrifluoride	180	11.276	11.276	0.000	100	386985	75.0	79.2	
97 Isopropylbenzene	105	11.380	11.380	0.000	100	1137215	75.0	81.7	
99 1,1,2,2-Tetrachloroethane	83	11.678	11.678	0.000	100	264462	75.0	76.7	
100 Bromobenzene	156	11.678	11.678	0.000	100	253502	75.0	74.9	
101 1,2,3-Trichloropropane	110	11.721	11.721	0.000	100	81225	75.0	73.0	
102 trans-1,4-Dichloro-2-butene	53	11.733	11.733	0.000	100	66879	75.0	72.3	
103 N-Propylbenzene	120	11.787	11.787	0.000	100	316980	75.0	75.9	
104 2-Chlorotoluene	126	11.873	11.873	0.000	100	262207	75.0	74.7	
105 3-Chlorotoluene	126	11.933	11.933	0.000	100	291288	75.0	74.3	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
106 1,3,5-Trimethylbenzene	105	11.964	11.964	0.000	100	921783	75.0	79.2	
107 4-Chlorotoluene	126	11.982	11.982	0.000	100	296950	75.0	78.2	
108 tert-Butylbenzene	119	12.286	12.286	0.000	100	794422	75.0	78.8	
110 1,2,4-Trimethylbenzene	105	12.335	12.335	0.000	100	951216	75.0	79.6	
111 1,2-dichloro-4-(trifluoromethyl)	214	12.402	12.402	0.000	100	303120	75.0	80.3	
112 sec-Butylbenzene	105	12.511	12.511	0.000	100	1138120	75.0	80.2	
113 1,3-Dichlorobenzene	146	12.621	12.621	0.000	100	480001	75.0	77.0	
114 4-Isopropyltoluene	119	12.651	12.651	0.000	100	948139	75.0	80.9	
115 1,4-Dichlorobenzene	146	12.706	12.706	0.000	100	484138	75.0	76.0	
116 2,4-Dichloro-1-(trifluoromethyl)	214	12.755	12.755	0.000	100	289446	75.0	81.8	
118 2,5-Dichlorobenzotrifluoride	214	12.803	12.803	0.000	100	309155	75.0	78.1	
120 n-Butylbenzene	91	13.065	13.065	0.000	100	861784	75.0	80.7	
121 1,2-Dichlorobenzene	146	13.083	13.083	0.000	100	440732	75.0	76.3	
122 1,2-Dibromo-3-Chloropropan	75	13.856	13.856	0.000	100	36318	75.0	76.9	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125	14.008	14.008	0.000	100	1058653	225.0	242.5	
125 2,3- & 3,4- Dichlorotoluene	125	14.428	14.428	0.000	100	694253	150.0	163.5	
126 1,2,4-Trichlorobenzene	180	14.695	14.695	0.000	100	240861	75.0	80.1	
127 Hexachlorobutadiene	225	14.866	14.866	0.000	100	112236	75.0	77.9	
128 Naphthalene	128	14.939	14.939	0.000	100	657935	75.0	83.4	
129 1,2,3-Trichlorobenzene	180	15.188	15.188	0.000	100	200345	75.0	81.3	
131 2,4,5-Trichlorotoluene	159	15.967	15.967	0.000	100	108037	75.0	81.5	
130 2,3,6-Trichlorotoluene	159	16.064	16.064	0.000	100	98974	75.0	82.7	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
S 133 Xylenes, Total	106				0		150.0	158.1	
S 134 1,2-Dichloroethene, Total	96				0		150.0	157.3	
S 135 1,3-Dichloropropene, Total	1				0		150.0	147.2	

### QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

### Reagents:

VOAACRPRI_00003	Amount Added: 7.00	Units: uL	
voaWEEpri Res_00003	Amount Added: 3.00	Units: uL	
voaWKetpri Re_00003	Amount Added: 3.00	Units: uL	
VOA8260SURR_00032	Amount Added: 3.00	Units: uL	
VOA8260VOAPRI_00105	Amount Added: 3.00	Units: uL	
VOAVAPRI_00005	Amount Added: 3.00	Units: uL	
VOA8260INT_00030	Amount Added: 2.00	Units: uL	Run Reagent

Report Date: 17-Mar-2015 10:59:27

Chrom Revision: 2.2 15-Jan-2015 13:05:58

TestAmerica Pittsburgh

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150316-6031.b\\50316006.D

Injection Date: 16-Mar-2015 13:29:30

Instrument ID: CHHP5

Operator ID: 001562

Lims ID: IC VSTD15

Worklist Smp#: 6

Client ID:

Purge Vol: 5.000 mL

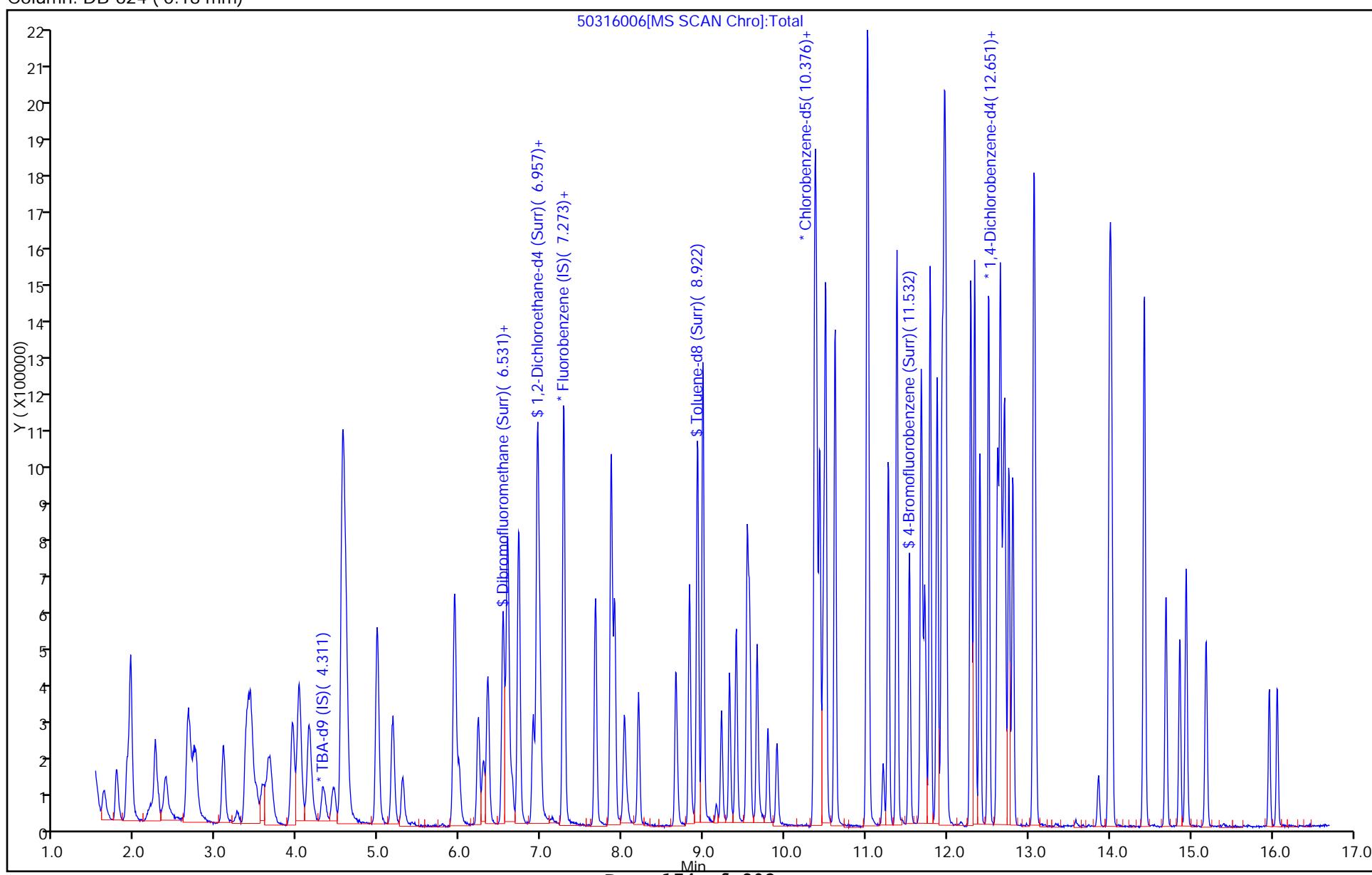
Dil. Factor: 1.0000

ALS Bottle#: 6

Method: MSVOA\_LL\_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 ( 0.18 mm)



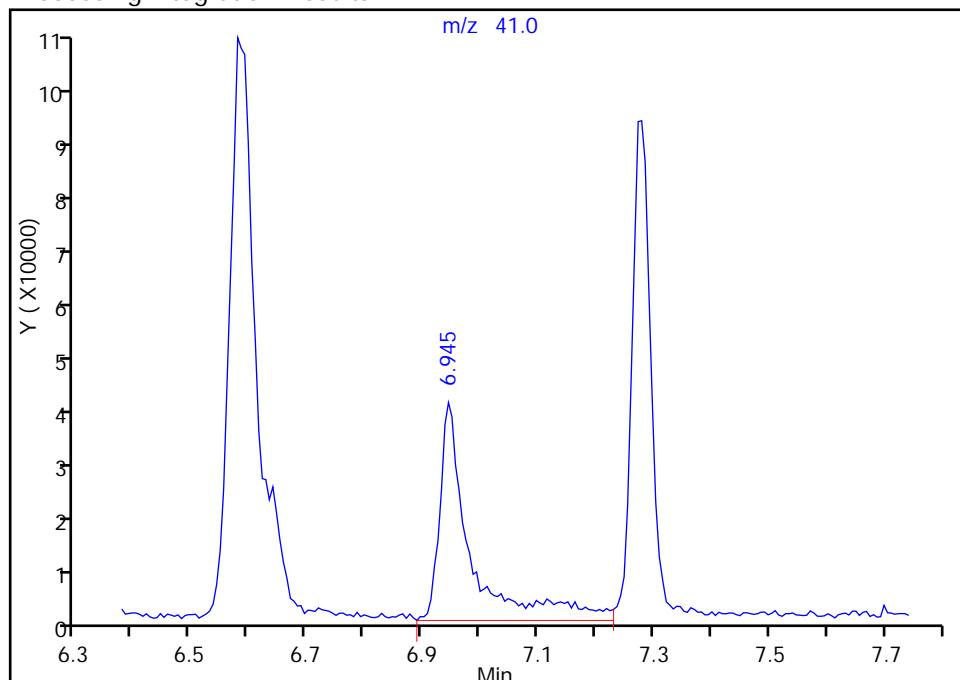
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316006.D  
 Injection Date: 16-Mar-2015 13:29:30 Instrument ID: CHHP5  
 Lims ID: IC VSTD15  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 6 Worklist Smp#: 6  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 57 Isobutyl alcohol, CAS: 78-83-1

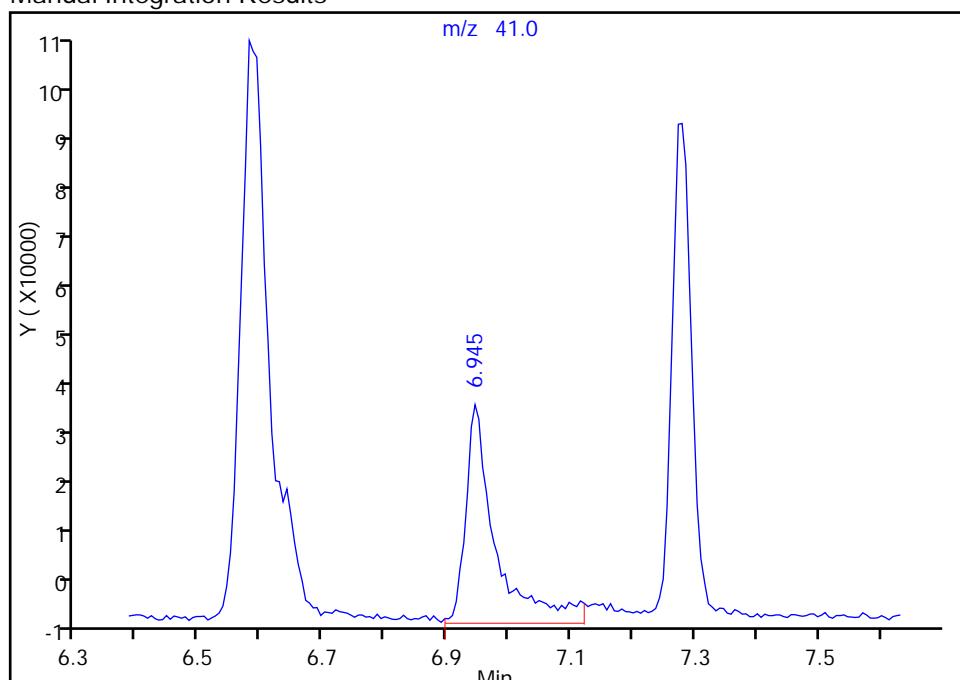
RT: 6.94  
 Area: 150922  
 Amount: 2067.3126  
 Amount Units: ng

## Processing Integration Results



RT: 6.94  
 Area: 137203  
 Amount: 1937.8985  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 17-Mar-2015 09:45:06

Audit Action: Manually Integrated

Audit Reason: Peak Tail

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316007.D  
 Lims ID: IC VSTD20  
 Client ID:  
 Sample Type: IC Calib Level: 5  
 Inject. Date: 16-Mar-2015 13:53:30 ALS Bottle#: 7 Worklist Smp#: 7  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: IC VSTD20  
 Misc. Info.: 180-0006031-007  
 Operator ID: 001562 Instrument ID: CHHP5  
 Sublist: chrom-MSVOA\_LL\_CHHP5\*sub4  
 Method: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 17-Mar-2015 10:59:28 Calib Date: 16-Mar-2015 16:17:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICal File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316013.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK012

First Level Reviewer: fergusond Date: 17-Mar-2015 09:48:25

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.326	4.305	0.021	86	154462	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.270	7.273	-0.003	99	558174	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.367	10.364	0.003	99	128898	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.685	12.682	0.003	99	188542	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.528	6.525	0.003	99	248750	100.0	98.0	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.899	6.902	-0.003	97	335757	100.0	100.3	
\$ 7 Toluene-d8 (Surr)	98	8.919	8.922	-0.003	100	1053927	100.0	102.6	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.529	11.532	-0.003	98	379740	100.0	102.6	
11 Dichlorodifluoromethane	85	1.619	1.622	-0.003	98	243823	100.0	101.9	
12 Chloromethane	50	1.777	1.768	0.009	100	316915	100.0	96.0	
13 Vinyl chloride	62	1.905	1.896	0.009	100	370271	100.0	100.3	
14 Butadiene	39	1.947	1.944	0.003	100	415323	100.0	98.5	
15 Bromomethane	94	2.251	2.249	0.002	99	192846	100.0	101.0	
16 Chloroethane	64	2.373	2.376	-0.003	99	245673	100.0	96.2	
17 Dichlorofluoromethane	67	2.653	2.644	0.009	100	548270	100.0	94.0	
18 Trichlorofluoromethane	101	2.702	2.723	-0.021	98	437688	100.0	98.9	
20 Ethyl ether	59	3.085	3.082	0.003	100	293889	100.0	100.7	
21 Acrolein	56	3.261	3.258	0.003	99	71073	200.0	200.4	
22 1,1-Dichloroethene	96	3.377	3.374	0.003	98	318457	100.0	98.9	
23 1,1,2-Trichloro-1,2,2-trif	101	3.432	3.423	0.009	98	319162	100.0	98.0	
24 Acetone	43	3.492	3.496	-0.004	100	217095	200.0	189.9	
25 Iodomethane	142	3.596	3.581	0.015	99	439512	100.0	98.3	
26 Carbon disulfide	76	3.669	3.660	0.009	100	772081	100.0	98.1	
28 3-Chloro-1-propene	76	3.930	3.934	-0.004	99	163875	100.0	96.3	
30 Methyl acetate	43	4.022	4.019	0.003	100	1321970	500.0	494.2	
31 Methylene Chloride	84	4.143	4.147	-0.004	98	345226	100.0	92.7	
32 2-Methyl-2-propanol	59	4.435	4.439	-0.004	99	175500	1000.0	964.6	
33 Acrylonitrile	53	4.551	4.554	-0.003	100	1363975	1000.0	991.2	
34 trans-1,2-Dichloroethene	96	4.557	4.560	-0.003	95	327278	100.0	98.3	
35 Methyl tert-butyl ether	73	4.594	4.597	-0.003	99	727030	100.0	98.8	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
36 Hexane	57	4.983	4.980	0.003	99	514868	100.0	96.8	
37 1,1-Dichloroethane	63	5.171	5.175	-0.004	100	595324	100.0	100.2	
38 Vinyl acetate	43	5.293	5.296	-0.003	100	419086	100.0	99.4	
44 2,2-Dichloropropane	77	5.926	5.929	-0.003	98	147216	100.0	99.1	
45 cis-1,2-Dichloroethene	96	5.938	5.941	-0.003	98	349805	100.0	99.7	
46 2-Butanone (MEK)	43	5.987	5.990	-0.003	100	371447	200.0	203.2	
49 Chlorobromomethane	128	6.230	6.227	0.003	98	150204	100.0	99.0	
51 Tetrahydrofuran	42	6.285	6.282	0.003	99	224920	200.0	196.4	
52 Chloroform	83	6.346	6.343	0.003	100	534362	100.0	99.0	
53 1,1,1-Trichloroethane	97	6.528	6.531	-0.003	99	344772	100.0	100.0	
54 Cyclohexane	56	6.589	6.586	0.003	99	649387	100.0	98.1	
56 Carbon tetrachloride	117	6.717	6.720	-0.003	98	274328	100.0	99.2	
55 1,1-Dichloropropene	75	6.723	6.726	-0.003	98	436454	100.0	97.5	
57 Isobutyl alcohol	41	6.942	6.945	-0.003	98	174166	2500.0	2337.7	M
58 Benzene	78	6.954	6.957	-0.003	99	1312435	100.0	99.2	
59 1,2-Dichloroethane	62	6.984	6.988	-0.004	99	429724	100.0	99.2	
62 n-Heptane	43	7.276	7.280	-0.004	99	443357	100.0	97.6	
64 Trichloroethene	130	7.666	7.669	-0.003	99	326599	100.0	98.5	
66 Methylcyclohexane	83	7.860	7.864	-0.004	100	583894	100.0	98.7	
67 1,2-Dichloropropane	63	7.903	7.906	-0.003	99	332279	100.0	101.6	
68 Dibromomethane	93	8.025	8.022	0.003	99	178905	100.0	101.6	
70 1,4-Dioxane	88	8.055	8.058	-0.003	99	66490	2000.0	1930.1	
71 Dichlorobromomethane	83	8.195	8.198	-0.003	98	363842	100.0	101.2	
74 cis-1,3-Dichloropropene	75	8.657	8.661	-0.004	99	345528	100.0	99.6	
75 4-Methyl-2-pentanone (MIBK)	43	8.822	8.825	-0.003	100	747218	200.0	214.2	
76 Toluene	91	8.992	8.989	0.003	100	1340817	100.0	101.5	
77 trans-1,3-Dichloropropene	75	9.217	9.220	-0.003	99	244258	100.0	102.4	
78 Ethyl methacrylate	69	9.314	9.318	-0.004	97	334858	100.0	107.6	
79 1,1,2-Trichloroethane	97	9.399	9.403	-0.004	99	252461	100.0	101.9	
80 Tetrachloroethene	164	9.539	9.537	0.002	99	261148	100.0	101.1	
81 1,3-Dichloropropane	76	9.564	9.561	0.003	100	467174	100.0	101.5	
82 2-Hexanone	43	9.655	9.658	-0.003	100	541680	200.0	203.2	
84 Chlorodibromomethane	129	9.789	9.792	-0.003	99	210013	100.0	106.2	
85 Ethylene Dibromide	107	9.898	9.902	-0.004	100	245946	100.0	104.0	
86 3-Chlorobenzotrifluoride	180	10.373	10.370	0.003	97	511845	100.0	101.6	
87 Chlorobenzene	112	10.391	10.388	0.003	100	845046	100.0	101.0	
88 4-Chlorobenzotrifluoride	180	10.428	10.431	-0.003	99	511237	100.0	104.9	
89 1,1,1,2-Tetrachloroethane	131	10.476	10.473	0.003	95	233228	100.0	107.9	
90 Ethylbenzene	106	10.501	10.504	-0.003	100	488611	100.0	101.8	
91 m-Xylene & p-Xylene	106	10.616	10.619	-0.003	100	592135	100.0	100.8	
92 o-Xylene	106	11.012	11.009	0.003	100	585609	100.0	101.9	
93 Styrene	104	11.024	11.027	-0.003	95	966850	100.0	104.4	
94 Bromoform	173	11.212	11.209	0.003	98	126605	100.0	103.7	
96 2-Chlorobenzotrifluoride	180	11.273	11.276	-0.003	99	521379	100.0	103.6	
97 Isopropylbenzene	105	11.377	11.380	-0.003	100	1474178	100.0	102.8	
99 1,1,2,2-Tetrachloroethane	83	11.675	11.678	-0.003	98	351798	100.0	99.0	
100 Bromobenzene	156	11.681	11.678	0.003	99	346996	100.0	99.4	
101 1,2,3-Trichloropropane	110	11.717	11.721	-0.003	97	111668	100.0	97.4	
102 trans-1,4-Dichloro-2-butene	53	11.729	11.733	-0.004	97	92761	100.0	97.3	
103 N-Propylbenzene	120	11.784	11.787	-0.003	100	419888	100.0	97.5	
104 2-Chlorotoluene	126	11.875	11.873	0.002	100	351403	100.0	97.2	
105 3-Chlorotoluene	126	11.936	11.933	0.003	98	415463	100.0	102.8	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
106 1,3,5-Trimethylbenzene	105	11.961	11.964	-0.003	100	1188743	100.0	99.0	
107 4-Chlorotoluene	126	11.985	11.982	0.003	97	377870	100.0	96.5	
108 tert-Butylbenzene	119	12.289	12.286	0.003	99	1020106	100.0	98.1	
110 1,2,4-Trimethylbenzene	105	12.338	12.335	0.003	100	1214438	100.0	98.6	
111 1,2-dichloro-4-(trifluoromethyl)	214	12.399	12.402	-0.003	100	396211	100.0	101.8	
112 sec-Butylbenzene	105	12.508	12.511	-0.003	100	1462842	100.0	99.9	
113 1,3-Dichlorobenzene	146	12.618	12.621	-0.003	99	630675	100.0	98.1	
114 4-Isopropyltoluene	119	12.648	12.651	-0.003	100	1195021	100.0	98.9	
115 1,4-Dichlorobenzene	146	12.709	12.706	0.003	99	642365	100.0	97.8	
116 2,4-Dichloro-1-(trifluoromethyl)	214	12.758	12.755	0.003	98	358539	100.0	98.3	
118 2,5-Dichlorobenzotrifluoride	214	12.806	12.803	0.003	99	406971	100.0	99.8	
120 n-Butylbenzene	91	13.062	13.065	-0.003	100	1093564	100.0	99.4	
121 1,2-Dichlorobenzene	146	13.080	13.083	-0.003	99	595901	100.0	100.1	
122 1,2-Dibromo-3-Chloropropan	75	13.859	13.856	0.003	95	47067	100.0	96.6	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125	14.005	14.008	-0.003	100	1333690	300.0	296.3	
125 2,3- & 3,4- Dichlorotoluene	125	14.424	14.428	-0.004	100	866884	200.0	198.1	
126 1,2,4-Trichlorobenzene	180	14.692	14.695	-0.003	99	295444	100.0	95.3	
127 Hexachlorobutadiene	225	14.862	14.866	-0.004	98	140410	100.0	94.5	
128 Naphthalene	128	14.942	14.939	0.003	100	789643	100.0	97.0	
129 1,2,3-Trichlorobenzene	180	15.185	15.188	-0.003	98	242534	100.0	95.4	
131 2,4,5-Trichlorotoluene	159	15.964	15.967	-0.003	98	123791	100.0	90.6	
130 2,3,6-Trichlorotoluene	159	16.061	16.064	-0.003	98	110702	100.0	89.7	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
S 133 Xylenes, Total	106				0		200.0	202.8	
S 134 1,2-Dichloroethene, Total	96				0		200.0	198.0	
S 135 1,3-Dichloropropene, Total	1				0		200.0	202.0	

### QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

### Reagents:

VOAVAPRI_00005	Amount Added: 4.00	Units: uL	
VOA8260VOAPRI_00105	Amount Added: 4.00	Units: uL	
VOA8260SURR_00032	Amount Added: 4.00	Units: uL	
voaWEEpri Res_00003	Amount Added: 4.00	Units: uL	
voaWKetpri Re_00003	Amount Added: 4.00	Units: uL	
VOAACRPRI_00003	Amount Added: 8.00	Units: uL	
VOA8260INT_00030	Amount Added: 2.00	Units: uL	Run Reagent

Report Date: 17-Mar-2015 10:59:29

Chrom Revision: 2.2 15-Jan-2015 13:05:58

TestAmerica Pittsburgh

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150316-6031.b\\50316007.D

Injection Date: 16-Mar-2015 13:53:30

Instrument ID: CHHP5

Operator ID: 001562

Lims ID: IC VSTD20

Worklist Smp#: 7

Client ID:

Purge Vol: 5.000 mL

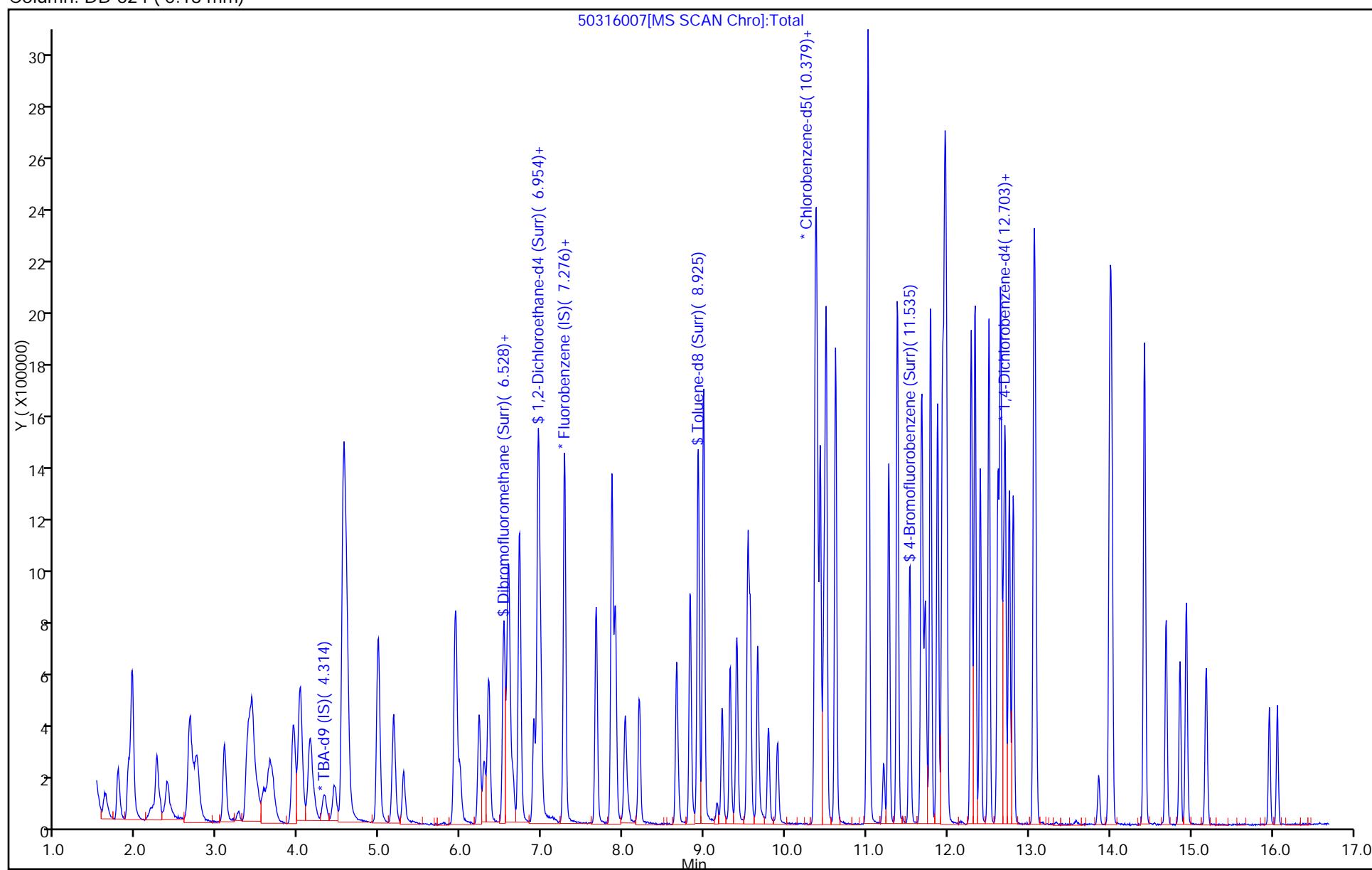
Dil. Factor: 1.0000

ALS Bottle#: 7

Method: MSVOA\_LL\_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 ( 0.18 mm)



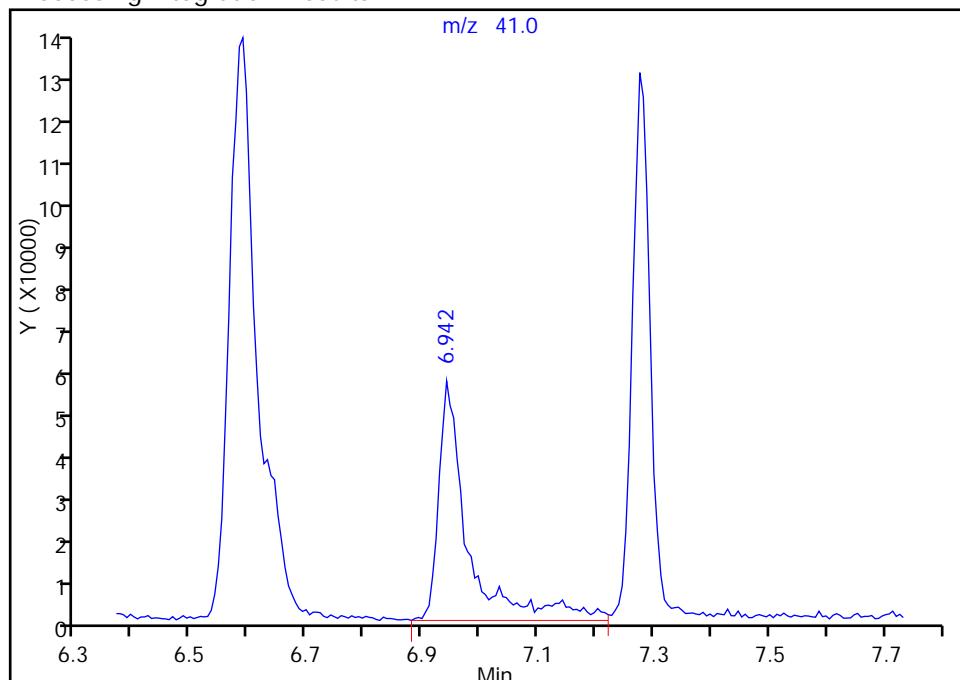
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316007.D  
 Injection Date: 16-Mar-2015 13:53:30 Instrument ID: CHHP5  
 Lims ID: IC VSTD20  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 7 Worklist Smp#: 7  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 57 Isobutyl alcohol, CAS: 78-83-1

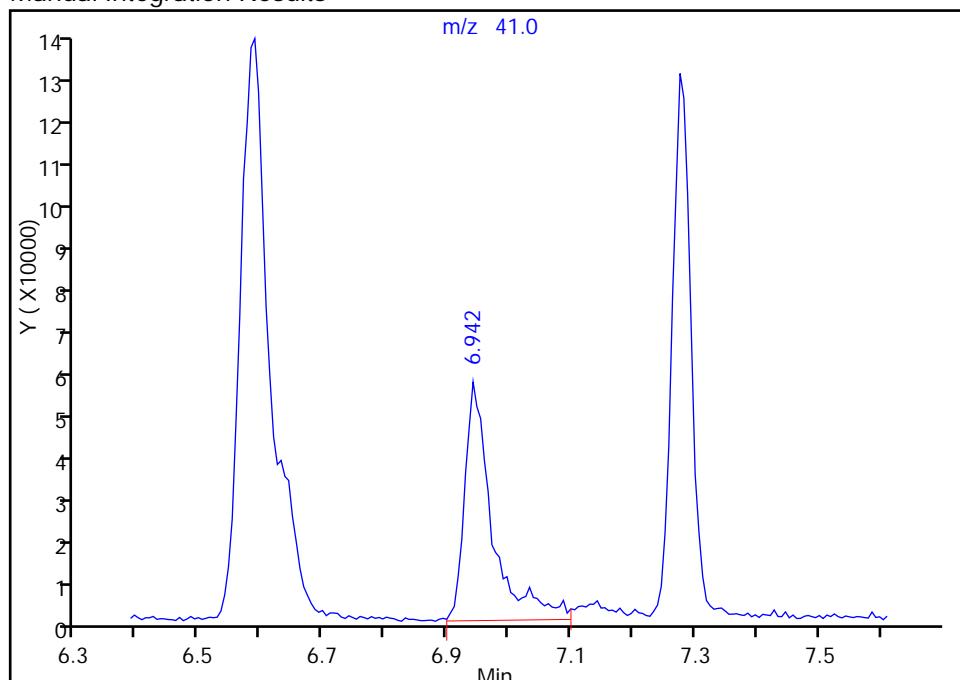
RT: 6.94  
 Area: 197796  
 Amount: 2559.7908  
 Amount Units: ng

## Processing Integration Results



RT: 6.94  
 Area: 174166  
 Amount: 2337.6542  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 17-Mar-2015 09:48:25

Audit Action: Manually Integrated

Audit Reason: Peak Tail

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316008.D  
 Lims ID: IC VSTD35  
 Client ID:  
 Sample Type: IC Calib Level: 6  
 Inject. Date: 16-Mar-2015 14:17:30 ALS Bottle#: 8 Worklist Smp#: 8  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: IC VSTD35  
 Misc. Info.: 180-0006031-008  
 Operator ID: 001562 Instrument ID: CHHP5  
 Sublist: chrom-MSVOA\_LL\_CHHP5\*sub4  
 Method: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 17-Mar-2015 10:59:29 Calib Date: 16-Mar-2015 16:17:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICal File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316013.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK012

First Level Reviewer: fergusond Date: 17-Mar-2015 09:49:39

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.309	4.305	0.004	95	172412	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.271	7.273	-0.002	99	562344	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.362	10.364	-0.002	92	147916	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.686	12.682	0.004	95	201448	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.529	6.525	0.004	99	435320	175.0	170.2	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.900	6.902	-0.002	98	589491	175.0	174.8	
\$ 7 Toluene-d8 (Surr)	98	8.920	8.922	-0.002	99	1858068	175.0	157.6	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.530	11.532	-0.002	98	701915	175.0	165.3	
11 Dichlorodifluoromethane	85	1.614	1.622	-0.008	99	432190	175.0	179.3	
12 Chloromethane	50	1.772	1.768	0.004	100	573343	175.0	172.3	
13 Vinyl chloride	62	1.906	1.896	0.010	100	624000	175.0	167.8	
14 Butadiene	39	1.948	1.944	0.004	99	709784	175.0	167.1	
15 Bromomethane	94	2.252	2.249	0.003	100	307964	175.0	162.9	
16 Chloroethane	64	2.380	2.376	0.004	98	455903	175.0	177.2	
17 Dichlorofluoromethane	67	2.648	2.644	0.004	100	974888	175.0	166.0	
18 Trichlorofluoromethane	101	2.703	2.723	-0.020	98	772293	175.0	173.1	
20 Ethyl ether	59	3.086	3.082	0.004	99	519119	175.0	176.5	
21 Acrolein	56	3.250	3.258	-0.008	100	81646	225.0	228.5	
22 1,1-Dichloroethene	96	3.372	3.374	-0.002	98	562804	175.0	173.6	
23 1,1,2-Trichloro-1,2,2-trif	101	3.427	3.423	0.004	98	577719	175.0	176.1	
24 Acetone	43	3.493	3.496	-0.003	100	429781	350.0	373.1	
25 Iodomethane	142	3.573	3.581	-0.008	100	784350	175.0	174.1	
26 Carbon disulfide	76	3.652	3.660	-0.008	100	1381152	175.0	174.2	
28 3-Chloro-1-propene	76	3.931	3.934	-0.003	100	314052	175.0	183.2	
30 Methyl acetate	43	4.017	4.019	-0.002	100	2407305	875.0	893.2	
31 Methylene Chloride	84	4.138	4.147	-0.009	97	597904	175.0	159.4	
32 2-Methyl-2-propanol	59	4.442	4.439	0.003	99	351016	1750.0	1728.4	
33 Acrylonitrile	53	4.546	4.554	-0.008	99	2446379	1750.0	1764.6	
34 trans-1,2-Dichloroethene	96	4.558	4.560	-0.002	92	581552	175.0	173.4	
35 Methyl tert-butyl ether	73	4.595	4.597	-0.002	98	1347848	175.0	181.8	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
36 Hexane	57	4.978	4.980	-0.002	99	929791	175.0	173.5	
37 1,1-Dichloroethane	63	5.166	5.175	-0.009	100	1052201	175.0	175.7	
38 Vinyl acetate	43	5.294	5.296	-0.002	100	831670	175.0	195.8	
44 2,2-Dichloropropane	77	5.927	5.929	-0.002	98	280515	175.0	187.4	
45 cis-1,2-Dichloroethene	96	5.933	5.941	-0.008	98	612812	175.0	173.4	
46 2-Butanone (MEK)	43	5.982	5.990	-0.008	100	665013	350.0	361.1	
49 Chlorobromomethane	128	6.225	6.227	-0.002	99	269375	175.0	176.2	
51 Tetrahydrofuran	42	6.286	6.282	0.004	100	415944	350.0	360.6	
52 Chloroform	83	6.341	6.343	-0.003	100	953676	175.0	175.3	
53 1,1,1-Trichloroethane	97	6.529	6.531	-0.002	99	639960	175.0	184.3	
54 Cyclohexane	56	6.584	6.586	-0.002	99	1161488	175.0	174.2	
56 Carbon tetrachloride	117	6.718	6.720	-0.002	100	504991	175.0	181.2	
55 1,1-Dichloropropene	75	6.724	6.726	-0.002	100	783682	175.0	173.7	
57 Isobutyl alcohol	41	6.943	6.945	-0.002	98	386141	4375.0	5144.3	
58 Benzene	78	6.955	6.957	-0.002	98	2286079	175.0	171.5	
59 1,2-Dichloroethane	62	6.985	6.988	-0.003	99	781760	175.0	179.2	
62 n-Heptane	43	7.277	7.280	-0.003	90	819785	175.0	179.1	
64 Trichloroethene	130	7.667	7.669	-0.002	99	586010	175.0	175.5	
66 Methylcyclohexane	83	7.861	7.864	-0.003	100	1055175	175.0	177.1	
67 1,2-Dichloropropane	63	7.904	7.906	-0.002	98	597514	175.0	181.3	
68 Dibromomethane	93	8.026	8.022	0.004	100	308441	175.0	173.8	
70 1,4-Dioxane	88	8.056	8.058	-0.002	97	132396	3500.0	3814.7	
71 Dichlorobromomethane	83	8.196	8.198	-0.002	100	663337	175.0	183.2	
74 cis-1,3-Dichloropropene	75	8.658	8.661	-0.003	100	681682	175.0	195.1	
75 4-Methyl-2-pentanone (MIBK)	43	8.823	8.825	-0.002	99	1390980	350.0	347.5	
76 Toluene	91	8.993	8.989	0.004	99	2347437	175.0	154.9	
77 trans-1,3-Dichloropropene	75	9.218	9.220	-0.002	98	502980	175.0	183.7	
78 Ethyl methacrylate	69	9.315	9.318	-0.003	98	654210	175.0	183.2	
79 1,1,2-Trichloroethane	97	9.400	9.403	-0.003	100	465584	175.0	163.8	
80 Tetrachloroethene	164	9.534	9.537	-0.003	99	477004	175.0	160.9	
81 1,3-Dichloropropane	76	9.565	9.561	0.004	99	854593	175.0	161.7	
82 2-Hexanone	43	9.656	9.658	-0.002	100	1103034	350.0	360.6	
84 Chlorodibromomethane	129	9.790	9.792	-0.002	99	406960	175.0	179.4	
85 Ethylene Dibromide	107	9.899	9.902	-0.003	99	461219	175.0	170.0	
86 3-Chlorobenzotrifluoride	180	10.374	10.370	0.004	87	925933	175.0	160.1	
87 Chlorobenzene	112	10.392	10.388	0.004	99	1507544	175.0	157.0	
88 4-Chlorobenzotrifluoride	180	10.429	10.431	-0.002	99	908777	175.0	162.5	
89 1,1,1,2-Tetrachloroethane	131	10.471	10.473	-0.002	95	439701	175.0	177.3	
90 Ethylbenzene	106	10.502	10.504	-0.002	99	889389	175.0	161.4	
91 m-Xylene & p-Xylene	106	10.617	10.619	-0.002	99	1092005	175.0	162.0	
92 o-Xylene	106	11.013	11.009	0.004	98	1059986	175.0	160.8	
93 Styrene	104	11.025	11.027	-0.002	93	1723778	175.0	162.3	
94 Bromoform	173	11.213	11.209	0.004	99	253560	175.0	180.9	
96 2-Chlorobenzotrifluoride	180	11.274	11.276	-0.002	99	922108	175.0	159.7	
97 Isopropylbenzene	105	11.378	11.380	-0.002	98	2580136	175.0	156.9	
99 1,1,2,2-Tetrachloroethane	83	11.676	11.678	-0.002	99	681581	175.0	167.2	
100 Bromobenzene	156	11.682	11.678	0.004	98	637569	175.0	171.0	
101 1,2,3-Trichloropropane	110	11.718	11.721	-0.002	98	214358	175.0	174.9	
102 trans-1,4-Dichloro-2-butene	53	11.730	11.733	-0.003	98	180624	175.0	177.3	
103 N-Propylbenzene	120	11.791	11.787	0.004	99	780243	175.0	169.6	
104 2-Chlorotoluene	126	11.876	11.873	0.003	99	666866	175.0	172.6	
105 3-Chlorotoluene	126	11.937	11.933	0.004	97	757051	175.0	175.3	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
106 1,3,5-Trimethylbenzene	105	11.962	11.964	-0.002	99	2136446	175.0	166.6	
107 4-Chlorotoluene	126	11.986	11.982	0.004	97	711885	175.0	170.2	
108 tert-Butylbenzene	119	12.290	12.286	0.004	100	1828125	175.0	164.5	
110 1,2,4-Trimethylbenzene	105	12.339	12.335	0.004	98	2187785	175.0	166.2	
111 1,2-dichloro-4-(trifluoromethyl)	214	12.400	12.402	-0.002	100	719294	175.0	172.9	
112 sec-Butylbenzene	105	12.509	12.511	-0.002	99	2565671	175.0	164.1	
113 1,3-Dichlorobenzene	146	12.619	12.621	-0.002	99	1159025	175.0	168.7	
114 4-Isopropyltoluene	119	12.655	12.651	0.004	98	2157955	175.0	167.2	
115 1,4-Dichlorobenzene	146	12.704	12.706	-0.002	99	1196958	175.0	170.6	
116 2,4-Dichloro-1-(trifluoromethyl)	214	12.759	12.755	0.004	99	675783	175.0	173.5	
118 2,5-Dichlorobenzotrifluoride	214	12.807	12.803	0.004	99	748317	175.0	171.7	
120 n-Butylbenzene	91	13.063	13.065	-0.002	99	1983203	175.0	168.7	
121 1,2-Dichlorobenzene	146	13.081	13.083	-0.002	99	1092014	175.0	171.7	
122 1,2-Dibromo-3-Chloropropan	75	13.860	13.856	0.004	95	97714	175.0	187.7	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125	14.006	14.008	-0.002	99	2487475	525.0	517.3	
125 2,3- & 3,4- Dichlorotoluene	125	14.425	14.428	-0.003	98	1639357	350.0	350.6	
126 1,2,4-Trichlorobenzene	180	14.693	14.695	-0.002	100	608110	175.0	183.6	
127 Hexachlorobutadiene	225	14.863	14.866	-0.003	98	274932	175.0	173.1	
128 Naphthalene	128	14.943	14.939	0.004	100	1599300	175.0	183.9	
129 1,2,3-Trichlorobenzene	180	15.186	15.188	-0.002	100	504504	175.0	185.8	
131 2,4,5-Trichlorotoluene	159	15.965	15.967	-0.002	99	273662	175.0	187.4	
130 2,3,6-Trichlorotoluene	159	16.062	16.064	-0.002	99	246163	175.0	186.7	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
S 133 Xylenes, Total	106				0		350.0	322.8	
S 134 1,2-Dichloroethene, Total	96				0		350.0	346.8	
S 135 1,3-Dichloropropene, Total	1				0		350.0	378.8	

### QC Flag Legend

#### Processing Flags

ND - Not Detected or Marked ND

#### Reagents:

VOAACRPRI_00003	Amount Added: 9.00	Units: uL	
voaWEEpri Res_00003	Amount Added: 7.00	Units: uL	
voaWKetpri Re_00003	Amount Added: 7.00	Units: uL	
VOA8260SURR_00032	Amount Added: 7.00	Units: uL	
VOA8260VOAPRI_00105	Amount Added: 7.00	Units: uL	
VOAVAPRI_00005	Amount Added: 7.00	Units: uL	
VOA8260INT_00030	Amount Added: 2.00	Units: uL	Run Reagent

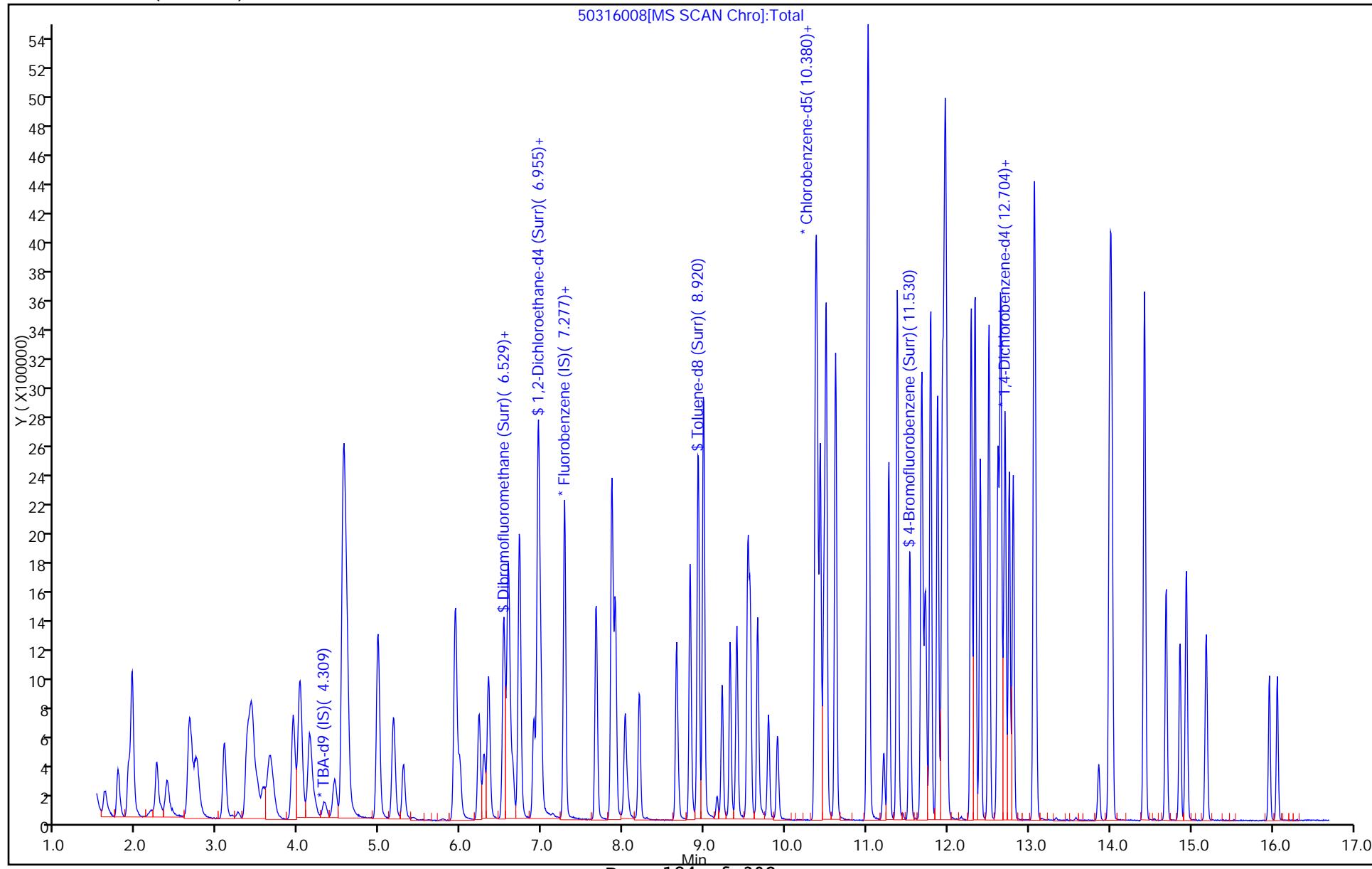
Report Date: 17-Mar-2015 10:59:30

Chrom Revision: 2.2 15-Jan-2015 13:05:58

TestAmerica Pittsburgh

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150316-6031.b\\50316008.D  
Injection Date: 16-Mar-2015 14:17:30 Instrument ID: CHHP5  
Lims ID: IC VSTD35 Operator ID: 001562  
Client ID:  
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 8  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)

Worklist Smp#: 8



TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316009.D  
 Lims ID: IC VSTD40  
 Client ID:  
 Sample Type: IC Calib Level: 7  
 Inject. Date: 16-Mar-2015 14:41:30 ALS Bottle#: 9 Worklist Smp#: 9  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: IC VSTD40  
 Misc. Info.: 180-0006031-009  
 Operator ID: 001562 Instrument ID: CHHP5  
 Sublist: chrom-MSVOA\_LL\_CHHP5\*sub4  
 Method: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 17-Mar-2015 10:59:31 Calib Date: 16-Mar-2015 16:17:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICal File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316013.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK012

First Level Reviewer: fergusond Date: 17-Mar-2015 09:50:54

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.327	4.305	0.022	86	183503	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.271	7.273	-0.002	99	592746	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.368	10.364	0.004	94	147746	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.685	12.682	0.003	94	203483	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.529	6.525	0.004	99	526164	200.0	195.2	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.900	6.902	-0.002	98	691002	200.0	194.4	
\$ 7 Toluene-d8 (Surr)	98	8.926	8.922	0.004	99	2153477	200.0	182.8	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.529	11.532	-0.003	98	798953	200.0	188.3	
11 Dichlorodifluoromethane	85	1.619	1.622	-0.003	99	522240	200.0	205.5	
12 Chloromethane	50	1.778	1.768	0.010	100	674845	200.0	192.4	
13 Vinyl chloride	62	1.905	1.896	0.009	100	767804	200.0	195.9	
14 Butadiene	39	1.948	1.944	0.004	98	840803	200.0	187.8	
15 Bromomethane	94	2.252	2.249	0.003	100	366671	200.0	184.6	
16 Chloroethane	64	2.374	2.376	-0.002	99	530813	200.0	195.7	
17 Dichlorofluoromethane	67	2.654	2.644	0.010	99	1188936	200.0	192.0	
18 Trichlorofluoromethane	101	2.733	2.723	0.010	98	946313	200.0	201.3	
20 Ethyl ether	59	3.092	3.082	0.010	100	592652	200.0	191.1	
21 Acrolein	56	3.250	3.258	-0.008	100	95028	250.0	252.3	
22 1,1-Dichloroethene	96	3.378	3.374	0.004	98	662050	200.0	193.7	
23 1,1,2-Trichloro-1,2,2-trif	101	3.432	3.423	0.009	98	684103	200.0	197.9	
24 Acetone	43	3.505	3.496	0.009	100	489133	400.0	402.8	
25 Iodomethane	142	3.597	3.581	0.016	100	945860	200.0	199.2	
26 Carbon disulfide	76	3.664	3.660	0.004	100	1643948	200.0	196.7	
28 3-Chloro-1-propene	76	3.931	3.934	-0.003	99	393345	200.0	217.7	
30 Methyl acetate	43	4.022	4.019	0.003	99	2810332	1000.0	989.3	
31 Methylene Chloride	84	4.144	4.147	-0.003	98	703059	200.0	177.9	
32 2-Methyl-2-propanol	59	4.448	4.439	0.009	99	399281	2000.0	1847.2	
33 Acrylonitrile	53	4.552	4.554	-0.002	99	2868164	2000.0	1962.8	
34 trans-1,2-Dichloroethene	96	4.564	4.560	0.004	96	692220	200.0	195.8	
35 Methyl tert-butyl ether	73	4.600	4.597	0.003	98	1581345	200.0	202.3	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
36 Hexane	57	4.978	4.980	-0.002	100	1096478	200.0	194.1	
37 1,1-Dichloroethane	63	5.172	5.175	-0.003	100	1250453	200.0	198.2	
38 Vinyl acetate	43	5.294	5.296	-0.002	100	1001771	200.0	223.8	
44 2,2-Dichloropropane	77	5.927	5.929	-0.002	98	338302	200.0	214.4	
45 cis-1,2-Dichloroethene	96	5.933	5.941	-0.008	98	721075	200.0	193.6	
46 2-Butanone (MEK)	43	5.987	5.990	-0.003	100	809232	400.0	416.8	
49 Chlorobromomethane	128	6.225	6.227	-0.002	98	311076	200.0	193.0	
51 Tetrahydrofuran	42	6.286	6.282	0.004	99	483324	400.0	397.5	
52 Chloroform	83	6.340	6.343	-0.003	100	1109416	200.0	193.5	
53 1,1,1-Trichloroethane	97	6.529	6.531	-0.002	99	768585	200.0	210.0	
54 Cyclohexane	56	6.584	6.586	-0.002	99	1366913	200.0	194.5	
56 Carbon tetrachloride	117	6.717	6.720	-0.003	99	612080	200.0	208.4	
55 1,1-Dichloropropene	75	6.724	6.726	-0.002	99	933326	200.0	196.3	
57 Isobutyl alcohol	41	6.949	6.945	0.004	98	433313	5000.0	5476.7	
58 Benzene	78	6.955	6.957	-0.002	97	2653105	200.0	188.9	
59 1,2-Dichloroethane	62	6.985	6.988	-0.003	99	907622	200.0	197.3	
62 n-Heptane	43	7.277	7.280	-0.003	88	940924	200.0	195.0	
64 Trichloroethene	130	7.666	7.669	-0.003	99	684010	200.0	194.4	
66 Methylcyclohexane	83	7.861	7.864	-0.003	100	1212427	200.0	193.1	
67 1,2-Dichloropropane	63	7.904	7.906	-0.002	98	700921	200.0	201.7	
68 Dibromomethane	93	8.025	8.022	0.003	99	370624	200.0	198.1	
70 1,4-Dioxane	88	8.062	8.058	0.004	98	146272	4000.0	3998.4	
71 Dichlorobromomethane	83	8.196	8.198	-0.002	100	773432	200.0	202.6	
74 cis-1,3-Dichloropropene	75	8.658	8.661	-0.003	99	829306	200.0	225.1	
75 4-Methyl-2-pentanone (MIBK)	43	8.822	8.825	-0.003	99	1617802	400.0	404.7	
76 Toluene	91	8.993	8.989	0.004	99	2714932	200.0	179.3	
77 trans-1,3-Dichloropropene	75	9.218	9.220	-0.002	98	613747	200.0	224.4	
78 Ethyl methacrylate	69	9.315	9.318	-0.003	98	782394	200.0	219.4	
79 1,1,2-Trichloroethane	97	9.400	9.403	-0.003	99	540864	200.0	190.5	
80 Tetrachloroethene	164	9.540	9.537	0.003	99	545517	200.0	184.2	
81 1,3-Dichloropropane	76	9.564	9.561	0.003	99	1001573	200.0	189.8	
82 2-Hexanone	43	9.656	9.658	-0.002	100	1305223	400.0	427.2	
84 Chlorodibromomethane	129	9.790	9.792	-0.002	99	473922	200.0	209.1	
85 Ethylene Dibromide	107	9.899	9.902	-0.003	100	534328	200.0	197.2	
86 3-Chlorobenzotrifluoride	180	10.374	10.370	0.004	88	1122812	200.0	194.4	
87 Chlorobenzene	112	10.392	10.388	0.004	99	1745676	200.0	182.0	
88 4-Chlorobenzotrifluoride	180	10.428	10.431	-0.003	99	1108797	200.0	198.5	
89 1,1,1,2-Tetrachloroethane	131	10.477	10.473	0.004	95	512980	200.0	207.1	
90 Ethylbenzene	106	10.501	10.504	-0.003	98	1044399	200.0	189.8	
91 m-Xylene & p-Xylene	106	10.617	10.619	-0.002	98	1256840	200.0	186.7	
92 o-Xylene	106	11.012	11.009	0.003	99	1214164	200.0	184.4	
93 Styrene	104	11.025	11.027	-0.002	97	1958961	200.0	184.6	
94 Bromoform	173	11.213	11.209	0.004	98	293938	200.0	210.0	
96 2-Chlorobenzotrifluoride	180	11.274	11.276	-0.002	99	1120386	200.0	194.2	
97 Isopropylbenzene	105	11.377	11.380	-0.003	98	2885608	200.0	175.6	
99 1,1,2,2-Tetrachloroethane	83	11.675	11.678	-0.003	98	772016	200.0	189.6	
100 Bromobenzene	156	11.682	11.678	0.004	99	740842	200.0	196.7	
101 1,2,3-Trichloropropane	110	11.718	11.721	-0.002	96	233938	200.0	189.0	
102 trans-1,4-Dichloro-2-butene	53	11.730	11.733	-0.003	98	211691	200.0	205.7	
103 N-Propylbenzene	120	11.791	11.787	0.004	98	887838	200.0	191.1	
104 2-Chlorotoluene	126	11.870	11.873	-0.003	99	756732	200.0	193.9	
105 3-Chlorotoluene	126	11.937	11.933	0.004	97	890638	200.0	204.2	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
106 1,3,5-Trimethylbenzene	105	11.961	11.964	-0.003	99	2387945	200.0	184.4	
107 4-Chlorotoluene	126	11.980	11.982	-0.002	96	795532	200.0	188.3	
108 tert-Butylbenzene	119	12.290	12.286	0.004	99	2060731	200.0	183.6	
110 1,2,4-Trimethylbenzene	105	12.339	12.335	0.004	99	2461131	200.0	185.1	
111 1,2-dichloro-4-(trifluoromethyl)	214	12.399	12.402	-0.003	99	832435	200.0	198.1	
112 sec-Butylbenzene	105	12.509	12.511	-0.002	99	2854173	200.0	180.7	
113 1,3-Dichlorobenzene	146	12.618	12.621	-0.003	99	1308081	200.0	188.5	
114 4-Isopropyltoluene	119	12.655	12.651	0.004	98	2408127	200.0	184.7	
115 1,4-Dichlorobenzene	146	12.710	12.706	0.004	99	1348596	200.0	190.3	
116 2,4-Dichloro-1-(trifluoromethyl)	214	12.758	12.755	0.003	99	786683	200.0	199.9	
118 2,5-Dichlorobenzotrifluoride	214	12.807	12.803	0.004	99	877059	200.0	199.2	
120 n-Butylbenzene	91	13.062	13.065	-0.003	98	2209671	200.0	186.1	
121 1,2-Dichlorobenzene	146	13.081	13.083	-0.002	99	1224311	200.0	190.6	
122 1,2-Dibromo-3-Chloropropan	75	13.859	13.856	0.003	94	112547	200.0	214.1	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125	14.005	14.008	-0.003	98	2860911	600.0	589.0	
125 2,3- & 3,4- Dichlorotoluene	125	14.425	14.428	-0.003	98	1868280	400.0	395.5	
126 1,2,4-Trichlorobenzene	180	14.693	14.695	-0.002	99	679520	200.0	203.2	
127 Hexachlorobutadiene	225	14.863	14.866	-0.003	99	307470	200.0	191.7	
128 Naphthalene	128	14.942	14.939	0.003	100	1786434	200.0	203.4	
129 1,2,3-Trichlorobenzene	180	15.186	15.188	-0.002	99	582911	200.0	212.5	
131 2,4,5-Trichlorotoluene	159	15.964	15.967	-0.003	98	315499	200.0	213.9	
130 2,3,6-Trichlorotoluene	159	16.062	16.064	-0.002	98	285573	200.0	214.4	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
S 133 Xylenes, Total	106				0		400.0	371.1	
S 134 1,2-Dichloroethene, Total	96				0		400.0	389.4	
S 135 1,3-Dichloropropene, Total	1				0		400.0	449.6	

### QC Flag Legend

#### Processing Flags

ND - Not Detected or Marked ND

#### Reagents:

VOAVAPRI_00005	Amount Added: 8.00	Units: uL	
VOA8260VOAPRI_00105	Amount Added: 8.00	Units: uL	
VOA8260SURR_00032	Amount Added: 8.00	Units: uL	
voaWEEpri Res_00003	Amount Added: 8.00	Units: uL	
voaWKetpri Re_00003	Amount Added: 8.00	Units: uL	
VOAACRPRI_00003	Amount Added: 10.00	Units: uL	
VOA8260INT_00030	Amount Added: 2.00	Units: uL	Run Reagent

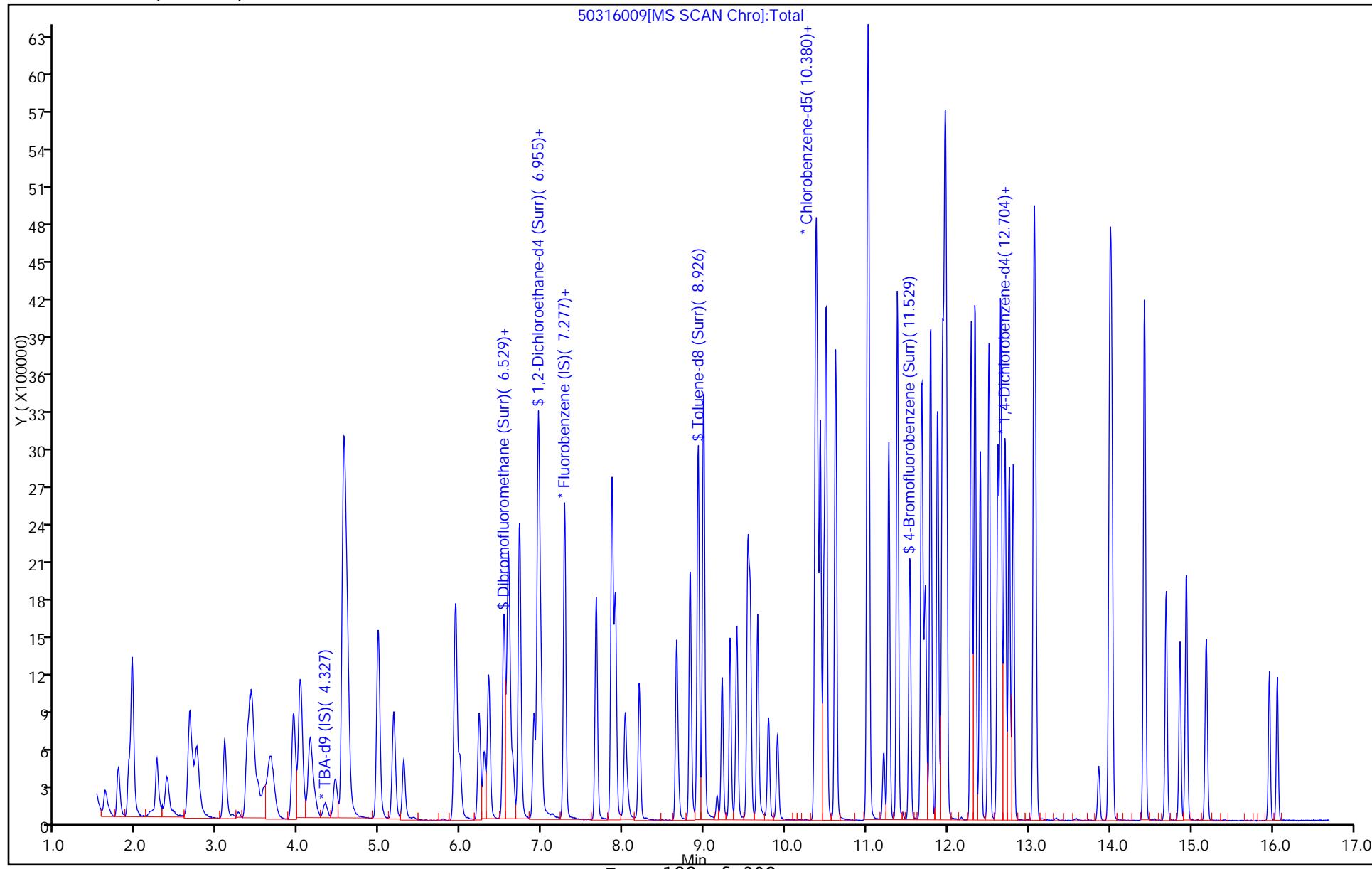
Report Date: 17-Mar-2015 10:59:31

Chrom Revision: 2.2 15-Jan-2015 13:05:58

TestAmerica Pittsburgh

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150316-6031.b\\50316009.D  
Injection Date: 16-Mar-2015 14:41:30 Instrument ID: CHHP5  
Lims ID: IC VSTD40 Operator ID: 001562  
Client ID:  
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 9  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)

Worklist Smp#: 9



TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316010.D  
 Lims ID: IC VSTD50  
 Client ID:  
 Sample Type: IC Calib Level: 8  
 Inject. Date: 16-Mar-2015 15:05:30 ALS Bottle#: 10 Worklist Smp#: 10  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: IC VSTD50  
 Misc. Info.: 180-0006031-010  
 Operator ID: 001562 Instrument ID: CHHP5  
 Sublist: chrom-MSVOA\_LL\_CHHP5\*sub4  
 Method: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 17-Mar-2015 10:59:32 Calib Date: 16-Mar-2015 16:17:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICal File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316013.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK012

First Level Reviewer: fergusond Date: 17-Mar-2015 09:55:54

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.336	4.305	0.031	85	202534	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.275	7.273	0.002	99	620293	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.359	10.364	-0.005	77	161503	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.683	12.682	0.001	92	212327	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.527	6.525	0.001	99	664693	250.0	235.6	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.898	6.902	-0.004	99	889045	250.0	239.0	
\$ 7 Toluene-d8 (Surr)	98	8.923	8.922	0.001	98	2632400	250.0	204.4	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.533	11.532	0.001	98	1045249	250.0	225.4	
11 Dichlorodifluoromethane	85	1.617	1.622	-0.005	99	640090	250.0	240.7	
12 Chloromethane	50	1.775	1.768	0.007	100	855933	250.0	233.2	
13 Vinyl chloride	62	1.909	1.896	0.013	100	924535	250.0	225.4	
14 Butadiene	39	1.946	1.944	0.002	99	1005925	250.0	214.7	
15 Bromomethane	94	2.250	2.249	0.001	100	461680	250.0	223.1	
16 Chloroethane	64	2.371	2.376	-0.005	99	700467	250.0	246.8	
17 Dichlorofluoromethane	67	2.651	2.644	0.007	100	1511714	250.0	233.3	
18 Trichlorofluoromethane	101	2.724	2.723	0.001	98	1178605	250.0	239.5	
20 Ethyl ether	59	3.083	3.082	0.001	99	792637	250.0	244.3	
21 Acrolein	56	3.254	3.258	-0.004	96	109180	275.0	277.0	
22 1,1-Dichloroethene	96	3.375	3.374	0.001	98	827120	250.0	231.2	
23 1,1,2-Trichloro-1,2,2-trif	101	3.424	3.423	0.001	99	834802	250.0	230.7	
24 Acetone	43	3.497	3.496	0.001	99	621064	500.0	488.8	
25 Iodomethane	142	3.594	3.581	0.013	100	1201056	250.0	241.7	
26 Carbon disulfide	76	3.655	3.660	-0.005	100	2031733	250.0	232.3	
28 3-Chloro-1-propene	76	3.935	3.934	0.001	99	482122	250.0	255.0	
30 Methyl acetate	43	4.014	4.019	-0.005	99	3718382	1250.0	1250.8	
31 Methylene Chloride	84	4.142	4.147	-0.005	98	919183	250.0	222.2	
32 2-Methyl-2-propanol	59	4.446	4.439	0.007	98	537174	2500.0	2251.6	
33 Acrylonitrile	53	4.549	4.554	-0.005	99	3721902	2500.0	2433.9	
34 trans-1,2-Dichloroethene	96	4.562	4.560	0.002	97	882651	250.0	238.6	
35 Methyl tert-butyl ether	73	4.598	4.597	0.001	98	2130684	250.0	260.5	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
36 Hexane	57	4.975	4.980	-0.005	99	1379168	250.0	233.3	
37 1,1-Dichloroethane	63	5.170	5.175	-0.005	99	1604398	250.0	242.9	
38 Vinyl acetate	43	5.292	5.296	-0.004	100	1337263	250.0	285.5	
44 2,2-Dichloropropane	77	5.924	5.929	-0.005	97	452022	250.0	273.8	
45 cis-1,2-Dichloroethene	96	5.936	5.941	-0.005	97	930230	250.0	238.7	
46 2-Butanone (MEK)	43	5.985	5.990	-0.005	100	1059138	500.0	521.3	
49 Chlorobromomethane	128	6.222	6.227	-0.005	99	404105	250.0	239.6	
51 Tetrahydrofuran	42	6.283	6.282	0.001	99	646482	500.0	508.1	
52 Chloroform	83	6.338	6.343	-0.005	100	1424461	250.0	237.4	
53 1,1,1-Trichloroethane	97	6.527	6.531	-0.005	99	971626	250.0	253.6	
54 Cyclohexane	56	6.581	6.586	-0.005	98	1669676	250.0	227.0	
56 Carbon tetrachloride	117	6.715	6.720	-0.005	99	790495	250.0	257.2	
55 1,1-Dichloropropene	75	6.721	6.726	-0.005	99	1159811	250.0	233.1	
57 Isobutyl alcohol	41	6.946	6.945	0.001	97	644697	6250.0	7786.6	
58 Benzene	78	6.952	6.957	-0.005	97	3351151	250.0	228.0	
59 1,2-Dichloroethane	62	6.983	6.988	-0.005	99	1159879	250.0	241.0	
62 n-Heptane	43	7.275	7.280	-0.005	86	1182643	250.0	234.2	
64 Trichloroethene	130	7.664	7.669	-0.005	99	860273	250.0	233.6	
66 Methylcyclohexane	83	7.859	7.864	-0.005	99	1519674	250.0	231.3	
67 1,2-Dichloropropane	63	7.907	7.906	0.001	99	918714	250.0	252.7	
68 Dibromomethane	93	8.023	8.022	0.001	99	479407	250.0	244.9	
70 1,4-Dioxane	88	8.053	8.058	-0.005	98	185631	5000.0	4848.9	
71 Dichlorobromomethane	83	8.199	8.198	0.001	100	1003399	250.0	251.2	
74 cis-1,3-Dichloropropene	75	8.656	8.661	-0.005	99	1098242	250.0	284.9	
75 4-Methyl-2-pentanone (MIBK)	43	8.820	8.825	-0.005	98	2109966	500.0	482.8	
76 Toluene	91	8.990	8.989	0.001	97	3368812	250.0	203.5	
77 trans-1,3-Dichloropropene	75	9.221	9.220	0.001	98	846559	250.0	283.2	
78 Ethyl methacrylate	69	9.319	9.318	0.001	98	1063861	250.0	272.9	
79 1,1,2-Trichloroethane	97	9.398	9.403	-0.005	99	706748	250.0	227.7	
80 Tetrachloroethene	164	9.538	9.537	0.001	99	690601	250.0	213.3	
81 1,3-Dichloropropane	76	9.562	9.561	0.001	100	1327847	250.0	230.2	
82 2-Hexanone	43	9.653	9.658	-0.005	99	1685534	500.0	504.7	
84 Chlorodibromomethane	129	9.787	9.792	-0.005	99	625118	250.0	252.3	
85 Ethylene Dibromide	107	9.903	9.902	0.001	99	713501	250.0	240.9	
86 3-Chlorobenzotrifluoride	180	10.371	10.370	0.001	87	1303041	250.0	206.4	
87 Chlorobenzene	112	10.390	10.388	0.002	98	2249414	250.0	214.5	
88 4-Chlorobenzotrifluoride	180	10.426	10.431	-0.005	99	1250140	250.0	204.8	
89 1,1,1,2-Tetrachloroethane	131	10.475	10.473	0.002	95	680608	250.0	251.4	
90 Ethylbenzene	106	10.499	10.504	-0.005	97	1329470	250.0	221.0	
91 m-Xylene & p-Xylene	106	10.621	10.619	0.002	97	1614511	250.0	219.4	
92 o-Xylene	106	11.010	11.009	0.001	94	1557898	250.0	216.4	
93 Styrene	104	11.022	11.027	-0.005	91	2525667	250.0	217.8	
94 Bromoform	173	11.211	11.209	0.002	99	395201	250.0	258.3	
96 2-Chlorobenzotrifluoride	180	11.272	11.276	-0.004	99	1298335	250.0	205.9	
97 Isopropylbenzene	105	11.381	11.380	0.001	97	3554151	250.0	197.9	
99 1,1,2,2-Tetrachloroethane	83	11.673	11.678	-0.005	98	1003707	250.0	225.5	
100 Bromobenzene	156	11.685	11.678	0.007	99	956763	250.0	243.5	
101 1,2,3-Trichloropropane	110	11.716	11.721	-0.004	97	325768	250.0	252.3	
102 trans-1,4-Dichloro-2-butene	53	11.728	11.733	-0.005	98	286166	250.0	266.5	
103 N-Propylbenzene	120	11.789	11.787	0.002	97	1131297	250.0	233.4	
104 2-Chlorotoluene	126	11.874	11.873	0.001	97	963573	250.0	236.6	
105 3-Chlorotoluene	126	11.935	11.933	0.002	96	1053875	250.0	231.5	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
106 1,3,5-Trimethylbenzene	105	11.965	11.964	0.001	98	2983647	250.0	220.7	
107 4-Chlorotoluene	126	11.983	11.982	0.001	97	1062581	250.0	241.0	
108 tert-Butylbenzene	119	12.288	12.286	0.002	98	2516209	250.0	214.9	
110 1,2,4-Trimethylbenzene	105	12.336	12.335	0.001	97	3068942	250.0	221.2	
111 1,2-dichloro-4-(trifluoromethyl)	214	12.403	12.402	0.001	99	991010	250.0	226.1	
112 sec-Butylbenzene	105	12.507	12.511	-0.004	98	3463106	250.0	210.1	
113 1,3-Dichlorobenzene	146	12.616	12.621	-0.005	98	1687649	250.0	233.1	
114 4-Isopropyltoluene	119	12.653	12.651	0.002	97	2970922	250.0	218.3	
115 1,4-Dichlorobenzene	146	12.707	12.706	0.001	98	1736319	250.0	234.8	
116 2,4-Dichloro-1-(trifluoromethyl)	214	12.756	12.755	0.001	98	909481	250.0	221.5	
118 2,5-Dichlorobenzotrifluoride	214	12.811	12.803	0.008	99	1042359	250.0	226.9	
120 n-Butylbenzene	91	13.060	13.065	-0.005	97	2715831	250.0	219.2	
121 1,2-Dichlorobenzene	146	13.078	13.083	-0.005	99	1565775	250.0	233.6	
122 1,2-Dibromo-3-Chloropropan	75	13.863	13.856	0.007	94	147059	250.0	268.1	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125	14.003	14.008	-0.005	98	3379751	750.0	666.8	
125 2,3- & 3,4- Dichlorotoluene	125	14.423	14.428	-0.005	97	2218229	500.0	450.0	
126 1,2,4-Trichlorobenzene	180	14.691	14.695	-0.004	99	825772	250.0	236.6	
127 Hexachlorobutadiene	225	14.861	14.866	-0.005	99	367792	250.0	219.8	
128 Naphthalene	128	14.940	14.939	0.001	99	2220927	250.0	242.4	
129 1,2,3-Trichlorobenzene	180	15.189	15.188	0.001	99	697862	250.0	243.8	
131 2,4,5-Trichlorotoluene	159	15.962	15.967	-0.005	99	364223	250.0	236.6	
130 2,3,6-Trichlorotoluene	159	16.065	16.064	0.001	98	323920	250.0	233.1	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
S 133 Xylenes, Total	106				0		500.0	435.9	
S 134 1,2-Dichloroethene, Total	96				0		500.0	477.2	
S 135 1,3-Dichloropropene, Total	1				0		500.0	568.1	

### QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

### Reagents:

VOAACRPRI_00003	Amount Added: 11.00	Units: uL	
voaWEEpri Res_00003	Amount Added: 10.00	Units: uL	
voaWKetpri Re_00003	Amount Added: 10.00	Units: uL	
VOA8260SURR_00032	Amount Added: 10.00	Units: uL	
VOA8260VOAPRI_00105	Amount Added: 10.00	Units: uL	
VOAVAPRI_00005	Amount Added: 10.00	Units: uL	
VOA8260INT_00030	Amount Added: 2.00	Units: uL	Run Reagent

Report Date: 17-Mar-2015 10:59:33

Chrom Revision: 2.2 15-Jan-2015 13:05:58

TestAmerica Pittsburgh

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150316-6031.b\\50316010.D

Injection Date: 16-Mar-2015 15:05:30

Instrument ID: CHHP5

Operator ID: 001562

Lims ID: IC VSTD50

Worklist Smp#: 10

Client ID:

Purge Vol: 5.000 mL

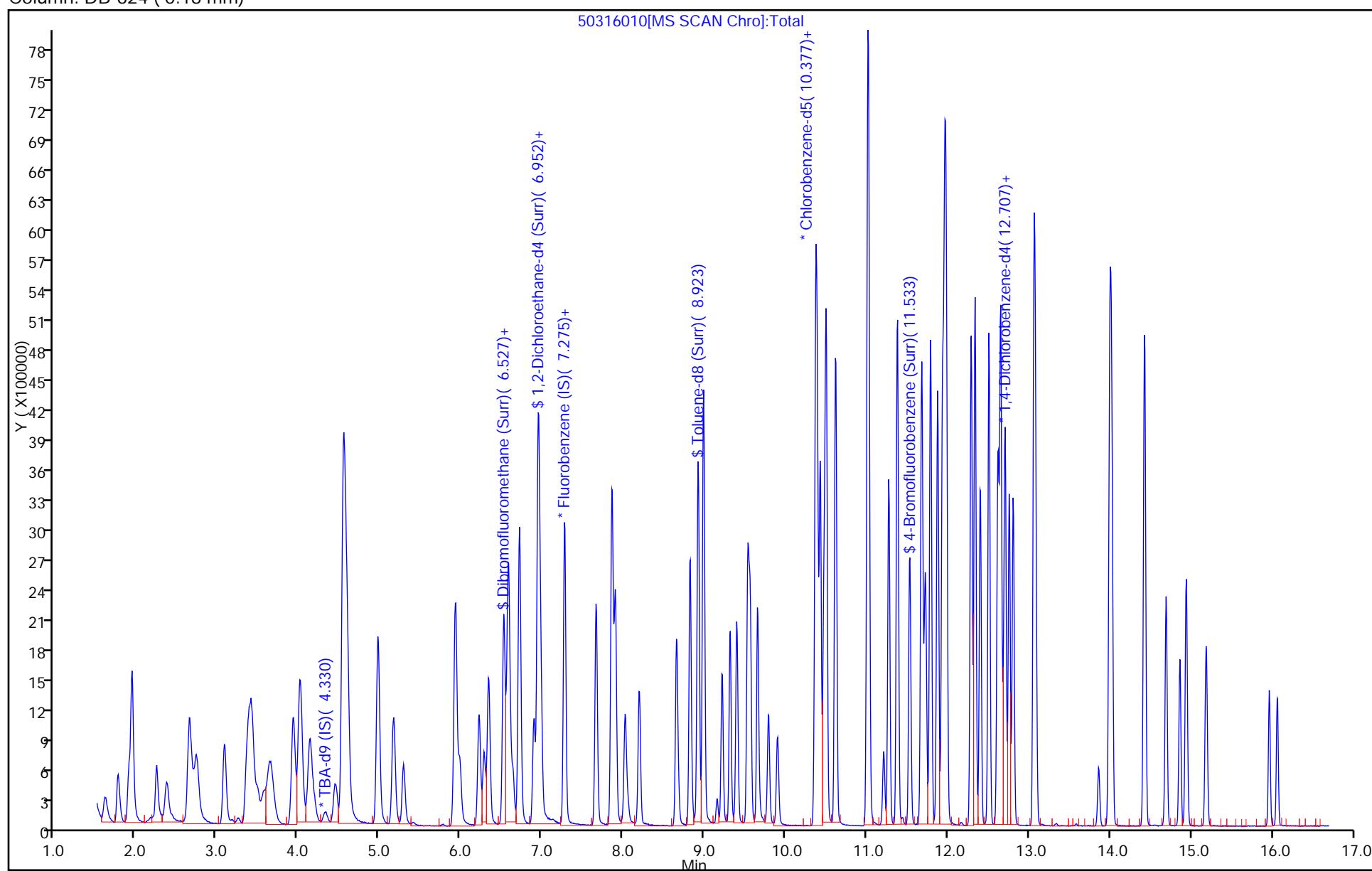
Dil. Factor: 1.0000

ALS Bottle#: 10

Method: MSVOA\_LL\_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 ( 0.18 mm)



TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316013.D  
 Lims ID: IC VSTD1  
 Client ID:  
 Sample Type: IC Calib Level: 1  
 Inject. Date: 16-Mar-2015 16:17:30 ALS Bottle#: 13 Worklist Smp#: 13  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: IC VSTD1  
 Misc. Info.: 180-0006031-013  
 Operator ID: 001562 Instrument ID: CHHP5  
 Sublist: chrom-MSVOA\_LL\_CHHP5\*sub4  
 Method: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 17-Mar-2015 10:59:33 Calib Date: 16-Mar-2015 16:17:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICal File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316013.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK012

First Level Reviewer: fergusond Date: 17-Mar-2015 10:01:36

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.317	4.305	0.012	83	148007	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.274	7.273	0.001	99	568509	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.364	10.364	0.000	74	121234	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.682	12.682	0.000	96	175081	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.526	6.525	0.001	93	14193	5.00	5.49	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.903	6.902	0.001	96	17152	5.00	5.03	
\$ 7 Toluene-d8 (Surr)	98	8.923	8.922	0.001	98	54935	5.00	5.68	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.532	11.532	0.000	92	19061	5.00	5.48	
11 Dichlorodifluoromethane	85	1.616	1.622	-0.006	96	11265	5.00	4.62	
12 Chloromethane	50	1.768	1.768	0.000	97	17972	5.00	5.34	M
13 Vinyl chloride	62	1.908	1.896	0.012	96	18981	5.00	5.05	
14 Butadiene	39	1.951	1.944	0.007	98	24095	5.00	5.61	
15 Bromomethane	94	2.249	2.249	0.000	85	18060	5.00	4.90	
16 Chloroethane	64	2.377	2.376	0.001	53	13187	5.00	5.07	
17 Dichlorofluoromethane	67	2.644	2.644	0.000	99	34297	5.00	5.78	
18 Trichlorofluoromethane	101	2.711	2.723	-0.012	92	20521	5.00	4.55	
20 Ethyl ether	59	3.082	3.082	0.000	94	16416	5.00	5.52	
21 Acrolein	56	3.247	3.258	-0.011	96	35289	100.0	97.7	M
22 1,1-Dichloroethene	96	3.368	3.374	-0.006	97	18234	5.00	5.56	
23 1,1,2-Trichloro-1,2,2-trif	101	3.435	3.423	0.012	90	16567	5.00	5.00	
24 Acetone	43	3.490	3.496	-0.006	93	29674	25.0	25.5	
25 Iodomethane	142	3.581	3.581	0.000	97	22824	5.00	5.01	
26 Carbon disulfide	76	3.648	3.660	-0.012	98	41336	5.00	5.16	
28 3-Chloro-1-propene	76	3.940	3.934	0.006	95	8006	5.00	4.62	
30 Methyl acetate	43	4.031	4.019	0.012	100	71022	25.0	26.1	
31 Methylene Chloride	84	4.135	4.147	-0.012	96	27978	5.00	7.38	
32 2-Methyl-2-propanol	59	4.433	4.439	-0.006	73	10830	50.0	62.1	
33 Acrylonitrile	53	4.555	4.554	0.001	99	71728	50.0	51.2	
34 trans-1,2-Dichloroethene	96	4.555	4.560	-0.005	57	17111	5.00	5.05	
35 Methyl tert-butyl ether	73	4.603	4.597	0.006	94	40058	5.00	5.34	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
36 Hexane	57	4.974	4.980	-0.006	96	29021	5.00	5.36	
37 1,1-Dichloroethane	63	5.169	5.175	-0.006	99	29622	5.00	4.89	
38 Vinyl acetate	43	5.297	5.296	0.001	79	19067	5.00	4.44	
44 2,2-Dichloropropane	77	5.936	5.929	0.007	87	6267	5.00	4.14	
45 cis-1,2-Dichloroethene	96	5.936	5.941	-0.005	95	18951	5.00	5.30	
46 2-Butanone (MEK)	43	5.996	5.990	0.006	99	42054	25.0	22.6	
49 Chlorobromomethane	128	6.234	6.227	0.007	95	8619	5.00	5.58	
51 Tetrahydrofuran	42	6.288	6.282	0.006	75	11913	10.0	10.2	
52 Chloroform	83	6.343	6.343	0.000	97	29168	5.00	5.30	
53 1,1,1-Trichloroethane	97	6.532	6.531	0.001	93	15663	5.00	4.46	
54 Cyclohexane	56	6.580	6.586	-0.006	94	36280	5.00	5.38	
56 Carbon tetrachloride	117	6.720	6.720	0.000	97	13013	5.00	4.62	
55 1,1-Dichloropropene	75	6.720	6.726	-0.006	97	24060	5.00	5.28	
57 Isobutyl alcohol	41	6.958	6.945	0.013	95	8820	125.0	116.2	
58 Benzene	78	6.958	6.957	0.001	96	73700	5.00	5.47	
59 1,2-Dichloroethane	62	6.976	6.988	-0.012	98	22108	5.00	5.01	
62 n-Heptane	43	7.274	7.280	-0.006	58	23490	5.00	5.08	
64 Trichloroethene	130	7.669	7.669	0.000	96	18397	5.00	5.45	
66 Methylcyclohexane	83	7.858	7.864	-0.006	94	29934	5.00	4.97	
67 1,2-Dichloropropane	63	7.907	7.906	0.001	90	16916	5.00	5.08	
68 Dibromomethane	93	8.022	8.022	0.000	93	9562	5.00	5.33	
70 1,4-Dioxane	88	8.047	8.058	-0.012	33	3746	100.0	106.8	
71 Dichlorobromomethane	83	8.193	8.198	-0.006	98	16863	5.00	4.61	
74 cis-1,3-Dichloropropene	75	8.655	8.661	-0.006	98	15462	5.00	4.38	
75 4-Methyl-2-pentanone (MIBK)	43	8.831	8.825	0.006	99	75787	25.0	23.1	
76 Toluene	91	8.989	8.989	0.000	99	72597	5.00	5.84	
77 trans-1,3-Dichloropropene	75	9.208	9.220	-0.012	92	10481	5.00	4.67	
78 Ethyl methacrylate	69	9.318	9.318	0.000	94	13336	5.00	4.56	
79 1,1,2-Trichloroethane	97	9.403	9.403	0.000	95	13086	5.00	5.62	
80 Tetrachloroethene	164	9.531	9.537	-0.006	96	13716	5.00	5.64	
81 1,3-Dichloropropane	76	9.567	9.561	0.006	97	23188	5.00	5.35	
82 2-Hexanone	43	9.659	9.658	0.001	98	53734	25.0	21.4	M
84 Chlorodibromomethane	129	9.786	9.792	-0.006	95	7988	5.00	4.30	
85 Ethylene Dibromide	107	9.902	9.902	0.000	96	11471	5.00	5.16	
86 3-Chlorobenzotrifluoride	180	10.370	10.370	0.000	67	26148	5.00	5.52	
87 Chlorobenzene	112	10.389	10.388	0.001	98	47481	5.00	6.03	
88 4-Chlorobenzotrifluoride	180	10.425	10.431	-0.006	97	25927	5.00	5.66	
89 1,1,1,2-Tetrachloroethane	131	10.474	10.473	0.001	87	9154	5.00	4.50	
90 Ethylbenzene	106	10.504	10.504	0.000	99	24142	5.00	5.35	
91 m-Xylene & p-Xylene	106	10.614	10.619	-0.005	98	30126	5.00	5.45	
92 o-Xylene	106	11.015	11.009	0.006	97	32009	5.00	5.92	
93 Styrene	104	11.027	11.027	0.000	95	47061	5.00	5.41	
94 Bromoform	173	11.216	11.209	0.007	32	5157	5.00	4.49	
96 2-Chlorobenzotrifluoride	180	11.277	11.276	0.001	98	25441	5.00	5.37	
97 Isopropylbenzene	105	11.380	11.380	0.000	99	75470	5.00	5.60	
99 1,1,2,2-Tetrachloroethane	83	11.672	11.678	-0.006	93	19128	5.00	5.73	
100 Bromobenzene	156	11.691	11.678	0.012	97	16809	5.00	5.19	
101 1,2,3-Trichloropropane	110	11.721	11.721	0.001	89	5918	5.00	5.56	
102 trans-1,4-Dichloro-2-butene	53	11.739	11.733	0.006	50	4503	5.00	5.09	M
103 N-Propylbenzene	120	11.788	11.787	0.001	99	21543	5.00	5.39	
104 2-Chlorotoluene	126	11.873	11.873	0.000	99	17942	5.00	5.34	
105 3-Chlorotoluene	126	11.934	11.933	0.001	98	20174	5.00	5.37	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
106 1,3,5-Trimethylbenzene	105	11.964	11.964	0.000	99	61438	5.00	5.51	
107 4-Chlorotoluene	126	11.983	11.982	0.000	94	19812	5.00	5.45	
108 tert-Butylbenzene	119	12.287	12.286	0.001	97	55729	5.00	5.77	
110 1,2,4-Trimethylbenzene	105	12.335	12.335	0.000	97	63098	5.00	5.52	
111 1,2-dichloro-4-(trifluoromethyl)	214	12.408	12.402	0.006	98	19333	5.00	5.35	
112 sec-Butylbenzene	105	12.506	12.511	-0.005	100	75379	5.00	5.55	
113 1,3-Dichlorobenzene	146	12.621	12.621	0.000	98	33497	5.00	5.61	
114 4-Isopropyltoluene	119	12.652	12.651	0.001	98	61054	5.00	5.44	
115 1,4-Dichlorobenzene	146	12.706	12.706	0.000	98	34596	5.00	5.67	
116 2,4-Dichloro-1-(trifluoromethyl)	214	12.755	12.755	0.000	94	17792	5.00	5.26	
118 2,5-Dichlorobenzotrifluoride	214	12.810	12.803	0.007	96	20678	5.00	5.46	
120 n-Butylbenzene	91	13.065	13.065	0.000	99	54758	5.00	5.36	
121 1,2-Dichlorobenzene	146	13.084	13.083	0.001	99	30414	5.00	5.50	
122 1,2-Dibromo-3-Chloropropan	75	13.874	13.856	0.018	18	2299	5.00	5.08	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125	13.996	14.008	-0.012	93	71584	15.0	17.1	
125 2,3- & 3,4- Dichlorotoluene	125	14.428	14.428	0.000	97	46257	10.0	11.4	
126 1,2,4-Trichlorobenzene	180	14.690	14.695	-0.005	94	17018	5.00	5.91	
127 Hexachlorobutadiene	225	14.866	14.866	0.000	90	8549	5.00	6.19	
128 Naphthalene	128	14.939	14.939	0.000	99	41842	5.00	5.54	
129 1,2,3-Trichlorobenzene	180	15.189	15.188	0.000	95	13823	5.00	5.86	
131 2,4,5-Trichlorotoluene	159	15.961	15.967	-0.006	94	8592	5.00	6.77	
130 2,3,6-Trichlorotoluene	159	16.058	16.064	-0.006	94	7658	5.00	6.68	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
S 133 Xylenes, Total	106				0		10.0	11.4	
S 134 1,2-Dichloroethene, Total	96				0		10.0	10.4	
S 135 1,3-Dichloropropene, Total	1				0		10.0	9.05	

### QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

### Reagents:

VOA8260SURR_00032	Amount Added: 0.20	Units: uL	
VOA8260VOAPRI_00105	Amount Added: 0.20	Units: uL	
VOAVAPRI_00005	Amount Added: 0.20	Units: uL	
voaWKetpri Re_00003	Amount Added: 0.80	Units: uL	
VOAACRPRI_00003	Amount Added: 4.00	Units: uL	
voaWEEpri Res_00003	Amount Added: 0.20	Units: uL	
VOA8260INT_00030	Amount Added: 2.00	Units: uL	Run Reagent

Report Date: 17-Mar-2015 10:59:34

Chrom Revision: 2.2 15-Jan-2015 13:05:58

TestAmerica Pittsburgh

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150316-6031.b\\50316013.D

Injection Date: 16-Mar-2015 16:17:30

Instrument ID: CHHP5

Operator ID: 001562

Lims ID: IC VSTD1

Worklist Smp#: 13

Client ID:

Purge Vol: 5.000 mL

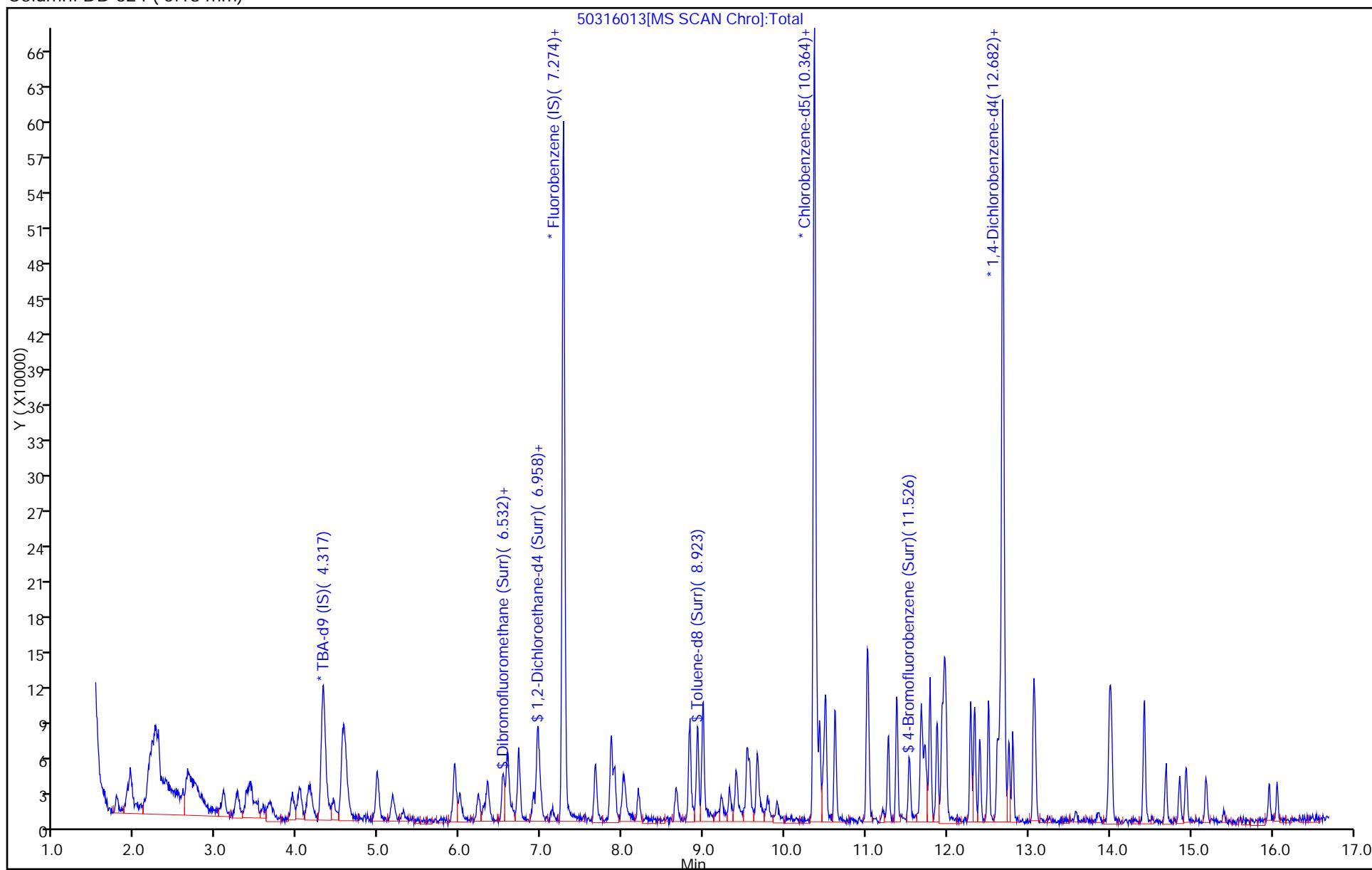
Dil. Factor: 1.0000

ALS Bottle#: 13

Method: MSVOA\_LL\_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 ( 0.18 mm)



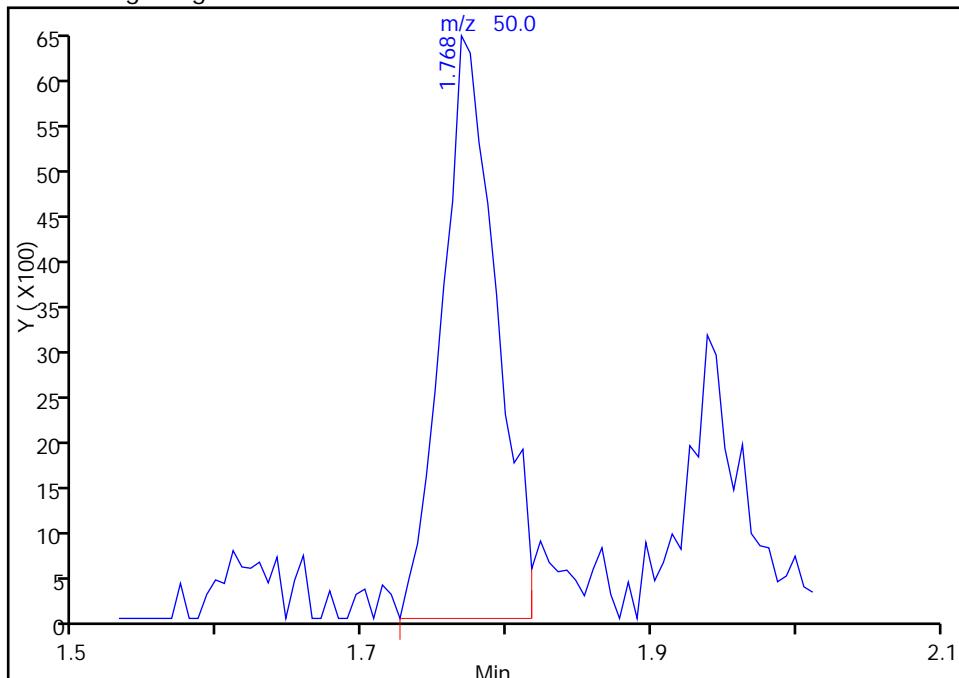
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316013.D  
 Injection Date: 16-Mar-2015 16:17:30 Instrument ID: CHHP5  
 Lims ID: IC VSTD1  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 13 Worklist Smp#: 13  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 12 Chloromethane, CAS: 74-87-3

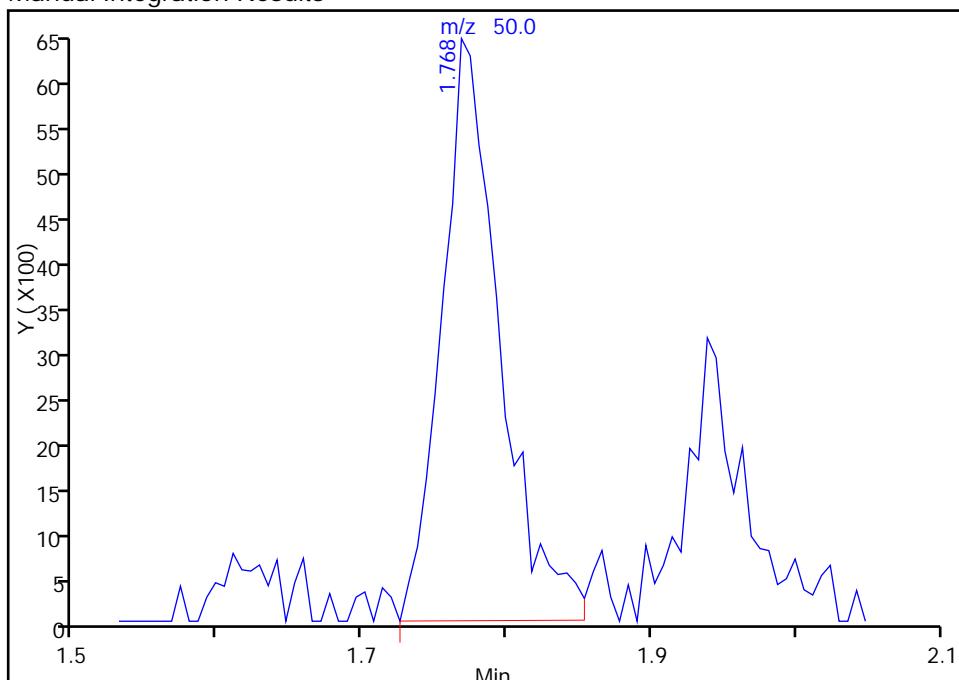
RT: 1.77  
 Area: 16860  
 Amount: 4.846171  
 Amount Units: ng

## Processing Integration Results



RT: 1.77  
 Area: 17972  
 Amount: 5.343308  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 17-Mar-2015 10:01:36

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

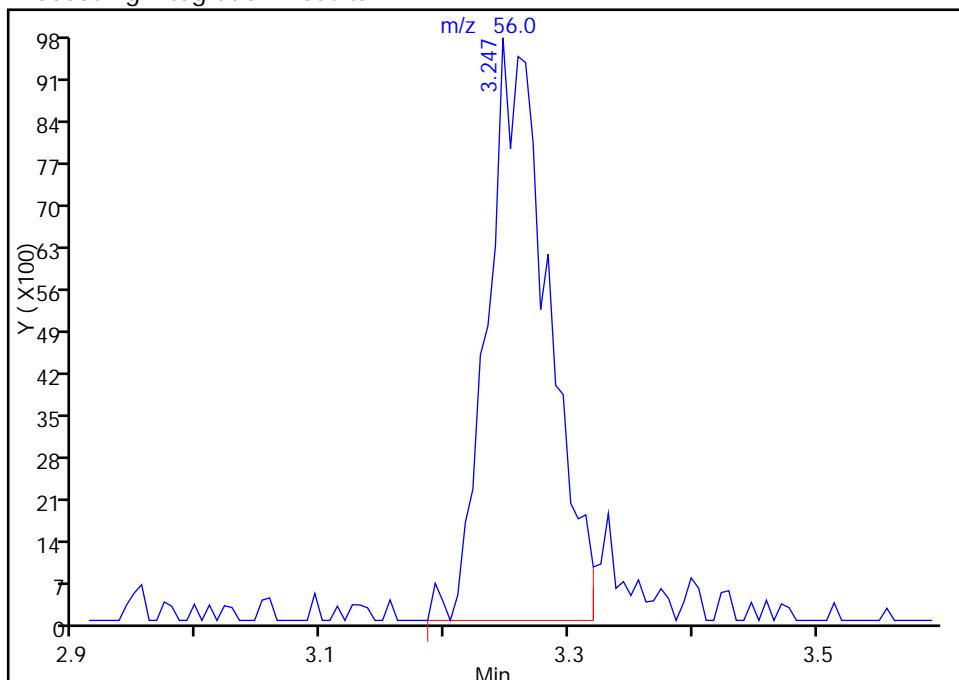
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316013.D  
 Injection Date: 16-Mar-2015 16:17:30 Instrument ID: CHHP5  
 Lims ID: IC VSTD1  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 13 Worklist Smp#: 13  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 21 Acrolein, CAS: 107-02-8

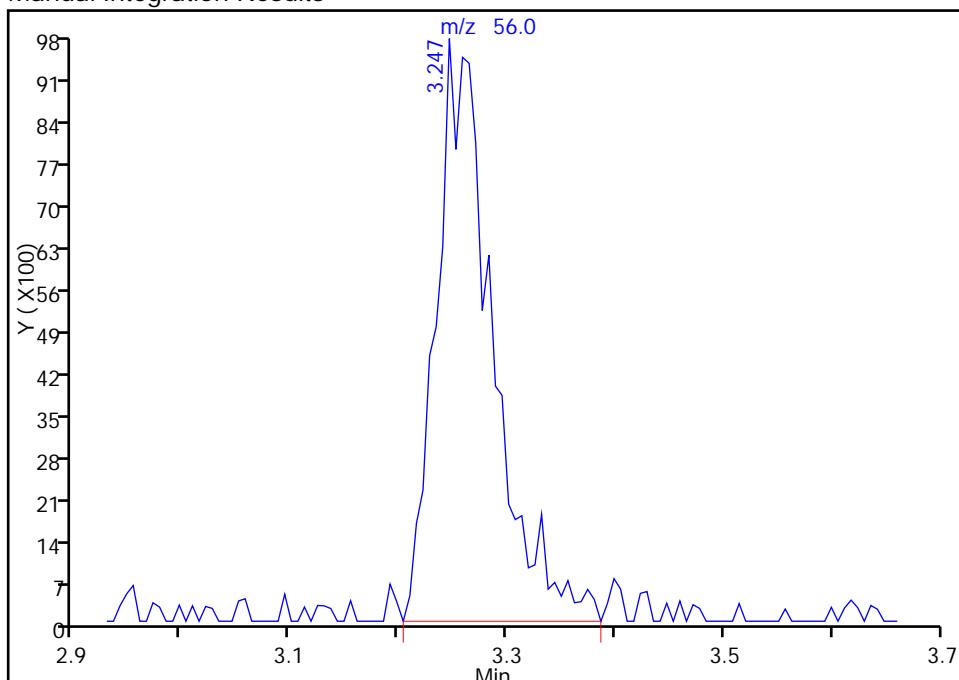
RT: 3.25  
 Area: 33235  
 Amount: 92.071591  
 Amount Units: ng

## Processing Integration Results



RT: 3.25  
 Area: 35289  
 Amount: 97.689446  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 17-Mar-2015 10:01:36

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

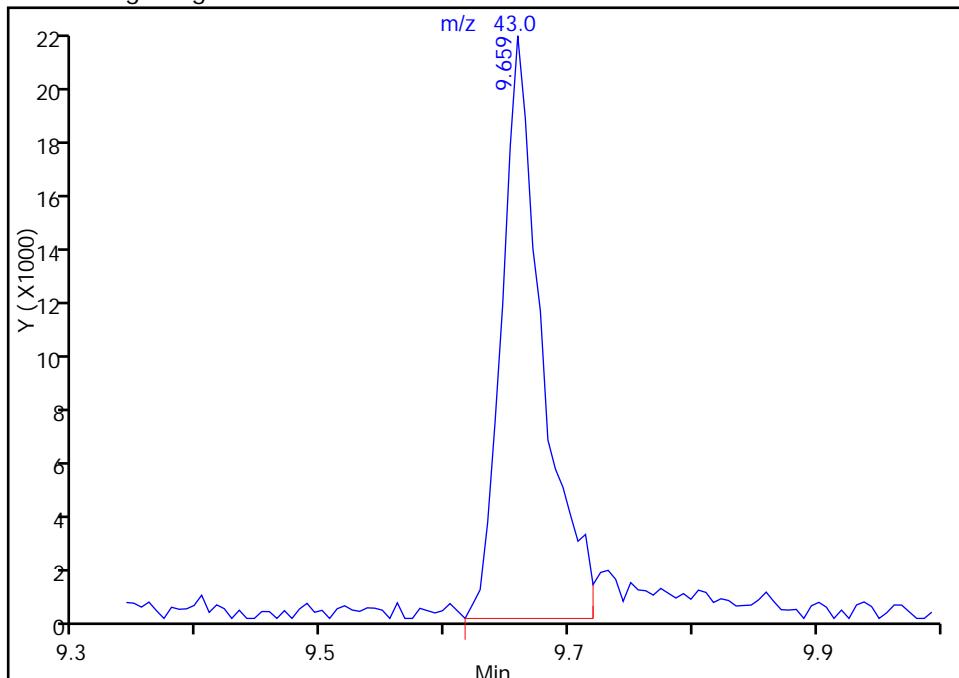
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316013.D  
 Injection Date: 16-Mar-2015 16:17:30 Instrument ID: CHHP5  
 Lims ID: IC VSTD1  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 13 Worklist Smp#: 13  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 82 2-Hexanone, CAS: 591-78-6

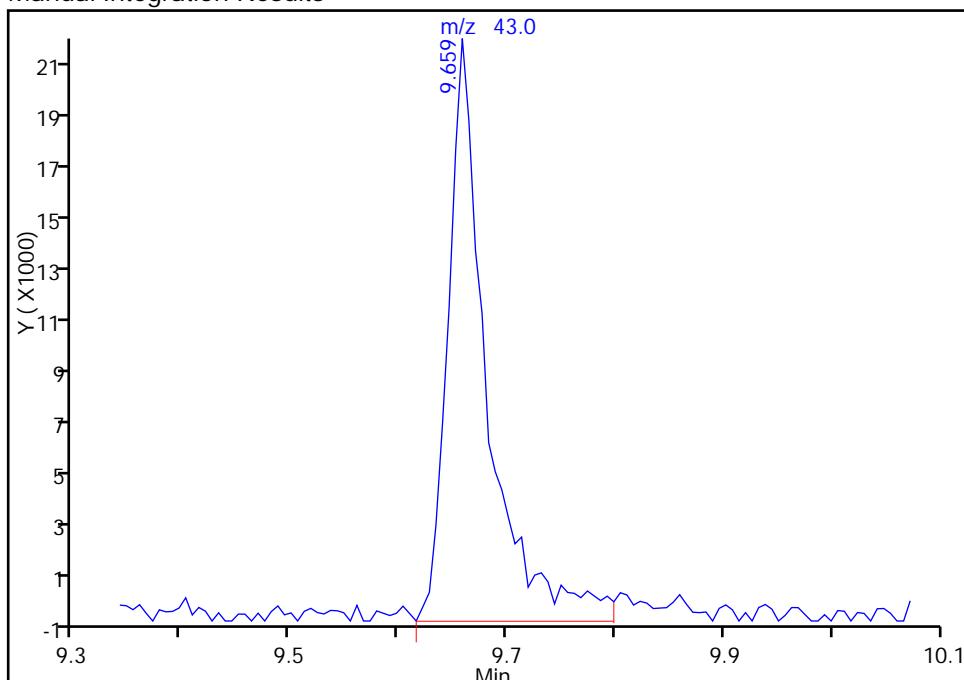
RT: 9.66  
 Area: 48498  
 Amount: 19.235523  
 Amount Units: ng

## Processing Integration Results



RT: 9.66  
 Area: 53734  
 Amount: 21.434406  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 17-Mar-2015 10:01:36

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

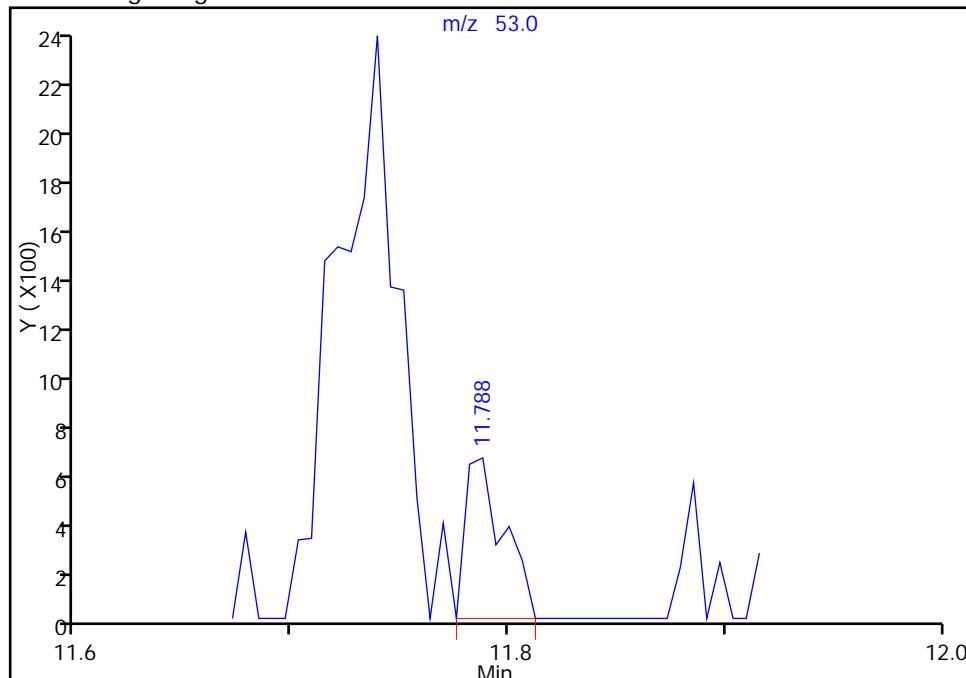
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316013.D  
 Injection Date: 16-Mar-2015 16:17:30 Instrument ID: CHHP5  
 Lims ID: IC VSTD1  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 13 Worklist Smp#: 13  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 102 trans-1,4-Dichloro-2-butene, CAS: 110-57-6

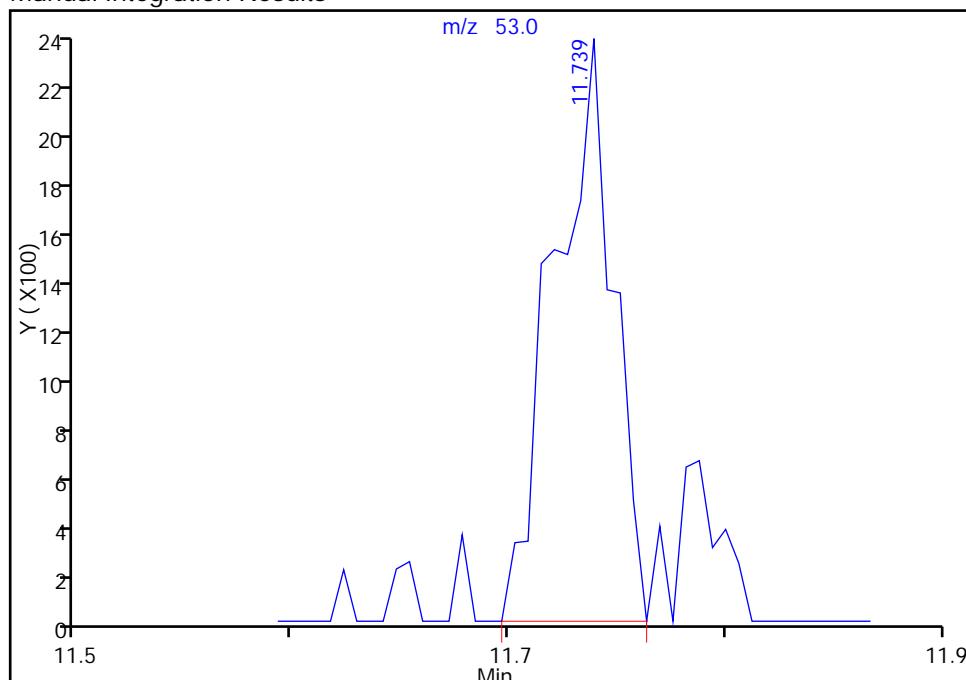
RT: 11.79  
 Area: 798  
 Amount: 0.892929  
 Amount Units: ng

## Processing Integration Results



RT: 11.74  
 Area: 4503  
 Amount: 5.086353  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 17-Mar-2015 10:01:36

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

FORM VII  
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Lab Sample ID: CCVIS 180-138583/2 Calibration Date: 04/15/2015 13:21

Instrument ID: CHHP5 Calib Start Date: 03/16/2015 12:41

GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 03/16/2015 16:17

Lab File ID: 50415002.D Conc. Units: ug/L Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Dichlorodifluoromethane	Ave	0.2143	0.2518	0.1000	11.7	10.0	17.5	20.0
Chloromethane	Ave	0.2958	0.2798	0.1000	9.46	10.0	-5.4	20.0
Vinyl chloride	Ave	0.3306	0.3289	0.1000	9.95	10.0	-0.5	20.0
Bromomethane	Lin2		0.1860	0.0500	10.4	10.0	4.3	20.0
Chloroethane	Ave	0.2287	0.2485	0.0500	10.9	10.0	8.6	20.0
Dichlorofluoromethane	Ave	0.5222	0.5361	0.0100	10.3	10.0	2.7	20.0
Trichlorofluoromethane	Ave	0.3966	0.3602	0.1000	9.08	10.0	-9.2	20.0
Ethyl ether	Ave	0.2615	0.2536	0.0100	9.69	10.0	-3.1	20.0
Acrolein	Ave	0.0318	0.0239	0.0100	22.6	30.0	-24.8*	20.0
1,1-Dichloroethene	Ave	0.2883	0.2670	0.1000	9.26	10.0	-7.4	20.0
1,1,2-Trichloro-1,2,2-trifluoroethane	Ave	0.2916	0.2998	0.1000	10.3	10.0	2.8	20.0
Acetone	Ave	0.1024	0.0962	0.0500	18.8	20.0	-6.1	20.0
Iodomethane	Ave	0.4005	0.3627	0.0100	9.06	10.0	-9.4	20.0
Carbon disulfide	Ave	0.7051	0.4663	0.1000	6.61	10.0	-33.9*	20.0
Allyl chloride	Ave	0.1524	0.1300	0.0100	8.53	10.0	-14.7	20.0
Methyl acetate	Ave	0.2396	0.2183	0.1000	45.5	50.0	-8.9	20.0
Methylene Chloride	Ave	0.3335	0.3112	0.1000	9.33	10.0	-6.7	20.0
tert-Butyl alcohol	Ave	1.178	1.091	0.0100	92.6	100	-7.4	20.0
Acrylonitrile	Ave	0.1233	0.1146	0.0100	93.0	100	-7.0	20.0
trans-1,2-Dichloroethene	Ave	0.2982	0.2866	0.1000	9.61	10.0	-3.9	20.0
Methyl tert-butyl ether	Ave	0.6593	0.5876	0.1000	8.91	10.0	-10.9	20.0
Hexane	Ave	0.4764	0.4043	0.0100	8.49	10.0	-15.1	20.0
1,1-Dichloroethane	Ave	0.5323	0.4988	0.2000	9.37	10.0	-6.3	20.0
Vinyl acetate	Ave	0.3776	0.3461	0.0100	9.17	10.0	-8.3	20.0
2,2-Dichloropropane	Ave	0.1331	0.1578	0.0100	11.9	10.0	18.6	20.0
cis-1,2-Dichloroethene	Ave	0.3142	0.2817	0.1000	8.97	10.0	-10.3	20.0
2-Butanone (MEK)	Ave	0.1638	0.1226	0.0500	15.0	20.0	-25.1*	20.0
Bromochloromethane	Ave	0.1360	0.1258	0.0100	9.26	10.0	-7.4	20.0
Tetrahydrofuran	Ave	0.1026	0.0790	0.0100	15.4	20.0	-23.0*	20.0
Chloroform	Ave	0.4836	0.4718	0.2000	9.76	10.0	-2.4	20.0
1,1,1-Trichloroethane	Ave	0.3088	0.3234	0.1000	10.5	10.0	4.7	20.0
Cyclohexane	Ave	0.5929	0.5002	0.1000	8.44	10.0	-15.6	20.0
1,1-Dichloropropene	Ave	0.4011	0.3568	0.0100	8.90	10.0	-11.0	20.0
Carbon tetrachloride	Ave	0.2478	0.2686	0.1000	10.8	10.0	8.4	20.0
Isobutyl alcohol	Ave	0.0067	0.0064*	0.0100	239	250	-4.2	20.0
Benzene	Ave	1.185	1.173	0.5000	9.90	10.0	-1.0	20.0
1,2-Dichloroethane	Ave	0.3880	0.3759	0.1000	9.69	10.0	-3.1	20.0
n-Heptane	Ave	0.4071	0.3444	0.0100	8.46	10.0	-15.4	20.0
Trichloroethene	Ave	0.2969	0.2729	0.2000	9.19	10.0	-8.1	20.0
Methylcyclohexane	Ave	0.5297	0.4520	0.1000	8.53	10.0	-14.7	20.0

FORM VII  
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Lab Sample ID: CCVIS 180-138583/2 Calibration Date: 04/15/2015 13:21

Instrument ID: CHHP5 Calib Start Date: 03/16/2015 12:41

GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 03/16/2015 16:17

Lab File ID: 50415002.D Conc. Units: ug/L Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,2-Dichloropropane	Ave	0.2931	0.2824	0.1000	9.64	10.0	-3.6	20.0
Dibromomethane	Ave	0.1578	0.1504	0.0100	9.53	10.0	-4.7	20.0
1,4-Dioxane	Ave	0.0031	0.0025*	0.0100	160	200	-19.8	20.0
Bromodichloromethane	Ave	0.3220	0.2962	0.2000	9.20	10.0	-8.0	20.0
cis-1,3-Dichloropropene	Ave	0.3107	0.3007	0.2000	9.68	10.0	-3.2	20.0
4-Methyl-2-pentanone (MIBK)	Ave	1.353	1.116	0.1000	16.5	20.0	-17.5	20.0
Toluene	Ave	5.124	5.448	0.4000	10.6	10.0	6.3	20.0
trans-1,3-Dichloropropene	Ave	0.9254	1.030	0.1000	11.1	10.0	11.3	20.0
Ethyl methacrylate	Ave	1.207	1.133	0.0100	9.39	10.0	-6.1	20.0
1,1,2-Trichloroethane	Ave	0.9609	1.013	0.1000	10.5	10.0	5.4	20.0
Tetrachloroethylene	Ave	1.002	1.050	0.2000	10.5	10.0	4.8	20.0
1,3-Dichloropropane	Ave	1.786	1.850	0.0100	10.4	10.0	3.6	20.0
2-Hexanone	Ave	1.034	0.8554	0.1000	16.5	20.0	-17.3	20.0
Dibromochloromethane	Ave	0.7670	0.7638	0.1000	9.96	10.0	-0.4	20.0
1,2-Dibromoethane (EDB)	Ave	0.9169	0.9378	0.1000	10.2	10.0	2.3	20.0
3-Chlorobenzotrifluoride	Ave	1.955	1.939	0.0100	9.92	10.0	-0.8	20.0
Chlorobenzene	Ave	3.246	3.291	0.5000	10.1	10.0	1.4	20.0
4-Chlorobenzotrifluoride	Ave	1.890	1.894	0.0100	10.0	10.0	0.2	20.0
1,1,1,2-Tetrachloroethane	Ave	0.8382	0.9457	0.0100	11.3	10.0	12.8	20.0
Ethylbenzene	Ave	1.863	1.848	0.1000	9.92	10.0	-0.8	20.0
m-Xylene & p-Xylene	Ave	2.278	2.231	0.1000	9.80	10.0	-2.0	20.0
o-Xylene	Ave	2.228	2.059	0.3000	9.24	10.0	-7.6	20.0
Styrene	Ave	3.591	3.484	0.3000	9.70	10.0	-3.0	20.0
Bromoform	Ave	0.4737	0.4576	0.1000	9.66	10.0	-3.4	20.0
2-Chlorobenzotrifluoride	Ave	1.952	1.890	0.0100	9.68	10.0	-3.2	20.0
Isopropylbenzene	Ave	5.560	5.189	0.1000	9.33	10.0	-6.7	20.0
1,1,2,2-Tetrachloroethane	Ave	1.378	1.373	0.3000	9.96	10.0	-0.4	20.0
Bromobenzene	Ave	0.9254	0.8607	0.0100	9.30	10.0	-7.0	20.0
1,2,3-Trichloropropane	Ave	0.3041	0.2961	0.0100	9.74	10.0	-2.6	20.0
trans-1,4-Dichloro-2-butene	Ave	0.2528	0.2371	0.0100	9.38	10.0	-6.2	20.0
N-Propylbenzene	Ave	1.142	1.020	0.0100	8.93	10.0	-10.7	20.0
2-Chlorotoluene	Ave	0.9591	0.8515	0.0100	8.88	10.0	-11.2	20.0
3-Chlorotoluene	Ave	1.072	1.077	0.0100	10.0	10.0	0.5	20.0
1,3,5-Trimethylbenzene	Ave	3.183	2.949	0.0100	9.27	10.0	-7.3	20.0
4-Chlorotoluene	Ave	1.038	0.9342	0.0100	9.00	10.0	-10.0	20.0
tert-Butylbenzene	Ave	2.758	2.342	0.0100	8.49	10.0	-15.1	20.0
1,2,4-Trimethylbenzene	Ave	3.267	2.935	0.0100	8.98	10.0	-10.2	20.0
3,4-Dichlorobenzotrifluoride	Ave	1.032	0.9283	0.0100	8.99	10.0	-10.1	20.0
sec-Butylbenzene	Ave	3.881	3.431	0.0100	8.84	10.0	-11.6	20.0
1,3-Dichlorobenzene	Ave	1.705	1.556	0.6000	9.13	10.0	-8.7	20.0
4-Isopropyltoluene	Ave	3.204	2.813	0.0100	8.78	10.0	-12.2	20.0

FORM VII  
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Lab Sample ID: CCVIS 180-138583/2 Calibration Date: 04/15/2015 13:21

Instrument ID: CHHP5 Calib Start Date: 03/16/2015 12:41

GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 03/16/2015 16:17

Lab File ID: 50415002.D Conc. Units: ug/L Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,4-Dichlorobenzene	Ave	1.741	1.627	0.5000	9.35	10.0	-6.5	20.0
2,4-Dichlorobenzotrifluoride	Ave	0.9669	0.8754	0.0100	9.05	10.0	-9.5	20.0
2,5-Dichlorobenzotrifluoride	Ave	1.082	0.9433	0.0100	8.72	10.0	-12.8	20.0
n-Butylbenzene	Ave	2.918	2.450	0.0100	8.40	10.0	-16.0	20.0
1,2-Dichlorobenzene	Ave	1.579	1.474	0.4000	9.34	10.0	-6.6	20.0
1,2-Dibromo-3-Chloropropane	Ave	0.1292	0.1170	0.0500	9.06	10.0	-9.4	20.0
2,4- & 2,5- & 2,6- Dichlorotoluene	Ave	1.194	0.9314	0.0100	23.4	30.0	-22.0*	20.0
2,3- & 3,4- Dichlorotoluene	Ave	1.161	0.8311	0.0100	14.3	20.0	-28.4*	20.0
1,2,4-Trichlorobenzene	Ave	0.8219	0.6160	0.2000	7.50	10.0	-25.0*	20.0
Hexachlorobutadiene	Ave	0.3941	0.3301	0.0100	8.37	10.0	-16.3	20.0
Naphthalene	Ave	2.158	1.333	0.0100	6.17	10.0	-38.3*	20.0
1,2,3-Trichlorobenzene	Ave	0.6740	0.4722	0.0100	7.01	10.0	-29.9*	20.0
2,4,5-Trichlorotoluene	Ave	0.3624	0.1610	0.0100	4.44	10.0	-55.6*	20.0
2,3,6-Trichlorotoluene	Ave	0.3273	0.1582	0.0100	4.83	10.0	-51.7*	20.0
Dibromofluoromethane (Surr)	Ave	0.2274	0.2125		9.35	10.0	-6.5	20.0
1,2-Dichloroethane-d4 (Surr)	Ave	0.2998	0.2832		9.44	10.0	-5.6	20.0
Toluene-d8 (Surr)	Ave	3.986	4.120		10.3	10.0	3.4	20.0
4-Bromofluorobenzene (Surr)	Ave	1.436	1.376		9.58	10.0	-4.2	20.0

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415002.D  
 Lims ID: CCVIS  
 Client ID:  
 Sample Type: CCVIS  
 Inject. Date: 15-Apr-2015 13:21:30 ALS Bottle#: 2 Worklist Smp#: 2  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: CCVIS  
 Misc. Info.: 180-0006480-002  
 Operator ID: 001562 Instrument ID: CHHP5  
 Sublist: chrom-MSVOA\_LL\_CHHP5\*sub11  
 Method: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 07:27:54 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICal File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 15-Apr-2015 13:44:51

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.314	4.314	0.000	0	144875	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.270	7.270	0.000	98	546873	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.361	10.361	0.000	89	119760	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.678	12.678	0.000	97	175806	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.528	6.528	0.000	80	116226	50.0	46.7	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.899	6.899	0.000	0	154857	50.0	47.2	
\$ 7 Toluene-d8 (Surr)	98	8.919	8.919	0.000	94	493409	50.0	51.7	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.529	11.529	0.000	88	164750	50.0	47.9	
11 Dichlorodifluoromethane	85	1.619	1.619	0.000	98	137708	50.0	58.7	
12 Chloromethane	50	1.789	1.789	0.000	99	153024	50.0	47.3	
13 Vinyl chloride	62	1.917	1.917	0.000	97	179867	50.0	49.7	
14 Butadiene	39	1.959	1.959	0.000	96	198236	50.0	48.0	
15 Bromomethane	94	2.264	2.264	0.000	91	101714	50.0	52.1	
16 Chloroethane	64	2.416	2.416	0.000	99	135875	50.0	54.3	
17 Dichlorofluoromethane	67	2.671	2.671	0.000	98	293175	50.0	51.3	
18 Trichlorofluoromethane	101	2.726	2.726	0.000	93	196991	50.0	45.4	
20 Ethyl ether	59	3.091	3.091	0.000	94	138662	50.0	48.5	
21 Acrolein	56	3.261	3.261	0.000	97	39180	150.0	112.8	
22 1,1-Dichloroethene	96	3.395	3.395	0.000	88	146016	50.0	46.3	
23 1,1,2-Trichloro-1,2,2-trif	101	3.450	3.450	0.000	95	163962	50.0	51.4	
24 Acetone	43	3.499	3.499	0.000	97	105244	100.0	93.9	
25 Iodomethane	142	3.626	3.626	0.000	100	198350	50.0	45.3	
26 Carbon disulfide	76	3.675	3.675	0.000	100	254978	50.0	33.1	
28 3-Chloro-1-propene	76	3.943	3.943	0.000	88	71081	50.0	42.6	
30 Methyl acetate	43	4.022	4.022	0.000	98	596762	250.0	227.7	
31 Methylene Chloride	84	4.143	4.143	0.000	98	170172	50.0	46.7	
32 2-Methyl-2-propanol	59	4.441	4.441	0.000	82	79043	500.0	463.2	
33 Acrylonitrile	53	4.551	4.551	0.000	99	626664	500.0	464.8	
34 trans-1,2-Dichloroethene	96	4.563	4.563	0.000	61	156721	50.0	48.0	
35 Methyl tert-butyl ether	73	4.600	4.600	0.000	95	321366	50.0	44.6	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
36 Hexane	57	4.983	4.983	0.000	95	221113	50.0	42.4	
37 1,1-Dichloroethane	63	5.165	5.165	0.000	96	272794	50.0	46.9	
38 Vinyl acetate	43	5.299	5.299	0.000	98	189250	50.0	45.8	
44 2,2-Dichloropropane	77	5.926	5.926	0.000	60	86294	50.0	59.3	
45 cis-1,2-Dichloroethene	96	5.938	5.938	0.000	85	154072	50.0	44.8	
46 2-Butanone (MEK)	43	5.987	5.987	0.000	100	134136	100.0	74.9	
49 Chlorobromomethane	128	6.224	6.224	0.000	97	68813	50.0	46.3	
51 Tetrahydrofuran	42	6.279	6.279	0.000	89	86407	100.0	77.0	
52 Chloroform	83	6.339	6.339	0.000	94	257998	50.0	48.8	
53 1,1,1-Trichloroethane	97	6.528	6.528	0.000	96	176859	50.0	52.4	
54 Cyclohexane	56	6.583	6.583	0.000	91	273529	50.0	42.2	
56 Carbon tetrachloride	117	6.723	6.723	0.000	73	146900	50.0	54.2	
55 1,1-Dichloropropene	75	6.723	6.723	0.000	92	195113	50.0	44.5	
57 Isobutyl alcohol	41	6.942	6.942	0.000	44	87391	1250.0	1197.2	
58 Benzene	78	6.954	6.954	0.000	97	641716	50.0	49.5	
59 1,2-Dichloroethane	62	6.984	6.984	0.000	97	205591	50.0	48.4	
62 n-Heptane	43	7.276	7.276	0.000	89	188333	50.0	42.3	
64 Trichloroethene	130	7.666	7.666	0.000	98	149241	50.0	46.0	
66 Methylcyclohexane	83	7.860	7.860	0.000	91	247182	50.0	42.7	
67 1,2-Dichloropropane	63	7.897	7.897	0.000	93	154456	50.0	48.2	
68 Dibromomethane	93	8.025	8.025	0.000	95	82243	50.0	47.7	
70 1,4-Dioxane	88	8.049	8.049	0.000	97	27073	1000.0	802.1	M
71 Dichlorobromomethane	83	8.195	8.195	0.000	98	161964	50.0	46.0	
73 2-Chloroethyl vinyl ether	63	8.517	8.517	0.000	93	165677	100.0	91.7	
74 cis-1,3-Dichloropropene	75	8.651	8.651	0.000	92	164416	50.0	48.4	
75 4-Methyl-2-pentanone (MIBK)	43	8.822	8.822	0.000	97	267388	100.0	82.5	
76 Toluene	91	8.986	8.986	0.000	98	652396	50.0	53.2	
77 trans-1,3-Dichloropropene	75	9.217	9.217	0.000	97	123387	50.0	55.7	
78 Ethyl methacrylate	69	9.314	9.314	0.000	90	135736	50.0	46.9	
79 1,1,2-Trichloroethane	97	9.399	9.399	0.000	92	121282	50.0	52.7	
80 Tetrachloroethene	164	9.533	9.533	0.000	97	125750	50.0	52.4	
81 1,3-Dichloropropane	76	9.558	9.558	0.000	95	221501	50.0	51.8	
82 2-Hexanone	43	9.655	9.655	0.000	96	204885	100.0	82.7	
84 Chlorodibromomethane	129	9.789	9.789	0.000	90	91471	50.0	49.8	
85 Ethylene Dibromide	107	9.898	9.898	0.000	98	112306	50.0	51.1	
86 3-Chlorobenzotrifluoride	180	10.373	10.373	0.000	95	232172	50.0	49.6	
87 Chlorobenzene	112	10.391	10.391	0.000	93	394179	50.0	50.7	
88 4-Chlorobenzotrifluoride	180	10.428	10.428	0.000	94	226858	50.0	50.1	
89 1,1,1,2-Tetrachloroethane	131	10.470	10.470	0.000	92	113256	50.0	56.4	
90 Ethylbenzene	106	10.501	10.501	0.000	99	221268	50.0	49.6	
91 m-Xylene & p-Xylene	106	10.616	10.616	0.000	0	267242	50.0	49.0	
92 o-Xylene	106	11.012	11.012	0.000	97	246631	50.0	46.2	
93 Styrene	104	11.024	11.024	0.000	96	417299	50.0	48.5	
94 Bromoform	173	11.212	11.212	0.000	97	54802	50.0	48.3	
96 2-Chlorobenzotrifluoride	180	11.273	11.273	0.000	97	226296	50.0	48.4	
97 Isopropylbenzene	105	11.377	11.377	0.000	97	621433	50.0	46.7	
99 1,1,2,2-Tetrachloroethane	83	11.675	11.675	0.000	95	164434	50.0	49.8	
100 Bromobenzene	156	11.681	11.681	0.000	95	151309	50.0	46.5	
101 1,2,3-Trichloropropane	110	11.717	11.717	0.000	85	52052	50.0	48.7	
102 trans-1,4-Dichloro-2-buten	53	11.729	11.729	0.000	70	41676	50.0	46.9	
103 N-Propylbenzene	120	11.784	11.784	0.000	99	179242	50.0	44.7	
104 2-Chlorotoluene	126	11.869	11.869	0.000	96	149703	50.0	44.4	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
105 3-Chlorotoluene	126	11.930	11.930	0.000	95	189348	50.0	50.2	
106 1,3,5-Trimethylbenzene	105	11.961	11.961	0.000	94	518496	50.0	46.3	
107 4-Chlorotoluene	126	11.985	11.985	0.000	98	164244	50.0	45.0	
108 tert-Butylbenzene	119	12.283	12.283	0.000	94	411758	50.0	42.5	
110 1,2,4-Trimethylbenzene	105	12.332	12.332	0.000	95	516034	50.0	44.9	
111 1,2-dichloro-4-(trifluoromethyl)	214	12.399	12.399	0.000	98	163198	50.0	45.0	
112 sec-Butylbenzene	105	12.502	12.502	0.000	95	603202	50.0	44.2	
113 1,3-Dichlorobenzene	146	12.618	12.618	0.000	98	273567	50.0	45.6	
114 4-Isopropyltoluene	119	12.648	12.648	0.000	96	494548	50.0	43.9	
115 1,4-Dichlorobenzene	146	12.703	12.703	0.000	93	286091	50.0	46.7	
116 2,4-Dichloro-1-(trifluoromethyl)	214	12.758	12.758	0.000	98	153894	50.0	45.3	
118 2,5-Dichlorobenzotrifluoride	214	12.806	12.806	0.000	0	165845	50.0	43.6	
120 n-Butylbenzene	91	13.062	13.062	0.000	98	430684	50.0	42.0	
121 1,2-Dichlorobenzene	146	13.080	13.080	0.000	96	259203	50.0	46.7	
122 1,2-Dibromo-3-Chloropropan	75	13.853	13.853	0.000	76	20568	50.0	45.3	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125	14.005	14.005	0.000	0	491254	150.0	117.1	
124 1,3,5-Trichlorobenzene	180	14.066	14.066	0.000	97	150831	50.0	44.8	
125 2,3- & 3,4- Dichlorotoluene	125	14.424	14.424	0.000	0	292209	100.0	71.6	
126 1,2,4-Trichlorobenzene	180	14.686	14.686	0.000	93	108295	50.0	37.5	
127 Hexachlorobutadiene	225	14.856	14.856	0.000	95	58029	50.0	41.9	
128 Naphthalene	128	14.942	14.942	0.000	97	234261	50.0	30.9	
129 1,2,3-Trichlorobenzene	180	15.185	15.185	0.000	93	83023	50.0	35.0	
131 2,4,5-Trichlorotoluene	159	15.964	15.964	0.000	0	28298	50.0	22.2	
130 2,3,6-Trichlorotoluene	159	16.061	16.061	0.000	94	27811	50.0	24.2	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
S 133 Xylenes, Total	106				0		100.0	95.2	
S 134 1,2-Dichloroethene, Total	96				0		100.0	92.9	
S 135 1,3-Dichloropropene, Total	1				0		100.0	104.0	

**QC Flag Legend**

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

**Reagents:**

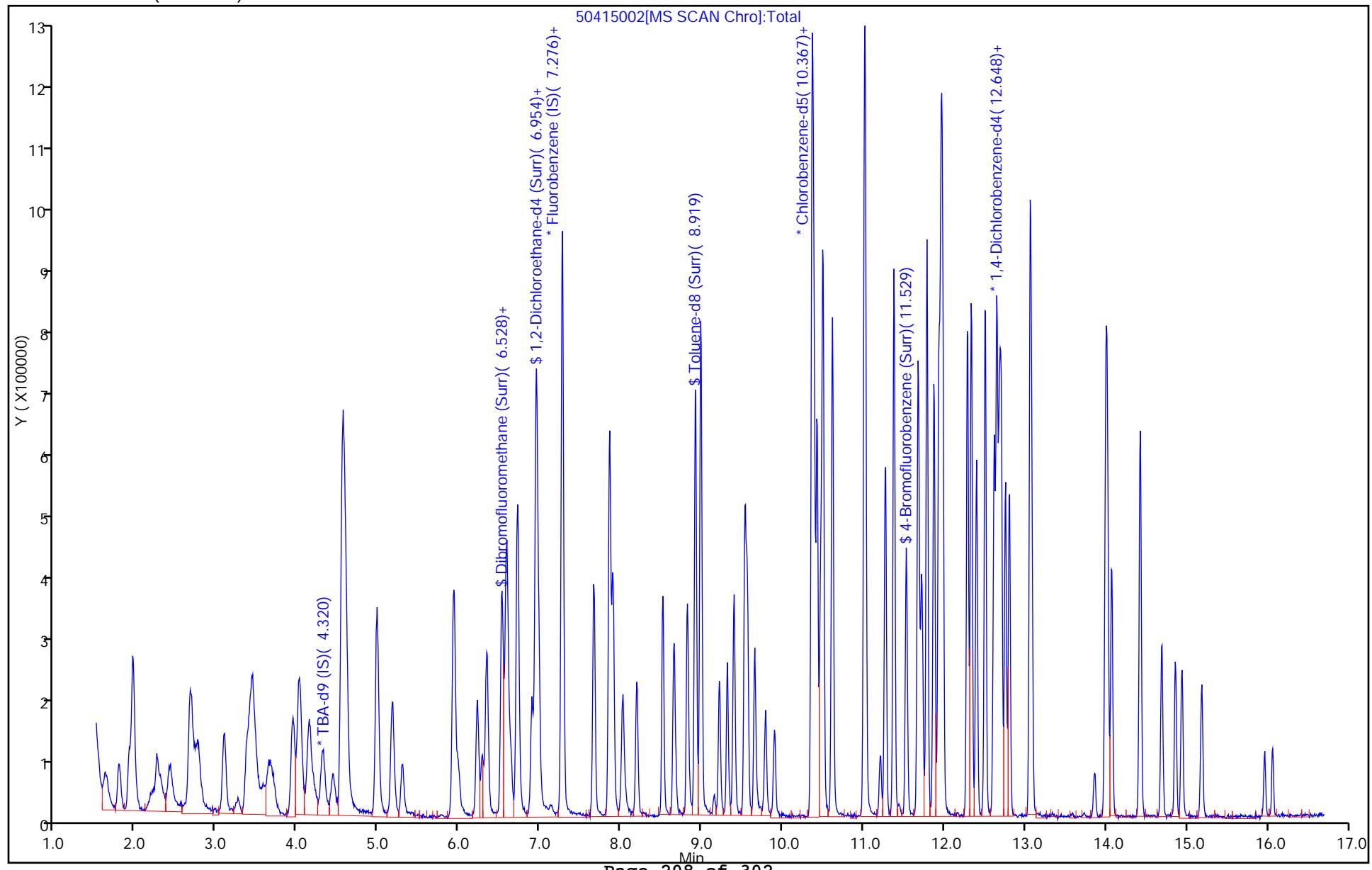
voaW 135tcb a_00001	Amount Added: 2.00	Units: uL	
VOA8260VOAPRI_00110	Amount Added: 2.00	Units: uL	
voaWKetpri Re_00004	Amount Added: 2.00	Units: uL	
voaWeemixPRI_00002	Amount Added: 2.00	Units: uL	
voaW VA pri R_00005	Amount Added: 2.00	Units: uL	
voaW2-cl 2ndR_00002	Amount Added: 2.00	Units: uL	
VOAACRPRI_00005	Amount Added: 6.00	Units: uL	
VOA8260INT_00031	Amount Added: 2.00	Units: uL	Run Reagent
VOA8260SURR_00033	Amount Added: 2.00	Units: uL	Run Reagent

Report Date: 16-Apr-2015 07:27:56

Chrom Revision: 2.2 13-Mar-2015 11:20:44

## TestAmerica Pittsburgh

Data File: \PITCHROM\ChromData\CHHP5\20150415-6480.b\50415002.D  
Injection Date: 15-Apr-2015 13:21:30 Instrument ID: CHHP5  
Lims ID: CCVIS Operator ID: 001562  
Client ID:  
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 2  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)



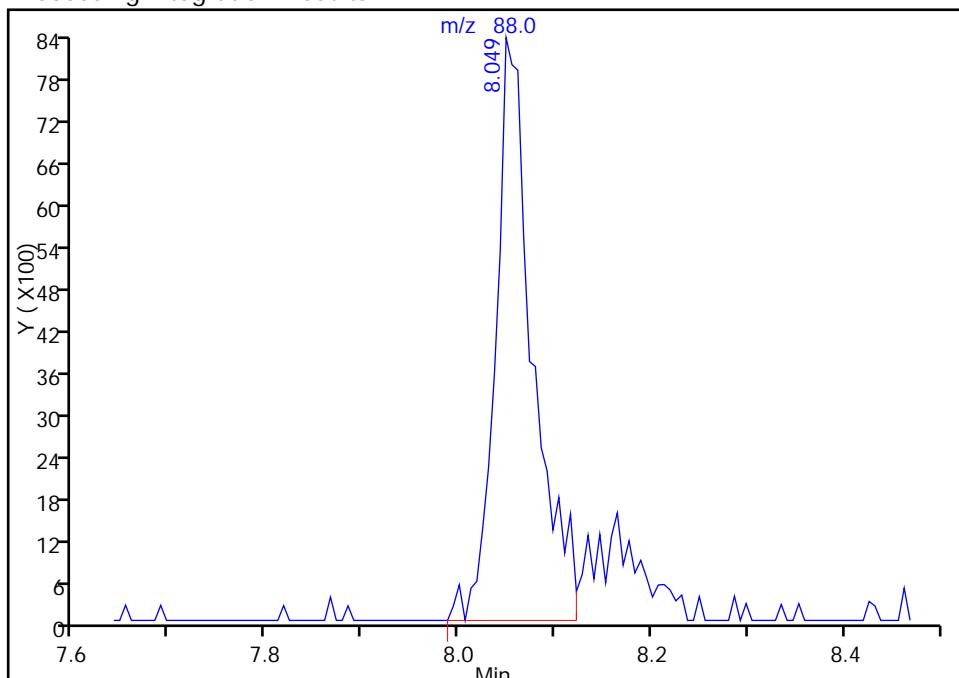
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415002.D  
 Injection Date: 15-Apr-2015 13:21:30 Instrument ID: CHHP5  
 Lims ID: CCVIS  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 2 Worklist Smp#: 2  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 70 1,4-Dioxane, CAS: 123-91-1

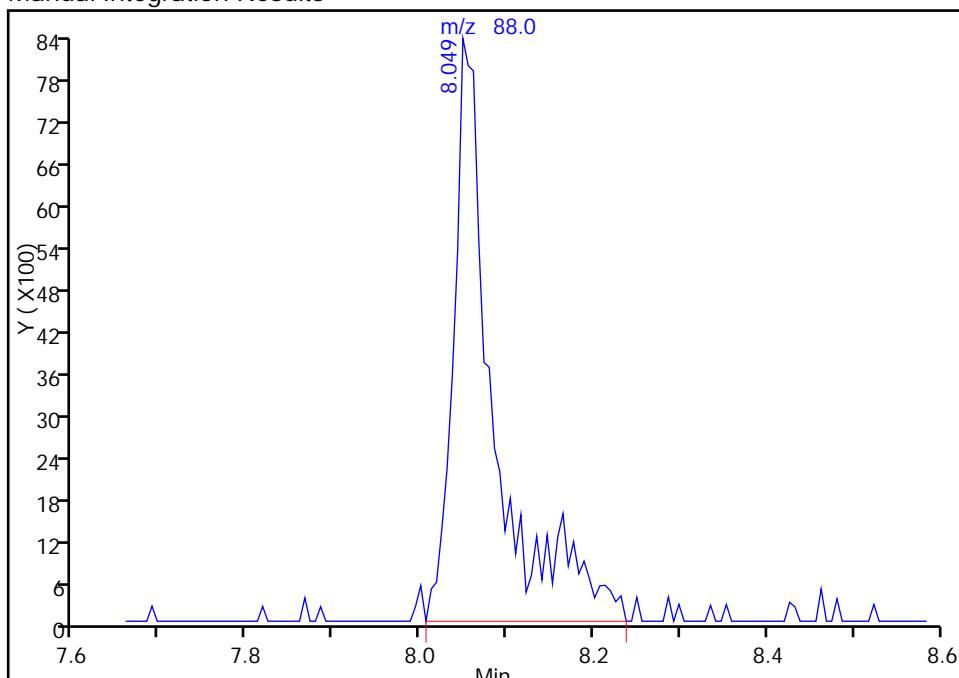
RT: 8.05  
 Area: 22422  
 Amount: 664.3240  
 Amount Units: ng

## Processing Integration Results



RT: 8.05  
 Area: 27073  
 Amount: 802.1248  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 15-Apr-2015 13:44:51

Audit Action: Manually Integrated

Audit Reason: Peak Tail

FORM VII  
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Lab Sample ID: CCVIS 180-138583/2 Calibration Date: 04/15/2015 13:21  
Instrument ID: CHHP5 Calib Start Date: 03/18/2015 13:31  
GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 03/18/2015 16:19  
Lab File ID: 50415002.D Conc. Units: ug/L Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
2-Chloroethyl vinyl ether	Ave	0.1652	0.1515	0.0100	18.3	20.0	-8.3	20.0
1,3,5-Trichlorobenzene	Ave	0.9577	0.8579	0.0100	8.96	10.0	-10.4	20.0

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415002.D  
 Lims ID: CCVIS  
 Client ID:  
 Sample Type: CCVIS  
 Inject. Date: 15-Apr-2015 13:21:30 ALS Bottle#: 2 Worklist Smp#: 2  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: CCVIS  
 Misc. Info.: 180-0006480-002  
 Operator ID: 001562 Instrument ID: CHHP5  
 Sublist: chrom-MSVOA\_LL\_CHHP5\*sub11  
 Method: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 07:27:54 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICal File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 15-Apr-2015 13:44:51

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.314	4.314	0.000	0	144875	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.270	7.270	0.000	98	546873	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.361	10.361	0.000	89	119760	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.678	12.678	0.000	97	175806	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.528	6.528	0.000	80	116226	50.0	46.7	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.899	6.899	0.000	0	154857	50.0	47.2	
\$ 7 Toluene-d8 (Surr)	98	8.919	8.919	0.000	94	493409	50.0	51.7	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.529	11.529	0.000	88	164750	50.0	47.9	
11 Dichlorodifluoromethane	85	1.619	1.619	0.000	98	137708	50.0	58.7	
12 Chloromethane	50	1.789	1.789	0.000	99	153024	50.0	47.3	
13 Vinyl chloride	62	1.917	1.917	0.000	97	179867	50.0	49.7	
14 Butadiene	39	1.959	1.959	0.000	96	198236	50.0	48.0	
15 Bromomethane	94	2.264	2.264	0.000	91	101714	50.0	52.1	
16 Chloroethane	64	2.416	2.416	0.000	99	135875	50.0	54.3	
17 Dichlorofluoromethane	67	2.671	2.671	0.000	98	293175	50.0	51.3	
18 Trichlorofluoromethane	101	2.726	2.726	0.000	93	196991	50.0	45.4	
20 Ethyl ether	59	3.091	3.091	0.000	94	138662	50.0	48.5	
21 Acrolein	56	3.261	3.261	0.000	97	39180	150.0	112.8	
22 1,1-Dichloroethene	96	3.395	3.395	0.000	88	146016	50.0	46.3	
23 1,1,2-Trichloro-1,2,2-trif	101	3.450	3.450	0.000	95	163962	50.0	51.4	
24 Acetone	43	3.499	3.499	0.000	97	105244	100.0	93.9	
25 Iodomethane	142	3.626	3.626	0.000	100	198350	50.0	45.3	
26 Carbon disulfide	76	3.675	3.675	0.000	100	254978	50.0	33.1	
28 3-Chloro-1-propene	76	3.943	3.943	0.000	88	71081	50.0	42.6	
30 Methyl acetate	43	4.022	4.022	0.000	98	596762	250.0	227.7	
31 Methylene Chloride	84	4.143	4.143	0.000	98	170172	50.0	46.7	
32 2-Methyl-2-propanol	59	4.441	4.441	0.000	82	79043	500.0	463.2	
33 Acrylonitrile	53	4.551	4.551	0.000	99	626664	500.0	464.8	
34 trans-1,2-Dichloroethene	96	4.563	4.563	0.000	61	156721	50.0	48.0	
35 Methyl tert-butyl ether	73	4.600	4.600	0.000	95	321366	50.0	44.6	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
36 Hexane	57	4.983	4.983	0.000	95	221113	50.0	42.4	
37 1,1-Dichloroethane	63	5.165	5.165	0.000	96	272794	50.0	46.9	
38 Vinyl acetate	43	5.299	5.299	0.000	98	189250	50.0	45.8	
44 2,2-Dichloropropane	77	5.926	5.926	0.000	60	86294	50.0	59.3	
45 cis-1,2-Dichloroethene	96	5.938	5.938	0.000	85	154072	50.0	44.8	
46 2-Butanone (MEK)	43	5.987	5.987	0.000	100	134136	100.0	74.9	
49 Chlorobromomethane	128	6.224	6.224	0.000	97	68813	50.0	46.3	
51 Tetrahydrofuran	42	6.279	6.279	0.000	89	86407	100.0	77.0	
52 Chloroform	83	6.339	6.339	0.000	94	257998	50.0	48.8	
53 1,1,1-Trichloroethane	97	6.528	6.528	0.000	96	176859	50.0	52.4	
54 Cyclohexane	56	6.583	6.583	0.000	91	273529	50.0	42.2	
56 Carbon tetrachloride	117	6.723	6.723	0.000	73	146900	50.0	54.2	
55 1,1-Dichloropropene	75	6.723	6.723	0.000	92	195113	50.0	44.5	
57 Isobutyl alcohol	41	6.942	6.942	0.000	44	87391	1250.0	1197.2	
58 Benzene	78	6.954	6.954	0.000	97	641716	50.0	49.5	
59 1,2-Dichloroethane	62	6.984	6.984	0.000	97	205591	50.0	48.4	
62 n-Heptane	43	7.276	7.276	0.000	89	188333	50.0	42.3	
64 Trichloroethene	130	7.666	7.666	0.000	98	149241	50.0	46.0	
66 Methylcyclohexane	83	7.860	7.860	0.000	91	247182	50.0	42.7	
67 1,2-Dichloropropane	63	7.897	7.897	0.000	93	154456	50.0	48.2	
68 Dibromomethane	93	8.025	8.025	0.000	95	82243	50.0	47.7	
70 1,4-Dioxane	88	8.049	8.049	0.000	97	27073	1000.0	802.1	M
71 Dichlorobromomethane	83	8.195	8.195	0.000	98	161964	50.0	46.0	
73 2-Chloroethyl vinyl ether	63	8.517	8.517	0.000	93	165677	100.0	91.7	
74 cis-1,3-Dichloropropene	75	8.651	8.651	0.000	92	164416	50.0	48.4	
75 4-Methyl-2-pentanone (MIBK)	43	8.822	8.822	0.000	97	267388	100.0	82.5	
76 Toluene	91	8.986	8.986	0.000	98	652396	50.0	53.2	
77 trans-1,3-Dichloropropene	75	9.217	9.217	0.000	97	123387	50.0	55.7	
78 Ethyl methacrylate	69	9.314	9.314	0.000	90	135736	50.0	46.9	
79 1,1,2-Trichloroethane	97	9.399	9.399	0.000	92	121282	50.0	52.7	
80 Tetrachloroethene	164	9.533	9.533	0.000	97	125750	50.0	52.4	
81 1,3-Dichloropropane	76	9.558	9.558	0.000	95	221501	50.0	51.8	
82 2-Hexanone	43	9.655	9.655	0.000	96	204885	100.0	82.7	
84 Chlorodibromomethane	129	9.789	9.789	0.000	90	91471	50.0	49.8	
85 Ethylene Dibromide	107	9.898	9.898	0.000	98	112306	50.0	51.1	
86 3-Chlorobenzotrifluoride	180	10.373	10.373	0.000	95	232172	50.0	49.6	
87 Chlorobenzene	112	10.391	10.391	0.000	93	394179	50.0	50.7	
88 4-Chlorobenzotrifluoride	180	10.428	10.428	0.000	94	226858	50.0	50.1	
89 1,1,1,2-Tetrachloroethane	131	10.470	10.470	0.000	92	113256	50.0	56.4	
90 Ethylbenzene	106	10.501	10.501	0.000	99	221268	50.0	49.6	
91 m-Xylene & p-Xylene	106	10.616	10.616	0.000	0	267242	50.0	49.0	
92 o-Xylene	106	11.012	11.012	0.000	97	246631	50.0	46.2	
93 Styrene	104	11.024	11.024	0.000	96	417299	50.0	48.5	
94 Bromoform	173	11.212	11.212	0.000	97	54802	50.0	48.3	
96 2-Chlorobenzotrifluoride	180	11.273	11.273	0.000	97	226296	50.0	48.4	
97 Isopropylbenzene	105	11.377	11.377	0.000	97	621433	50.0	46.7	
99 1,1,2,2-Tetrachloroethane	83	11.675	11.675	0.000	95	164434	50.0	49.8	
100 Bromobenzene	156	11.681	11.681	0.000	95	151309	50.0	46.5	
101 1,2,3-Trichloropropane	110	11.717	11.717	0.000	85	52052	50.0	48.7	
102 trans-1,4-Dichloro-2-buten	53	11.729	11.729	0.000	70	41676	50.0	46.9	
103 N-Propylbenzene	120	11.784	11.784	0.000	99	179242	50.0	44.7	
104 2-Chlorotoluene	126	11.869	11.869	0.000	96	149703	50.0	44.4	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
105 3-Chlorotoluene	126	11.930	11.930	0.000	95	189348	50.0	50.2	
106 1,3,5-Trimethylbenzene	105	11.961	11.961	0.000	94	518496	50.0	46.3	
107 4-Chlorotoluene	126	11.985	11.985	0.000	98	164244	50.0	45.0	
108 tert-Butylbenzene	119	12.283	12.283	0.000	94	411758	50.0	42.5	
110 1,2,4-Trimethylbenzene	105	12.332	12.332	0.000	95	516034	50.0	44.9	
111 1,2-dichloro-4-(trifluoromethyl)	214	12.399	12.399	0.000	98	163198	50.0	45.0	
112 sec-Butylbenzene	105	12.502	12.502	0.000	95	603202	50.0	44.2	
113 1,3-Dichlorobenzene	146	12.618	12.618	0.000	98	273567	50.0	45.6	
114 4-Isopropyltoluene	119	12.648	12.648	0.000	96	494548	50.0	43.9	
115 1,4-Dichlorobenzene	146	12.703	12.703	0.000	93	286091	50.0	46.7	
116 2,4-Dichloro-1-(trifluoromethyl)	214	12.758	12.758	0.000	98	153894	50.0	45.3	
118 2,5-Dichlorobenzotrifluoride	214	12.806	12.806	0.000	0	165845	50.0	43.6	
120 n-Butylbenzene	91	13.062	13.062	0.000	98	430684	50.0	42.0	
121 1,2-Dichlorobenzene	146	13.080	13.080	0.000	96	259203	50.0	46.7	
122 1,2-Dibromo-3-Chloropropan	75	13.853	13.853	0.000	76	20568	50.0	45.3	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125	14.005	14.005	0.000	0	491254	150.0	117.1	
124 1,3,5-Trichlorobenzene	180	14.066	14.066	0.000	97	150831	50.0	44.8	
125 2,3- & 3,4- Dichlorotoluene	125	14.424	14.424	0.000	0	292209	100.0	71.6	
126 1,2,4-Trichlorobenzene	180	14.686	14.686	0.000	93	108295	50.0	37.5	
127 Hexachlorobutadiene	225	14.856	14.856	0.000	95	58029	50.0	41.9	
128 Naphthalene	128	14.942	14.942	0.000	97	234261	50.0	30.9	
129 1,2,3-Trichlorobenzene	180	15.185	15.185	0.000	93	83023	50.0	35.0	
131 2,4,5-Trichlorotoluene	159	15.964	15.964	0.000	0	28298	50.0	22.2	
130 2,3,6-Trichlorotoluene	159	16.061	16.061	0.000	94	27811	50.0	24.2	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
S 133 Xylenes, Total	106				0		100.0	95.2	
S 134 1,2-Dichloroethene, Total	96				0		100.0	92.9	
S 135 1,3-Dichloropropene, Total	1				0		100.0	104.0	

**QC Flag Legend**

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

**Reagents:**

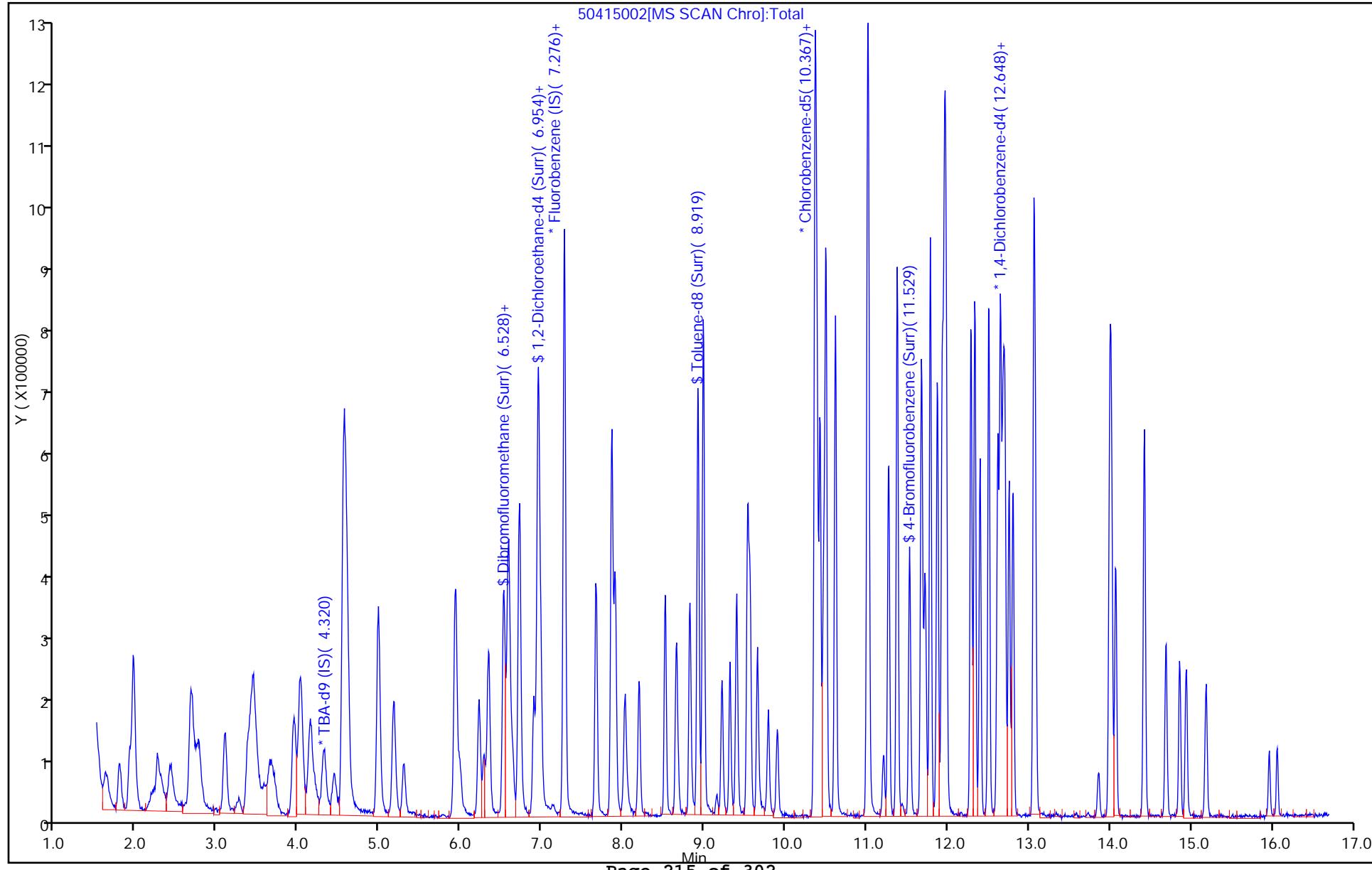
voaW 135tcb a_00001	Amount Added: 2.00	Units: uL	
VOA8260VOAPRI_00110	Amount Added: 2.00	Units: uL	
voaWKetpri Re_00004	Amount Added: 2.00	Units: uL	
voaWeemixPRI_00002	Amount Added: 2.00	Units: uL	
voaW VA pri R_00005	Amount Added: 2.00	Units: uL	
voaW2-cl 2ndR_00002	Amount Added: 2.00	Units: uL	
VOAACRPRI_00005	Amount Added: 6.00	Units: uL	
VOA8260INT_00031	Amount Added: 2.00	Units: uL	Run Reagent
VOA8260SURR_00033	Amount Added: 2.00	Units: uL	Run Reagent

Report Date: 16-Apr-2015 07:27:56

Chrom Revision: 2.2 13-Mar-2015 11:20:44

## TestAmerica Pittsburgh

Data File: \PITCHROM\ChromData\CHHP5\20150415-6480.b\50415002.D  
Injection Date: 15-Apr-2015 13:21:30 Instrument ID: CHHP5  
Lims ID: CCVIS Operator ID: 001562  
Client ID:  
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 2  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)



FORM VII  
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Lab Sample ID: CCVIS 180-138685/2 Calibration Date: 04/16/2015 10:19  
Instrument ID: CHHP5 Calib Start Date: 03/16/2015 12:41  
GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 03/16/2015 16:17  
Lab File ID: 50416002.D Conc. Units: ug/L Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Dichlorodifluoromethane	Ave	0.2143	0.2605	0.1000	12.2	10.0	21.6*	20.0
Chloromethane	Ave	0.2958	0.3144	0.1000	10.6	10.0	6.3	20.0
Vinyl chloride	Ave	0.3306	0.3604	0.1000	10.9	10.0	9.0	20.0
Bromomethane	Lin2		0.1972	0.0500	11.1	10.0	11.1	20.0
Chloroethane	Ave	0.2287	0.2541	0.0500	11.1	10.0	11.1	20.0
Dichlorofluoromethane	Ave	0.5222	0.5748	0.0100	11.0	10.0	10.1	20.0
Trichlorofluoromethane	Ave	0.3966	0.3900	0.1000	9.83	10.0	-1.7	20.0
Ethyl ether	Ave	0.2615	0.2632	0.0100	10.1	10.0	0.6	20.0
Acrolein	Ave	0.0318	0.0285	0.0100	26.9	30.0	-10.3	20.0
1,1-Dichloroethene	Ave	0.2883	0.2850	0.1000	9.88	10.0	-1.2	20.0
1,1,2-Trichloro-1,2,2-trifluoroethane	Ave	0.2916	0.3070	0.1000	10.5	10.0	5.3	20.0
Acetone	Ave	0.1024	0.1191	0.0500	23.3	20.0	16.3	20.0
Iodomethane	Ave	0.4005	0.3841	0.0100	9.59	10.0	-4.1	20.0
Carbon disulfide	Ave	0.7051	0.5468	0.1000	7.76	10.0	-22.4*	20.0
Allyl chloride	Ave	0.1524	0.1509	0.0100	9.90	10.0	-1.0	20.0
Methyl acetate	Ave	0.2396	0.2343	0.1000	48.9	50.0	-2.2	20.0
Methylene Chloride	Ave	0.3335	0.3345	0.1000	10.0	10.0	0.3	20.0
tert-Butyl alcohol	Ave	1.178	1.133	0.0100	96.2	100	-3.8	20.0
Acrylonitrile	Ave	0.1233	0.1214	0.0100	98.5	100	-1.5	20.0
trans-1,2-Dichloroethene	Ave	0.2982	0.2982	0.1000	10.0	10.0	-0.0	20.0
Methyl tert-butyl ether	Ave	0.6593	0.6333	0.1000	9.61	10.0	-3.9	20.0
Hexane	Ave	0.4764	0.4195	0.0100	8.80	10.0	-12.0	20.0
1,1-Dichloroethane	Ave	0.5323	0.5436	0.2000	10.2	10.0	2.1	20.0
Vinyl acetate	Ave	0.3776	0.3407	0.0100	9.02	10.0	-9.8	20.0
2,2-Dichloropropane	Ave	0.1331	0.1712	0.0100	12.9	10.0	28.6*	20.0
cis-1,2-Dichloroethene	Ave	0.3142	0.3021	0.1000	9.62	10.0	-3.8	20.0
2-Butanone (MEK)	Ave	0.1638	0.1479	0.0500	18.1	20.0	-9.7	20.0
Bromochloromethane	Ave	0.1360	0.1335	0.0100	9.82	10.0	-1.8	20.0
Tetrahydrofuran	Ave	0.1026	0.0860	0.0100	16.8	20.0	-16.1	20.0
Chloroform	Ave	0.4836	0.5039	0.2000	10.4	10.0	4.2	20.0
1,1,1-Trichloroethane	Ave	0.3088	0.3448	0.1000	11.2	10.0	11.7	20.0
Cyclohexane	Ave	0.5929	0.5352	0.1000	9.03	10.0	-9.7	20.0
1,1-Dichloropropene	Ave	0.4011	0.3792	0.0100	9.45	10.0	-5.5	20.0
Carbon tetrachloride	Ave	0.2478	0.2915	0.1000	11.8	10.0	17.6	20.0
Isobutyl alcohol	Ave	0.0067	0.0077*	0.0100	287	250	14.7	20.0
Benzene	Ave	1.185	1.229	0.5000	10.4	10.0	3.7	20.0
1,2-Dichloroethane	Ave	0.3880	0.3938	0.1000	10.1	10.0	1.5	20.0
n-Heptane	Ave	0.4071	0.3743	0.0100	9.19	10.0	-8.1	20.0
Trichloroethene	Ave	0.2969	0.2714	0.2000	9.14	10.0	-8.6	20.0
Methylcyclohexane	Ave	0.5297	0.4750	0.1000	8.97	10.0	-10.3	20.0

FORM VII  
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Lab Sample ID: CCVIS 180-138685/2 Calibration Date: 04/16/2015 10:19

Instrument ID: CHHP5 Calib Start Date: 03/16/2015 12:41

GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 03/16/2015 16:17

Lab File ID: 50416002.D Conc. Units: ug/L Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,2-Dichloropropane	Ave	0.2931	0.3027	0.1000	10.3	10.0	3.3	20.0
Dibromomethane	Ave	0.1578	0.1607	0.0100	10.2	10.0	1.8	20.0
1,4-Dioxane	Ave	0.0031	0.0026*	0.0100	171	200	-14.7	20.0
Bromodichloromethane	Ave	0.3220	0.3294	0.2000	10.2	10.0	2.3	20.0
cis-1,3-Dichloropropene	Ave	0.3107	0.3245	0.2000	10.4	10.0	4.4	20.0
4-Methyl-2-pentanone (MIBK)	Ave	1.353	1.131	0.1000	16.7	20.0	-16.4	20.0
Toluene	Ave	5.124	5.520	0.4000	10.8	10.0	7.7	20.0
trans-1,3-Dichloropropene	Ave	0.9254	1.078	0.1000	11.7	10.0	16.5	20.0
Ethyl methacrylate	Ave	1.207	1.111	0.0100	9.21	10.0	-7.9	20.0
1,1,2-Trichloroethane	Ave	0.9609	1.008	0.1000	10.5	10.0	4.9	20.0
Tetrachloroethylene	Ave	1.002	1.038	0.2000	10.4	10.0	3.5	20.0
1,3-Dichloropropane	Ave	1.786	1.843	0.0100	10.3	10.0	3.2	20.0
2-Hexanone	Ave	1.034	0.9355	0.1000	18.1	20.0	-9.5	20.0
Dibromochloromethane	Ave	0.7670	0.8273	0.1000	10.8	10.0	7.9	20.0
1,2-Dibromoethane (EDB)	Ave	0.9169	0.9609	0.1000	10.5	10.0	4.8	20.0
3-Chlorobenzotrifluoride	Ave	1.955	2.119	0.0100	10.8	10.0	8.4	20.0
Chlorobenzene	Ave	3.246	3.332	0.5000	10.3	10.0	2.6	20.0
4-Chlorobenzotrifluoride	Ave	1.890	1.983	0.0100	10.5	10.0	4.9	20.0
1,1,1,2-Tetrachloroethane	Ave	0.8382	0.9716	0.0100	11.6	10.0	15.9	20.0
Ethylbenzene	Ave	1.863	1.828	0.1000	9.82	10.0	-1.8	20.0
m-Xylene & p-Xylene	Ave	2.278	2.226	0.1000	9.77	10.0	-2.3	20.0
o-Xylene	Ave	2.228	2.084	0.3000	9.35	10.0	-6.5	20.0
Styrene	Ave	3.591	3.560	0.3000	9.92	10.0	-0.8	20.0
Bromoform	Ave	0.4737	0.4985	0.1000	10.5	10.0	5.2	20.0
2-Chlorobenzotrifluoride	Ave	1.952	2.000	0.0100	10.2	10.0	2.4	20.0
Isopropylbenzene	Ave	5.560	5.275	0.1000	9.49	10.0	-5.1	20.0
1,1,2,2-Tetrachloroethane	Ave	1.378	1.425	0.3000	10.3	10.0	3.4	20.0
Bromobenzene	Ave	0.9254	0.8662	0.0100	9.36	10.0	-6.4	20.0
1,2,3-Trichloropropene	Ave	0.3041	0.2800	0.0100	9.21	10.0	-7.9	20.0
trans-1,4-Dichloro-2-butene	Ave	0.2528	0.2305	0.0100	9.12	10.0	-8.8	20.0
N-Propylbenzene	Ave	1.142	1.034	0.0100	9.06	10.0	-9.4	20.0
2-Chlorotoluene	Ave	0.9591	0.8812	0.0100	9.19	10.0	-8.1	20.0
3-Chlorotoluene	Ave	1.072	1.088	0.0100	10.1	10.0	1.5	20.0
1,3,5-Trimethylbenzene	Ave	3.183	3.036	0.0100	9.54	10.0	-4.6	20.0
4-Chlorotoluene	Ave	1.038	0.9930	0.0100	9.56	10.0	-4.4	20.0
tert-Butylbenzene	Ave	2.758	2.345	0.0100	8.50	10.0	-15.0	20.0
1,2,4-Trimethylbenzene	Ave	3.267	3.033	0.0100	9.29	10.0	-7.1	20.0
3,4-Dichlorobenzotrifluoride	Ave	1.032	1.030	0.0100	9.98	10.0	-0.2	20.0
sec-Butylbenzene	Ave	3.881	3.528	0.0100	9.09	10.0	-9.1	20.0
1,3-Dichlorobenzene	Ave	1.705	1.639	0.6000	9.61	10.0	-3.9	20.0
4-Isopropyltoluene	Ave	3.204	2.865	0.0100	8.94	10.0	-10.6	20.0

FORM VII  
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Lab Sample ID: CCVIS 180-138685/2 Calibration Date: 04/16/2015 10:19

Instrument ID: CHHP5 Calib Start Date: 03/16/2015 12:41

GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 03/16/2015 16:17

Lab File ID: 50416002.D Conc. Units: ug/L Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,4-Dichlorobenzene	Ave	1.741	1.705	0.5000	9.79	10.0	-2.1	20.0
2,4-Dichlorobenzotrifluoride	Ave	0.9669	0.9241	0.0100	9.56	10.0	-4.4	20.0
2,5-Dichlorobenzotrifluoride	Ave	1.082	1.053	0.0100	9.73	10.0	-2.7	20.0
n-Butylbenzene	Ave	2.918	2.539	0.0100	8.70	10.0	-13.0	20.0
1,2-Dichlorobenzene	Ave	1.579	1.516	0.4000	9.61	10.0	-3.9	20.0
1,2-Dibromo-3-Chloropropane	Ave	0.1292	0.1147	0.0500	8.88	10.0	-11.2	20.0
2,4- & 2,5- & 2,6- Dichlorotoluene	Ave	1.194	1.010	0.0100	25.4	30.0	-15.4	20.0
2,3- & 3,4- Dichlorotoluene	Ave	1.161	0.9227	0.0100	15.9	20.0	-20.5*	20.0
1,2,4-Trichlorobenzene	Ave	0.8219	0.5962	0.2000	7.25	10.0	-27.5*	20.0
Hexachlorobutadiene	Ave	0.3941	0.3337	0.0100	8.47	10.0	-15.3	20.0
Naphthalene	Ave	2.158	1.314	0.0100	6.09	10.0	-39.1*	20.0
1,2,3-Trichlorobenzene	Ave	0.6740	0.4959	0.0100	7.36	10.0	-26.4*	20.0
2,4,5-Trichlorotoluene	Ave	0.3624	0.1879	0.0100	5.18	10.0	-48.2*	20.0
2,3,6-Trichlorotoluene	Ave	0.3273	0.1646	0.0100	5.03	10.0	-49.7*	20.0
Dibromofluoromethane (Surr)	Ave	0.2274	0.2206		9.70	10.0	-3.0	20.0
1,2-Dichloroethane-d4 (Surr)	Ave	0.2998	0.2865		9.56	10.0	-4.4	20.0
Toluene-d8 (Surr)	Ave	3.986	3.995		10.0	10.0	0.2	20.0
4-Bromofluorobenzene (Surr)	Ave	1.436	1.331		9.27	10.0	-7.3	20.0

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\50416002.D  
 Lims ID: CCVIS  
 Client ID:  
 Sample Type: CCVIS  
 Inject. Date: 16-Apr-2015 10:19:30 ALS Bottle#: 2 Worklist Smp#: 2  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: CCVIS  
 Misc. Info.: 180-0006494-002  
 Operator ID: 001562 Instrument ID: CHHP5  
 Sublist: chrom-MSVOA\_LL\_CHHP5\*sub11  
 Method: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 12:37:09 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: RT Order ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICal File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 16-Apr-2015 11:02:35

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.317	4.317	0.000	0	154447	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.268	7.268	0.000	95	501487	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.358	10.358	0.000	80	113593	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.682	12.682	0.000	92	167378	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.525	6.525	0.000	44	110629	50.0	48.5	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.897	6.897	0.000	0	143684	50.0	47.8	
\$ 7 Toluene-d8 (Surr)	98	8.916	8.916	0.000	81	453854	50.0	50.1	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.532	11.532	0.000	86	151139	50.0	46.3	
11 Dichlorodifluoromethane	85	1.628	1.628	0.000	68	130657	50.0	60.8	
12 Chloromethane	50	1.786	1.786	0.000	82	157687	50.0	53.1	
13 Vinyl chloride	62	1.914	1.914	0.000	80	180719	50.0	54.5	
14 Butadiene	39	1.957	1.957	0.000	95	195408	50.0	51.6	
15 Bromomethane	94	2.267	2.267	0.000	92	98878	50.0	55.6	
16 Chloroethane	64	2.413	2.413	0.000	94	127441	50.0	55.5	
17 Dichlorofluoromethane	67	2.669	2.669	0.000	82	288239	50.0	55.0	
18 Trichlorofluoromethane	101	2.735	2.735	0.000	81	195594	50.0	49.2	
20 Ethyl ether	59	3.094	3.094	0.000	90	131990	50.0	50.3	
21 Acrolein	56	3.265	3.265	0.000	87	42856	150.0	134.5	
22 1,1-Dichloroethene	96	3.386	3.386	0.000	97	142916	50.0	49.4	
23 1,1,2-Trichloro-1,2,2-trif	101	3.441	3.441	0.000	79	153956	50.0	52.6	
24 Acetone	43	3.496	3.496	0.000	92	119447	100.0	116.3	
25 Iodomethane	142	3.618	3.618	0.000	97	192601	50.0	47.9	M
26 Carbon disulfide	76	3.691	3.691	0.000	99	274206	50.0	38.8	
28 3-Chloro-1-propene	76	3.946	3.946	0.000	83	75667	50.0	49.5	
30 Methyl acetate	43	4.025	4.025	0.000	95	587601	250.0	244.5	
31 Methylene Chloride	84	4.147	4.147	0.000	87	167759	50.0	50.2	
32 2-Methyl-2-propanol	59	4.439	4.439	0.000	77	87509	500.0	481.0	
33 Acrylonitrile	53	4.548	4.548	0.000	97	608751	500.0	492.4	
34 trans-1,2-Dichloroethene	96	4.561	4.561	0.000	59	149546	50.0	50.0	
35 Methyl tert-butyl ether	73	4.591	4.591	0.000	87	317613	50.0	48.0	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
36 Hexane	57	4.986	4.986	0.000	94	210351	50.0	44.0	
37 1,1-Dichloroethane	63	5.175	5.175	0.000	86	272593	50.0	51.1	
38 Vinyl acetate	43	5.291	5.291	0.000	97	170839	50.0	45.1	
44 2,2-Dichloropropane	77	5.929	5.929	0.000	54	85847	50.0	64.3	
45 cis-1,2-Dichloroethene	96	5.935	5.935	0.000	70	151521	50.0	48.1	
46 2-Butanone (MEK)	43	5.990	5.990	0.000	100	148344	100.0	90.3	
49 Chlorobromomethane	128	6.227	6.227	0.000	79	66946	50.0	49.1	
51 Tetrahydrofuran	42	6.282	6.282	0.000	89	86258	100.0	83.9	
52 Chloroform	83	6.337	6.337	0.000	83	252719	50.0	52.1	
53 1,1,1-Trichloroethane	97	6.525	6.525	0.000	55	172915	50.0	55.8	
54 Cyclohexane	56	6.580	6.580	0.000	93	268386	50.0	45.1	
56 Carbon tetrachloride	117	6.720	6.720	0.000	63	146186	50.0	58.8	
55 1,1-Dichloropropene	75	6.720	6.720	0.000	90	190150	50.0	47.3	
57 Isobutyl alcohol	41	6.945	6.945	0.000	44	95937	1250.0	1433.2	
58 Benzene	78	6.951	6.951	0.000	97	616081	50.0	51.8	
59 1,2-Dichloroethane	62	6.982	6.982	0.000	93	197462	50.0	50.7	
62 n-Heptane	43	7.280	7.280	0.000	71	187695	50.0	46.0	
64 Trichloroethene	130	7.669	7.669	0.000	92	136100	50.0	45.7	
66 Methylcyclohexane	83	7.858	7.858	0.000	92	238211	50.0	44.8	
67 1,2-Dichloropropane	63	7.900	7.900	0.000	88	151791	50.0	51.6	
68 Dibromomethane	93	8.022	8.022	0.000	88	80573	50.0	50.9	
70 1,4-Dioxane	88	8.052	8.052	0.000	95	26394	1000.0	852.8	
71 Dichlorobromomethane	83	8.192	8.192	0.000	97	165176	50.0	51.1	
73 2-Chloroethyl vinyl ether	63	8.515	8.515	0.000	91	144323	100.0	87.1	
74 cis-1,3-Dichloropropene	75	8.655	8.655	0.000	90	162730	50.0	52.2	
75 4-Methyl-2-pentanone (MIBK)	43	8.819	8.819	0.000	96	256983	100.0	83.6	
76 Toluene	91	8.989	8.989	0.000	98	627029	50.0	53.9	
77 trans-1,3-Dichloropropene	75	9.220	9.220	0.000	83	122500	50.0	58.3	
78 Ethyl methacrylate	69	9.318	9.318	0.000	89	126215	50.0	46.0	
79 1,1,2-Trichloroethane	97	9.397	9.397	0.000	87	114530	50.0	52.5	
80 Tetrachloroethene	164	9.531	9.531	0.000	90	117891	50.0	51.8	
81 1,3-Dichloropropane	76	9.561	9.561	0.000	86	209372	50.0	51.6	
82 2-Hexanone	43	9.652	9.652	0.000	98	212529	100.0	90.5	
84 Chlorodibromomethane	129	9.786	9.786	0.000	86	93975	50.0	53.9	
85 Ethylene Dibromide	107	9.902	9.902	0.000	93	109155	50.0	52.4	
86 3-Chlorobenzotrifluoride	180	10.370	10.370	0.000	93	240685	50.0	54.2	
87 Chlorobenzene	112	10.388	10.388	0.000	89	378491	50.0	51.3	
88 4-Chlorobenzotrifluoride	180	10.425	10.425	0.000	91	225219	50.0	52.4	
89 1,1,1,2-Tetrachloroethane	131	10.474	10.474	0.000	84	110371	50.0	58.0	
90 Ethylbenzene	106	10.498	10.498	0.000	99	207689	50.0	49.1	
91 m-Xylene & p-Xylene	106	10.614	10.614	0.000	0	252888	50.0	48.9	
92 o-Xylene	106	11.009	11.009	0.000	92	236713	50.0	46.8	
93 Styrene	104	11.021	11.021	0.000	91	404434	50.0	49.6	
94 Bromoform	173	11.204	11.204	0.000	90	56620	50.0	52.6	
96 2-Chlorobenzotrifluoride	180	11.271	11.271	0.000	95	227159	50.0	51.2	
97 Isopropylbenzene	105	11.380	11.380	0.000	97	599241	50.0	47.4	
99 1,1,2,2-Tetrachloroethane	83	11.672	11.672	0.000	75	161913	50.0	51.7	
100 Bromobenzene	156	11.678	11.678	0.000	91	144988	50.0	46.8	
101 1,2,3-Trichloropropane	110	11.721	11.721	0.000	54	46868	50.0	46.0	
102 trans-1,4-Dichloro-2-buten	53	11.727	11.727	0.000	59	38579	50.0	45.6	
103 N-Propylbenzene	120	11.788	11.788	0.000	86	173083	50.0	45.3	
104 2-Chlorotoluene	126	11.873	11.873	0.000	94	147490	50.0	45.9	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
105 3-Chlorotoluene	126	11.934	11.934	0.000	64	182079	50.0	50.7	
106 1,3,5-Trimethylbenzene	105	11.958	11.958	0.000	92	508206	50.0	47.7	
107 4-Chlorotoluene	126	11.976	11.976	0.000	99	166210	50.0	47.8	
108 tert-Butylbenzene	119	12.287	12.287	0.000	71	392523	50.0	42.5	
110 1,2,4-Trimethylbenzene	105	12.335	12.335	0.000	88	507728	50.0	46.4	
111 1,2-dichloro-4-(trifluoromethyl)	214	12.402	12.402	0.000	96	172407	50.0	49.9	
112 sec-Butylbenzene	105	12.506	12.506	0.000	94	590451	50.0	45.4	
113 1,3-Dichlorobenzene	146	12.615	12.615	0.000	83	274288	50.0	48.1	
114 4-Isopropyltoluene	119	12.645	12.645	0.000	88	479532	50.0	44.7	
115 1,4-Dichlorobenzene	146	12.706	12.706	0.000	94	285346	50.0	49.0	
116 2,4-Dichloro-1-(trifluoromethyl)	214	12.755	12.755	0.000	93	154677	50.0	47.8	
118 2,5-Dichlorobenzotrifluoride	214	12.804	12.804	0.000	0	176236	50.0	48.7	
120 n-Butylbenzene	91	13.059	13.059	0.000	95	425011	50.0	43.5	
121 1,2-Dichlorobenzene	146	13.077	13.077	0.000	95	253815	50.0	48.0	
122 1,2-Dibromo-3-Chloropropan	75	13.862	13.862	0.000	69	19200	50.0	44.4	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125	14.008	14.008	0.000	0	507029	150.0	126.9	
124 1,3,5-Trichlorobenzene	180	14.069	14.069	0.000	95	149503	50.0	46.6	
125 2,3- & 3,4- Dichlorotoluene	125	14.422	14.422	0.000	0	308871	100.0	79.5	
126 1,2,4-Trichlorobenzene	180	14.690	14.690	0.000	92	99793	50.0	36.3	
127 Hexachlorobutadiene	225	14.860	14.860	0.000	89	55845	50.0	42.3	
128 Naphthalene	128	14.939	14.939	0.000	97	219958	50.0	30.4	
129 1,2,3-Trichlorobenzene	180	15.182	15.182	0.000	94	82996	50.0	36.8	
131 2,4,5-Trichlorotoluene	159	15.961	15.961	0.000	0	31451	50.0	25.9	
130 2,3,6-Trichlorotoluene	159	16.058	16.058	0.000	93	27555	50.0	25.2	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
S 133 Xylenes, Total	106				0		100.0	95.6	
S 134 1,2-Dichloroethene, Total	96				0		100.0	98.1	
S 135 1,3-Dichloropropene, Total	1				0		100.0	110.5	

**QC Flag Legend**

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

**Reagents:**

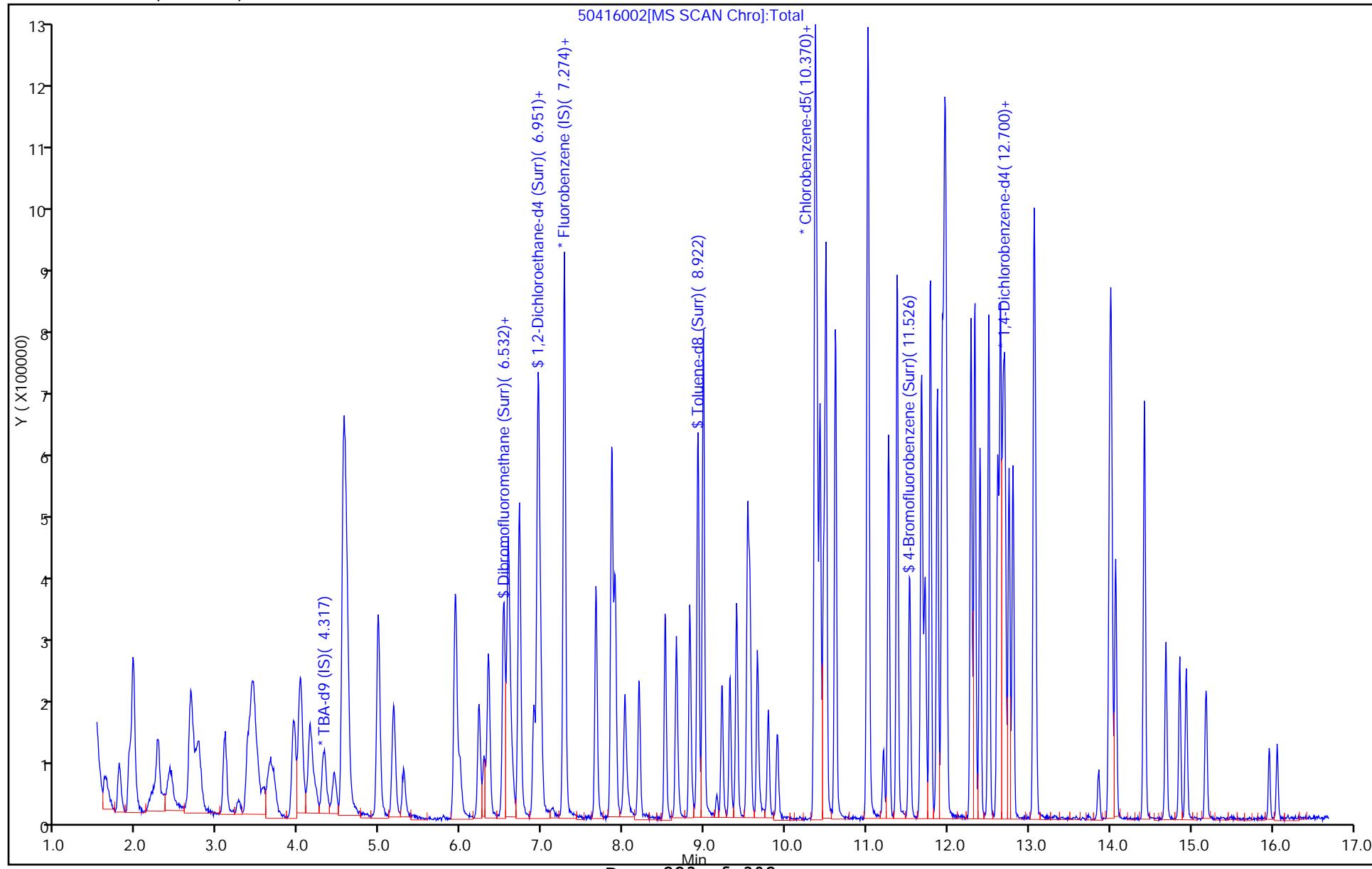
voaW 135tcb a_00001	Amount Added: 2.00	Units: uL	
VOA8260VOAPRI_00110	Amount Added: 2.00	Units: uL	
voaWKetpri Re_00004	Amount Added: 2.00	Units: uL	
voaWeemixPRI_00002	Amount Added: 2.00	Units: uL	
voaW VA pri R_00005	Amount Added: 2.00	Units: uL	
voaW2-cl 2ndR_00002	Amount Added: 2.00	Units: uL	
VOAACRPRI_00005	Amount Added: 6.00	Units: uL	
VOA8260INT_00031	Amount Added: 2.00	Units: uL	Run Reagent
VOA8260SURR_00033	Amount Added: 2.00	Units: uL	Run Reagent

Report Date: 16-Apr-2015 12:37:11

Chrom Revision: 2.2 13-Mar-2015 11:20:44

## TestAmerica Pittsburgh

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150416-6494.b\\50416002.D  
Injection Date: 16-Apr-2015 10:19:30 Instrument ID: CHHP5  
Lims ID: CCVIS Operator ID: 001562  
Client ID:  
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 2  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)



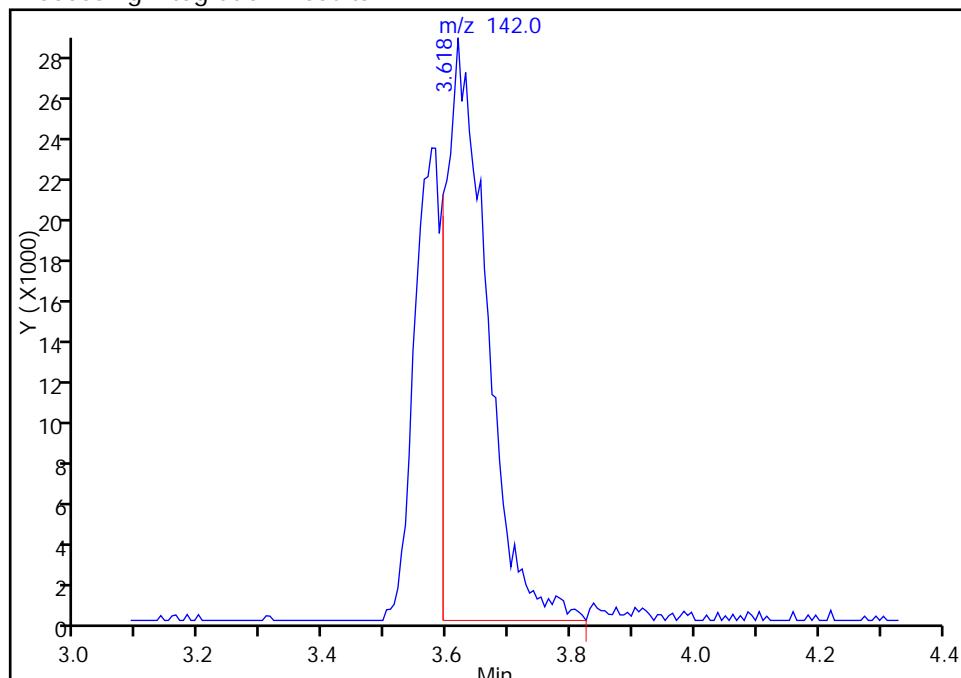
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\50416002.D  
 Injection Date: 16-Apr-2015 10:19:30 Instrument ID: CHHP5  
 Lims ID: CCVIS  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 2 Worklist Smp#: 2  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 25 Iodomethane, CAS: 74-88-4

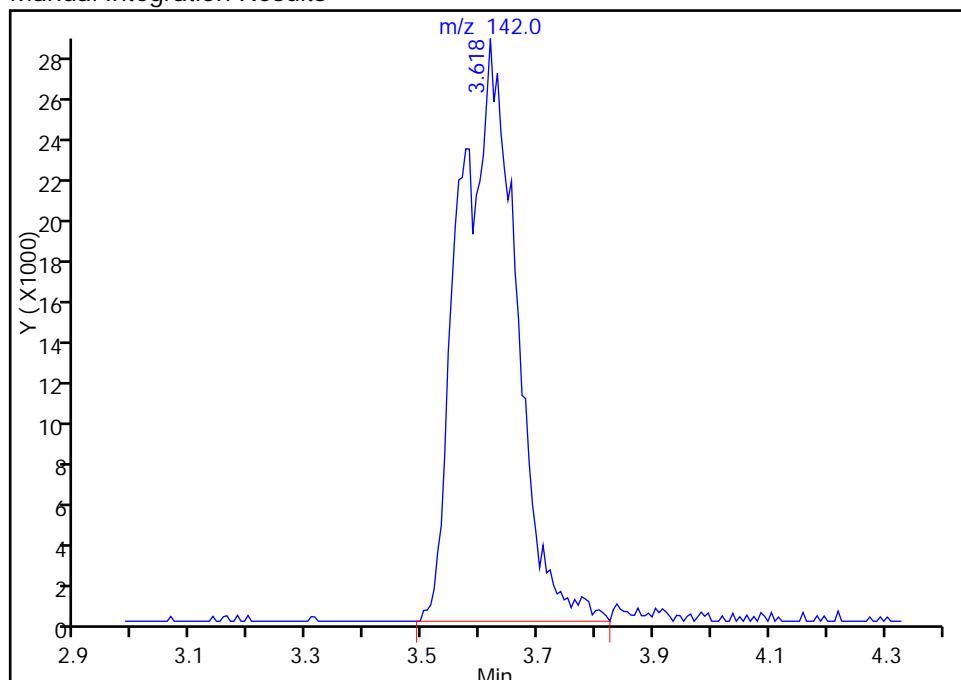
RT: 3.62  
 Area: 128815  
 Amount: 32.064526  
 Amount Units: ng

## Processing Integration Results



RT: 3.62  
 Area: 192601  
 Amount: 47.942086  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 16-Apr-2015 11:02:35

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

FORM VII  
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Lab Sample ID: CCVIS 180-138685/2 Calibration Date: 04/16/2015 10:19

Instrument ID: CHHP5 Calib Start Date: 03/18/2015 13:31

GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 03/18/2015 16:19

Lab File ID: 50416002.D Conc. Units: ug/L Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
2-Chloroethyl vinyl ether	Ave	0.1652	0.1439	0.0100	17.4	20.0	-12.9	20.0
1,3,5-Trichlorobenzene	Ave	0.9577	0.8932	0.0100	9.33	10.0	-6.7	20.0

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\50416002.D  
 Lims ID: CCVIS  
 Client ID:  
 Sample Type: CCVIS  
 Inject. Date: 16-Apr-2015 10:19:30 ALS Bottle#: 2 Worklist Smp#: 2  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: CCVIS  
 Misc. Info.: 180-0006494-002  
 Operator ID: 001562 Instrument ID: CHHP5  
 Sublist: chrom-MSVOA\_LL\_CHHP5\*sub11  
 Method: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 12:37:09 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: RT Order ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICal File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 16-Apr-2015 11:02:35

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.317	4.317	0.000	0	154447	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.268	7.268	0.000	95	501487	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.358	10.358	0.000	80	113593	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.682	12.682	0.000	92	167378	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.525	6.525	0.000	44	110629	50.0	48.5	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.897	6.897	0.000	0	143684	50.0	47.8	
\$ 7 Toluene-d8 (Surr)	98	8.916	8.916	0.000	81	453854	50.0	50.1	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.532	11.532	0.000	86	151139	50.0	46.3	
11 Dichlorodifluoromethane	85	1.628	1.628	0.000	68	130657	50.0	60.8	
12 Chloromethane	50	1.786	1.786	0.000	82	157687	50.0	53.1	
13 Vinyl chloride	62	1.914	1.914	0.000	80	180719	50.0	54.5	
14 Butadiene	39	1.957	1.957	0.000	95	195408	50.0	51.6	
15 Bromomethane	94	2.267	2.267	0.000	92	98878	50.0	55.6	
16 Chloroethane	64	2.413	2.413	0.000	94	127441	50.0	55.5	
17 Dichlorofluoromethane	67	2.669	2.669	0.000	82	288239	50.0	55.0	
18 Trichlorofluoromethane	101	2.735	2.735	0.000	81	195594	50.0	49.2	
20 Ethyl ether	59	3.094	3.094	0.000	90	131990	50.0	50.3	
21 Acrolein	56	3.265	3.265	0.000	87	42856	150.0	134.5	
22 1,1-Dichloroethene	96	3.386	3.386	0.000	97	142916	50.0	49.4	
23 1,1,2-Trichloro-1,2,2-trif	101	3.441	3.441	0.000	79	153956	50.0	52.6	
24 Acetone	43	3.496	3.496	0.000	92	119447	100.0	116.3	
25 Iodomethane	142	3.618	3.618	0.000	97	192601	50.0	47.9	M
26 Carbon disulfide	76	3.691	3.691	0.000	99	274206	50.0	38.8	
28 3-Chloro-1-propene	76	3.946	3.946	0.000	83	75667	50.0	49.5	
30 Methyl acetate	43	4.025	4.025	0.000	95	587601	250.0	244.5	
31 Methylene Chloride	84	4.147	4.147	0.000	87	167759	50.0	50.2	
32 2-Methyl-2-propanol	59	4.439	4.439	0.000	77	87509	500.0	481.0	
33 Acrylonitrile	53	4.548	4.548	0.000	97	608751	500.0	492.4	
34 trans-1,2-Dichloroethene	96	4.561	4.561	0.000	59	149546	50.0	50.0	
35 Methyl tert-butyl ether	73	4.591	4.591	0.000	87	317613	50.0	48.0	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
36 Hexane	57	4.986	4.986	0.000	94	210351	50.0	44.0	
37 1,1-Dichloroethane	63	5.175	5.175	0.000	86	272593	50.0	51.1	
38 Vinyl acetate	43	5.291	5.291	0.000	97	170839	50.0	45.1	
44 2,2-Dichloropropane	77	5.929	5.929	0.000	54	85847	50.0	64.3	
45 cis-1,2-Dichloroethene	96	5.935	5.935	0.000	70	151521	50.0	48.1	
46 2-Butanone (MEK)	43	5.990	5.990	0.000	100	148344	100.0	90.3	
49 Chlorobromomethane	128	6.227	6.227	0.000	79	66946	50.0	49.1	
51 Tetrahydrofuran	42	6.282	6.282	0.000	89	86258	100.0	83.9	
52 Chloroform	83	6.337	6.337	0.000	83	252719	50.0	52.1	
53 1,1,1-Trichloroethane	97	6.525	6.525	0.000	55	172915	50.0	55.8	
54 Cyclohexane	56	6.580	6.580	0.000	93	268386	50.0	45.1	
56 Carbon tetrachloride	117	6.720	6.720	0.000	63	146186	50.0	58.8	
55 1,1-Dichloropropene	75	6.720	6.720	0.000	90	190150	50.0	47.3	
57 Isobutyl alcohol	41	6.945	6.945	0.000	44	95937	1250.0	1433.2	
58 Benzene	78	6.951	6.951	0.000	97	616081	50.0	51.8	
59 1,2-Dichloroethane	62	6.982	6.982	0.000	93	197462	50.0	50.7	
62 n-Heptane	43	7.280	7.280	0.000	71	187695	50.0	46.0	
64 Trichloroethene	130	7.669	7.669	0.000	92	136100	50.0	45.7	
66 Methylcyclohexane	83	7.858	7.858	0.000	92	238211	50.0	44.8	
67 1,2-Dichloropropane	63	7.900	7.900	0.000	88	151791	50.0	51.6	
68 Dibromomethane	93	8.022	8.022	0.000	88	80573	50.0	50.9	
70 1,4-Dioxane	88	8.052	8.052	0.000	95	26394	1000.0	852.8	
71 Dichlorobromomethane	83	8.192	8.192	0.000	97	165176	50.0	51.1	
73 2-Chloroethyl vinyl ether	63	8.515	8.515	0.000	91	144323	100.0	87.1	
74 cis-1,3-Dichloropropene	75	8.655	8.655	0.000	90	162730	50.0	52.2	
75 4-Methyl-2-pentanone (MIBK)	43	8.819	8.819	0.000	96	256983	100.0	83.6	
76 Toluene	91	8.989	8.989	0.000	98	627029	50.0	53.9	
77 trans-1,3-Dichloropropene	75	9.220	9.220	0.000	83	122500	50.0	58.3	
78 Ethyl methacrylate	69	9.318	9.318	0.000	89	126215	50.0	46.0	
79 1,1,2-Trichloroethane	97	9.397	9.397	0.000	87	114530	50.0	52.5	
80 Tetrachloroethene	164	9.531	9.531	0.000	90	117891	50.0	51.8	
81 1,3-Dichloropropane	76	9.561	9.561	0.000	86	209372	50.0	51.6	
82 2-Hexanone	43	9.652	9.652	0.000	98	212529	100.0	90.5	
84 Chlorodibromomethane	129	9.786	9.786	0.000	86	93975	50.0	53.9	
85 Ethylene Dibromide	107	9.902	9.902	0.000	93	109155	50.0	52.4	
86 3-Chlorobenzotrifluoride	180	10.370	10.370	0.000	93	240685	50.0	54.2	
87 Chlorobenzene	112	10.388	10.388	0.000	89	378491	50.0	51.3	
88 4-Chlorobenzotrifluoride	180	10.425	10.425	0.000	91	225219	50.0	52.4	
89 1,1,1,2-Tetrachloroethane	131	10.474	10.474	0.000	84	110371	50.0	58.0	
90 Ethylbenzene	106	10.498	10.498	0.000	99	207689	50.0	49.1	
91 m-Xylene & p-Xylene	106	10.614	10.614	0.000	0	252888	50.0	48.9	
92 o-Xylene	106	11.009	11.009	0.000	92	236713	50.0	46.8	
93 Styrene	104	11.021	11.021	0.000	91	404434	50.0	49.6	
94 Bromoform	173	11.204	11.204	0.000	90	56620	50.0	52.6	
96 2-Chlorobenzotrifluoride	180	11.271	11.271	0.000	95	227159	50.0	51.2	
97 Isopropylbenzene	105	11.380	11.380	0.000	97	599241	50.0	47.4	
99 1,1,2,2-Tetrachloroethane	83	11.672	11.672	0.000	75	161913	50.0	51.7	
100 Bromobenzene	156	11.678	11.678	0.000	91	144988	50.0	46.8	
101 1,2,3-Trichloropropane	110	11.721	11.721	0.000	54	46868	50.0	46.0	
102 trans-1,4-Dichloro-2-buten	53	11.727	11.727	0.000	59	38579	50.0	45.6	
103 N-Propylbenzene	120	11.788	11.788	0.000	86	173083	50.0	45.3	
104 2-Chlorotoluene	126	11.873	11.873	0.000	94	147490	50.0	45.9	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
105 3-Chlorotoluene	126	11.934	11.934	0.000	64	182079	50.0	50.7	
106 1,3,5-Trimethylbenzene	105	11.958	11.958	0.000	92	508206	50.0	47.7	
107 4-Chlorotoluene	126	11.976	11.976	0.000	99	166210	50.0	47.8	
108 tert-Butylbenzene	119	12.287	12.287	0.000	71	392523	50.0	42.5	
110 1,2,4-Trimethylbenzene	105	12.335	12.335	0.000	88	507728	50.0	46.4	
111 1,2-dichloro-4-(trifluoromethyl)	214	12.402	12.402	0.000	96	172407	50.0	49.9	
112 sec-Butylbenzene	105	12.506	12.506	0.000	94	590451	50.0	45.4	
113 1,3-Dichlorobenzene	146	12.615	12.615	0.000	83	274288	50.0	48.1	
114 4-Isopropyltoluene	119	12.645	12.645	0.000	88	479532	50.0	44.7	
115 1,4-Dichlorobenzene	146	12.706	12.706	0.000	94	285346	50.0	49.0	
116 2,4-Dichloro-1-(trifluoromethyl)	214	12.755	12.755	0.000	93	154677	50.0	47.8	
118 2,5-Dichlorobenzotrifluoride	214	12.804	12.804	0.000	0	176236	50.0	48.7	
120 n-Butylbenzene	91	13.059	13.059	0.000	95	425011	50.0	43.5	
121 1,2-Dichlorobenzene	146	13.077	13.077	0.000	95	253815	50.0	48.0	
122 1,2-Dibromo-3-Chloropropan	75	13.862	13.862	0.000	69	19200	50.0	44.4	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125	14.008	14.008	0.000	0	507029	150.0	126.9	
124 1,3,5-Trichlorobenzene	180	14.069	14.069	0.000	95	149503	50.0	46.6	
125 2,3- & 3,4- Dichlorotoluene	125	14.422	14.422	0.000	0	308871	100.0	79.5	
126 1,2,4-Trichlorobenzene	180	14.690	14.690	0.000	92	99793	50.0	36.3	
127 Hexachlorobutadiene	225	14.860	14.860	0.000	89	55845	50.0	42.3	
128 Naphthalene	128	14.939	14.939	0.000	97	219958	50.0	30.4	
129 1,2,3-Trichlorobenzene	180	15.182	15.182	0.000	94	82996	50.0	36.8	
131 2,4,5-Trichlorotoluene	159	15.961	15.961	0.000	0	31451	50.0	25.9	
130 2,3,6-Trichlorotoluene	159	16.058	16.058	0.000	93	27555	50.0	25.2	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
S 133 Xylenes, Total	106				0		100.0	95.6	
S 134 1,2-Dichloroethene, Total	96				0		100.0	98.1	
S 135 1,3-Dichloropropene, Total	1				0		100.0	110.5	

**QC Flag Legend**

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

**Reagents:**

voaW 135tcb a_00001	Amount Added: 2.00	Units: uL	
VOA8260VOAPRI_00110	Amount Added: 2.00	Units: uL	
voaWKetpri Re_00004	Amount Added: 2.00	Units: uL	
voaWeemixPRI_00002	Amount Added: 2.00	Units: uL	
voaW VA pri R_00005	Amount Added: 2.00	Units: uL	
voaW2-cl 2ndR_00002	Amount Added: 2.00	Units: uL	
VOAACRPRI_00005	Amount Added: 6.00	Units: uL	
VOA8260INT_00031	Amount Added: 2.00	Units: uL	Run Reagent
VOA8260SURR_00033	Amount Added: 2.00	Units: uL	Run Reagent

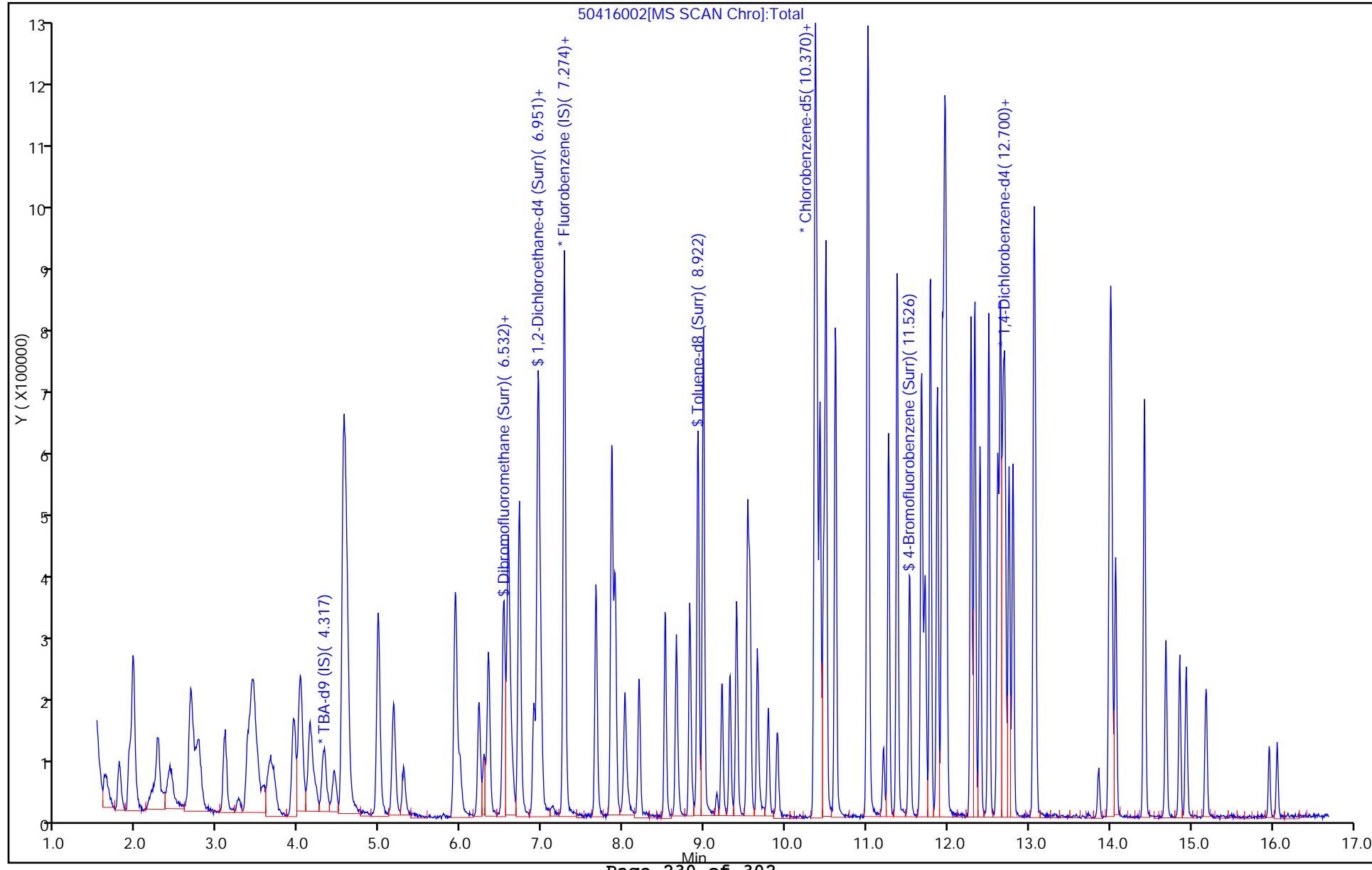
Report Date: 16-Apr-2015 12:37:11

Chrom Revision: 2.2 13-Mar-2015 11:20:44

## TestAmerica Pittsburgh

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150416-6494.b\\50416002.D  
Injection Date: 16-Apr-2015 10:19:30 Instrument ID: CHHP5  
Lims ID: CCVIS Operator ID: 001562  
Client ID:  
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 2  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)

Worklist Smp#: 2



TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316001.D  
 Lims ID: BFB  
 Client ID:  
 Sample Type: BFB  
 Inject. Date: 16-Mar-2015 10:49:30 ALS Bottle#: 1 Worklist Smp#: 1  
 Injection Vol: 5.0 mL Dil. Factor: 1.0000  
 Sample Info: BFB  
 Misc. Info.: 180-0006031-001  
 Operator ID: 001562 Instrument ID: CHHP5  
 Method: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 17-Mar-2015 10:59:24 Calib Date: 16-Mar-2015 16:17:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICAL File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316013.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK012

First Level Reviewer: fergusond Date: 16-Mar-2015 11:15:59

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
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\$ 10 BFB

95 8.341 8.341 0.000 0 133980

NR NR

### QC Flag Legend

Processing Flags

NR - Missing Quant Standard

### Reagents:

VOA BFB 25\_00001

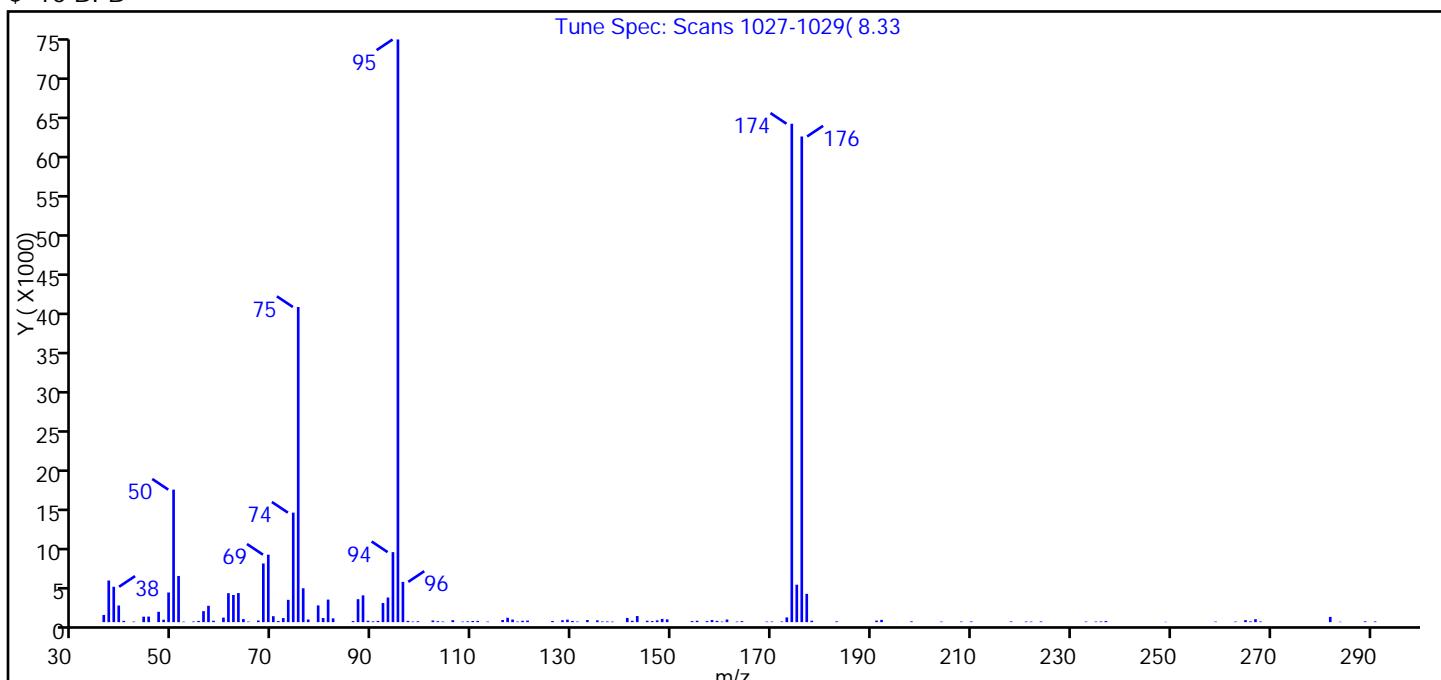
Amount Added: 1.00

Units: uL

## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316001.D  
 Injection Date: 16-Mar-2015 10:49:30 Instrument ID: CHHP5  
 Lims ID: BFB  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 1 Worklist Smp#: 1  
 Injection Vol: 5.0 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Tune Method: BFB Method 8260

\$ 10 BFB



m/z	Ion Abundance Criteria	% Relative Abundance
95	Base peak, 100% relative abundance	100.0
50	15 to 40% of m/z 95	22.7
75	30 to 60% of m/z 95	54.1
96	5 to 9% of m/z 95	6.9
173	Less than 2% of m/z 174	0.8 (0.9)
174	50 to 120% of m/z 95	85.5
175	5 to 9% of m/z 174	6.4 (7.5)
176	Greater than 95% but less than 101% of m/z 174	83.4 (97.4)
177	5 to 9% of m/z 176	4.9 (5.8)

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150316-6031.b\\50316001.D\\MSVOA\_LL\_CHHP5.rslt\\spectra.d  
 Injection Date: 16-Mar-2015 10:49:30  
 Spectrum: Tune Spec: Scans 1027-1029( 8.33  
 Base Peak: 95.00  
 Minimum % Base Peak: 0  
 Number of Points: 132

m/z	Y	m/z	Y	m/z	Y	m/z	Y
36.00	922	75.00	40336	119.00	71	173.00	604
37.00	5329	76.00	4335	120.00	170	174.00	63792
38.00	4528	77.00	339	121.00	203	175.00	4791
39.00	2130	79.00	2142	126.00	145	176.00	62160
40.00	163	80.00	527	128.00	241	177.00	3622
42.00	71	81.00	2886	129.00	320	178.00	182
44.00	700	82.00	482	130.00	150	183.00	99
45.00	713	86.00	138	131.00	72	191.00	196
47.00	1323	87.00	2939	133.00	273	192.00	286
48.00	310	88.00	3429	135.00	226	195.00	9
49.00	3792	89.00	182	136.00	81	198.00	98
50.00	16960	90.00	101	137.00	87	204.00	68
51.00	5912	91.00	160	138.00	71	208.00	75
52.00	63	92.00	2448	141.00	541	210.00	85
54.00	83	93.00	3152	142.00	172	218.00	87
55.00	155	94.00	8961	143.00	779	221.00	76
56.00	1409	95.00	74576	145.00	182	222.00	70
57.00	2093	96.00	5155	146.00	133	224.00	88
58.00	180	97.00	159	147.00	227	233.00	73
60.00	582	98.00	71	148.00	412	235.00	76
61.00	3707	99.00	112	149.00	352	236.00	88
62.00	3479	102.00	212	154.00	135	237.00	141
63.00	3721	103.00	120	155.00	179	249.00	43
64.00	392	104.00	75	157.00	135	259.00	70
65.00	71	106.00	253	158.00	274	263.00	71
67.00	207	108.00	68	159.00	163	265.00	262
68.00	7510	109.00	97	160.00	73	266.00	100
69.00	8635	110.00	146	161.00	334	267.00	377
70.00	764	111.00	161	163.00	71	268.00	85
71.00	139	113.00	71	164.00	125	282.00	672
72.00	524	116.00	278	169.00	70	284.00	50
73.00	2854	117.00	558	170.00	78	289.00	99
74.00	14015	118.00	332	172.00	82	291.00	87

m/z	Y	m/z	Y	m/z	Y	m/z	Y
36.00	922	75.00	40336	119.00	71	173.00	604
37.00	5329	76.00	4335	120.00	170	174.00	63792
38.00	4528	77.00	339	121.00	203	175.00	4791
39.00	2130	79.00	2142	126.00	145	176.00	62160
40.00	163	80.00	527	128.00	241	177.00	3622
42.00	71	81.00	2886	129.00	320	178.00	182
44.00	700	82.00	482	130.00	150	183.00	99
45.00	713	86.00	138	131.00	72	191.00	196
47.00	1323	87.00	2939	133.00	273	192.00	286
48.00	310	88.00	3429	135.00	226	195.00	9
49.00	3792	89.00	182	136.00	81	198.00	98
50.00	16960	90.00	101	137.00	87	204.00	68
51.00	5912	91.00	160	138.00	71	208.00	75
52.00	63	92.00	2448	141.00	541	210.00	85
54.00	83	93.00	3152	142.00	172	218.00	87
55.00	155	94.00	8961	143.00	779	221.00	76
56.00	1409	95.00	74576	145.00	182	222.00	70
57.00	2093	96.00	5155	146.00	133	224.00	88
58.00	180	97.00	159	147.00	227	233.00	73
60.00	582	98.00	71	148.00	412	235.00	76
61.00	3707	99.00	112	149.00	352	236.00	88
62.00	3479	102.00	212	154.00	135	237.00	141
63.00	3721	103.00	120	155.00	179	249.00	43
64.00	392	104.00	75	157.00	135	259.00	70
65.00	71	106.00	253	158.00	274	263.00	71
67.00	207	108.00	68	159.00	163	265.00	262
68.00	7510	109.00	97	160.00	73	266.00	100
69.00	8635	110.00	146	161.00	334	267.00	377
70.00	764	111.00	161	163.00	71	268.00	85
71.00	139	113.00	71	164.00	125	282.00	672
72.00	524	116.00	278	169.00	70	284.00	50
73.00	2854	117.00	558	170.00	78	289.00	99
74.00	14015	118.00	332	172.00	82	291.00	87

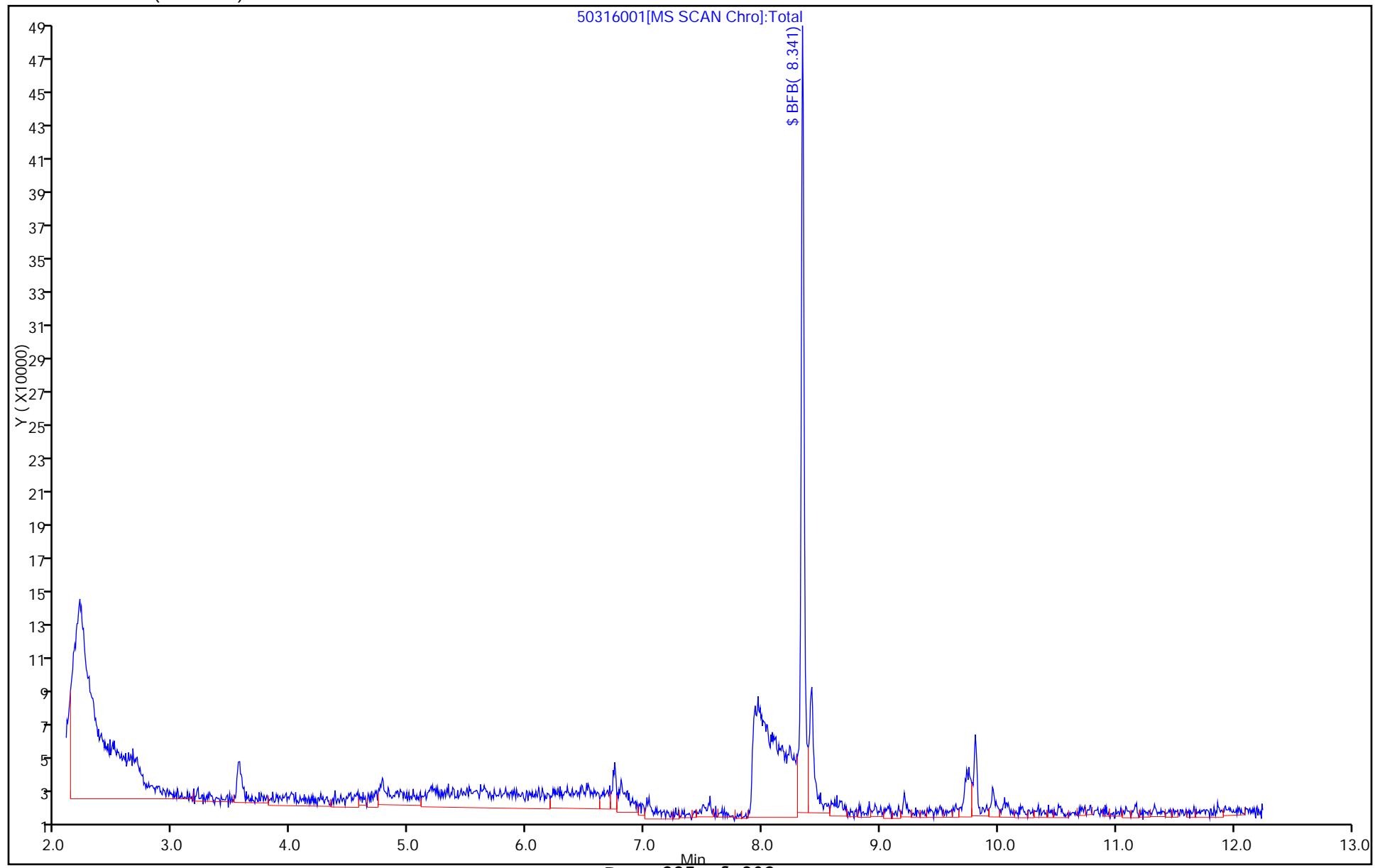
Report Date: 17-Mar-2015 10:59:25

Chrom Revision: 2.2 15-Jan-2015 13:05:58

TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150316-6031.b\50316001.D  
Injection Date: 16-Mar-2015 10:49:30 Instrument ID: CHHP5  
Lims ID: BFB Operator ID: 001562  
Client ID:  
Injection Vol: 5.0 mL Dil. Factor: 1.0000 ALS Bottle#: 1  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)

Worklist Smp#: 1



TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415001.D  
 Lims ID: BFB  
 Client ID:  
 Sample Type: BFB  
 Inject. Date: 15-Apr-2015 12:36:30 ALS Bottle#: 1 Worklist Smp#: 1  
 Injection Vol: 5.0 mL Dil. Factor: 1.0000  
 Sample Info: BFB  
 Misc. Info.: 180-0006480-001  
 Operator ID: 001562 Instrument ID: CHHP5  
 Method: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 07:27:53 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICAL File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 15-Apr-2015 12:53:36

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
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\$ 10 BFB

95 8.337 8.337 0.000 0 75621

NR NR

### QC Flag Legend

Processing Flags

NR - Missing Quant Standard

### Reagents:

VOABFB25\_00060

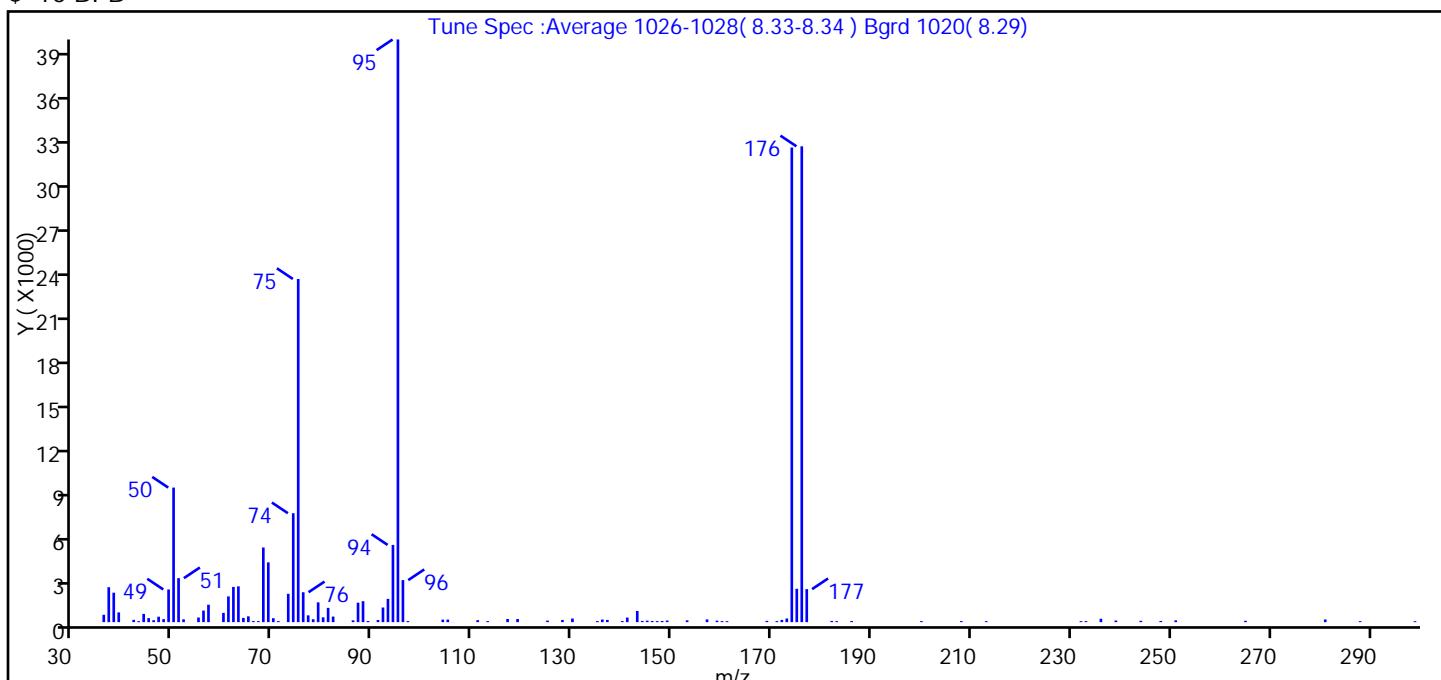
Amount Added: 1.00

Units: uL

## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415001.D  
 Injection Date: 15-Apr-2015 12:36:30 Instrument ID: CHHP5  
 Lims ID: BFB  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 1 Worklist Smp#: 1  
 Injection Vol: 5.0 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Tune Method: BFB Method 8260

\$ 10 BFB



m/z	Ion Abundance Criteria	% Relative Abundance
95	Base peak, 100% relative abundance	100.0
50	15 to 40% of m/z 95	23.1
75	30 to 60% of m/z 95	58.9
96	5 to 9% of m/z 95	7.2
173	Less than 2% of m/z 174	0.6 (0.8)
174	50 to 120% of m/z 95	81.5
175	5 to 9% of m/z 174	5.7 (7.0)
176	Greater than 95% but less than 101% of m/z 174	81.7 (100.2)
177	5 to 9% of m/z 176	5.7 (6.9)

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150415-6480.b\\50415001.D\\MSVOA\_LL\_CHHP5.rslt\\spectra.d  
 Injection Date: 15-Apr-2015 12:36:30  
 Spectrum: Tune Spec :Average 1026-1028( 8.33-8.34 ) Bgrd 1020( 8.29)  
 Base Peak: 95.00  
 Minimum % Base Peak: 0  
 Number of Points: 103

m/z	Y	m/z	Y	m/z	Y	m/z	Y
36.00	503	68.00	5092	105.00	175	169.00	88
37.00	2382	69.00	4076	106.00	5	171.00	73
38.00	2008	70.00	270	111.00	138	172.00	154
39.00	663	71.00	72	113.00	71	173.00	249
42.00	156	73.00	1932	117.00	217	174.00	32376
43.00	70	74.00	7430	119.00	213	175.00	2275
44.00	563	75.00	23408	125.00	106	176.00	32456
45.00	274	76.00	2040	128.00	148	177.00	2253
46.00	132	77.00	462	130.00	248	182.00	82
47.00	371	78.00	199	135.00	75	183.00	67
48.00	205	79.00	1349	136.00	179	186.00	77
49.00	2230	80.00	328	137.00	152	200.00	70
50.00	9178	81.00	975	140.00	72	208.00	73
51.00	3001	82.00	381	141.00	312	213.00	68
52.00	192	86.00	124	143.00	762	232.00	70
55.00	317	87.00	1326	144.00	82	233.00	73
56.00	789	88.00	1430	145.00	101	236.00	229
57.00	1190	89.00	78	146.00	79	239.00	112
60.00	633	91.00	148	147.00	71	244.00	92
61.00	1756	92.00	998	148.00	79	248.00	81
62.00	2401	93.00	1588	149.00	111	251.00	124
63.00	2441	94.00	5268	153.00	123	265.00	93
64.00	289	95.00	39744	157.00	186	281.00	182
65.00	397	96.00	2867	159.00	100	288.00	67
66.00	72	97.00	74	160.00	71	299.00	68
67.00	70	104.00	173	161.00	67		

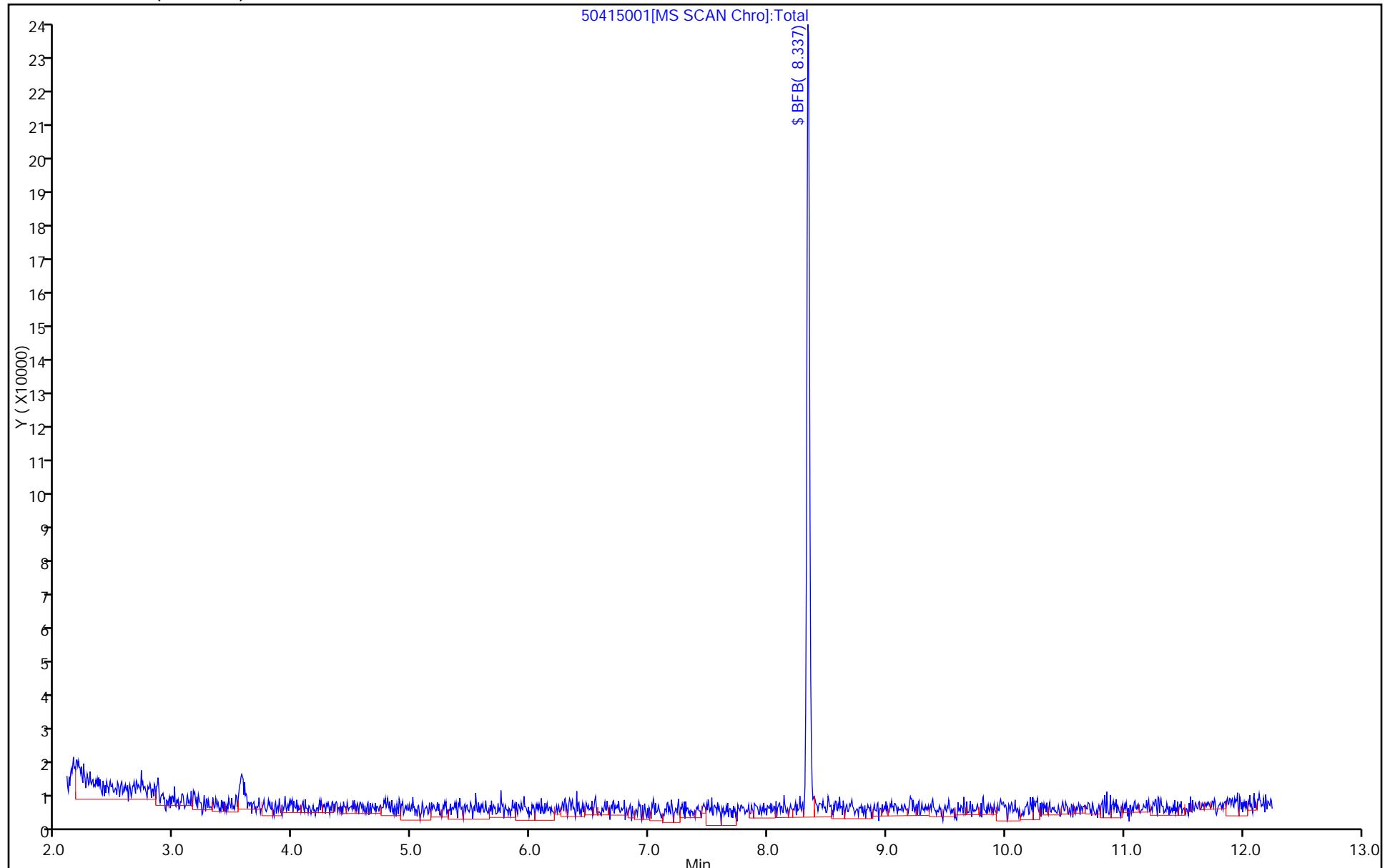
Report Date: 16-Apr-2015 07:27:54

Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Pittsburgh

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150415-6480.b\\50415001.D  
Injection Date: 15-Apr-2015 12:36:30 Instrument ID: CHHP5  
Lims ID: BFB Operator ID: 001562  
Client ID:  
Injection Vol: 5.0 mL Dil. Factor: 1.0000 ALS Bottle#: 1  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)

Worklist Smp#: 1



TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\50416001.D  
 Lims ID: BFB  
 Client ID:  
 Sample Type: BFB  
 Inject. Date: 16-Apr-2015 09:31:30 ALS Bottle#: 1 Worklist Smp#: 1  
 Injection Vol: 5.0 mL Dil. Factor: 1.0000  
 Sample Info: BFB  
 Misc. Info.: 180-0006494-001  
 Operator ID: 001562 Instrument ID: CHHP5  
 Method: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 12:37:07 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICal File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 16-Apr-2015 09:48:12

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
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\$ 10 BFB

95 8.342 8.342 0.000 0 75970

NR NR

### QC Flag Legend

Processing Flags

NR - Missing Quant Standard

### Reagents:

VOABFB25\_00060

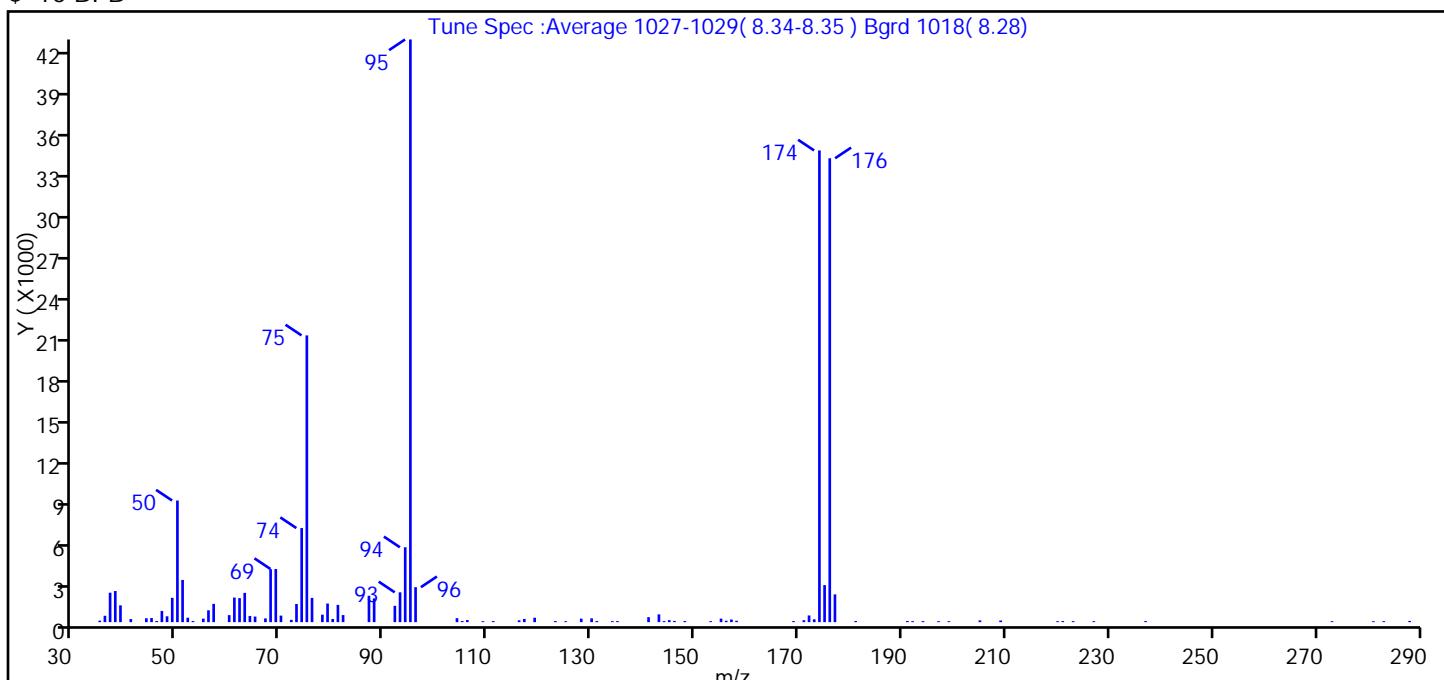
Amount Added: 1.00

Units: uL

## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\50416001.D  
 Injection Date: 16-Apr-2015 09:31:30 Instrument ID: CHHP5  
 Lims ID: BFB  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 1 Worklist Smp#: 1  
 Injection Vol: 5.0 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Tune Method: BFB Method 8260

\$ 10 BFB



m/z	Ion Abundance Criteria	% Relative Abundance
95	Base peak, 100% relative abundance	100.0
50	15 to 40% of m/z 95	20.9
75	30 to 60% of m/z 95	49.2
96	5 to 9% of m/z 95	6.0
173	Less than 2% of m/z 174	0.5 (0.6)
174	50 to 120% of m/z 95	81.0
175	5 to 9% of m/z 174	6.4 (7.9)
176	Greater than 95% but less than 101% of m/z 174	79.6 (98.3)
177	5 to 9% of m/z 176	4.8 (6.0)

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150416-6494.b\\50416001.D\\MSVOA\_LL\_CHHP5.rslt\\spectra.d  
 Injection Date: 16-Apr-2015 09:31:30  
 Spectrum: Tune Spec :Average 1027-1029( 8.34-8.35 ) Bgrd 1018( 8.28)  
 Base Peak: 95.00  
 Minimum % Base Peak: 0  
 Number of Points: 97

m/z	Y	m/z	Y	m/z	Y	m/z	Y
35.00	108	67.00	270	111.00	78	173.00	207
36.00	461	68.00	3840	116.00	133	174.00	34216
37.00	2135	69.00	3854	117.00	235	175.00	2688
38.00	2257	70.00	475	119.00	321	176.00	33648
39.00	1219	72.00	167	123.00	82	177.00	2012
41.00	222	73.00	1315	125.00	67	181.00	82
44.00	280	74.00	6824	128.00	253	191.00	83
45.00	298	75.00	20800	130.00	276	192.00	72
46.00	83	76.00	1759	131.00	82	194.00	75
47.00	812	78.00	539	134.00	73	197.00	70
48.00	423	79.00	1350	135.00	75	199.00	69
49.00	1762	80.00	228	141.00	360	205.00	127
50.00	8822	81.00	1249	143.00	564	209.00	121
51.00	3061	82.00	513	144.00	74	220.00	69
52.00	322	87.00	1913	145.00	146	221.00	79
53.00	78	88.00	1724	146.00	88	223.00	74
55.00	255	92.00	1185	148.00	81	227.00	72
56.00	860	93.00	2164	153.00	72	237.00	77
57.00	1325	94.00	5431	155.00	256	273.00	69
60.00	509	95.00	42264	156.00	103	281.00	67
61.00	1781	96.00	2536	157.00	192	283.00	70
62.00	1741	104.00	288	158.00	99	288.00	89
63.00	2122	105.00	81	169.00	76		
64.00	449	106.00	152	171.00	140		
65.00	405	109.00	67	172.00	486		

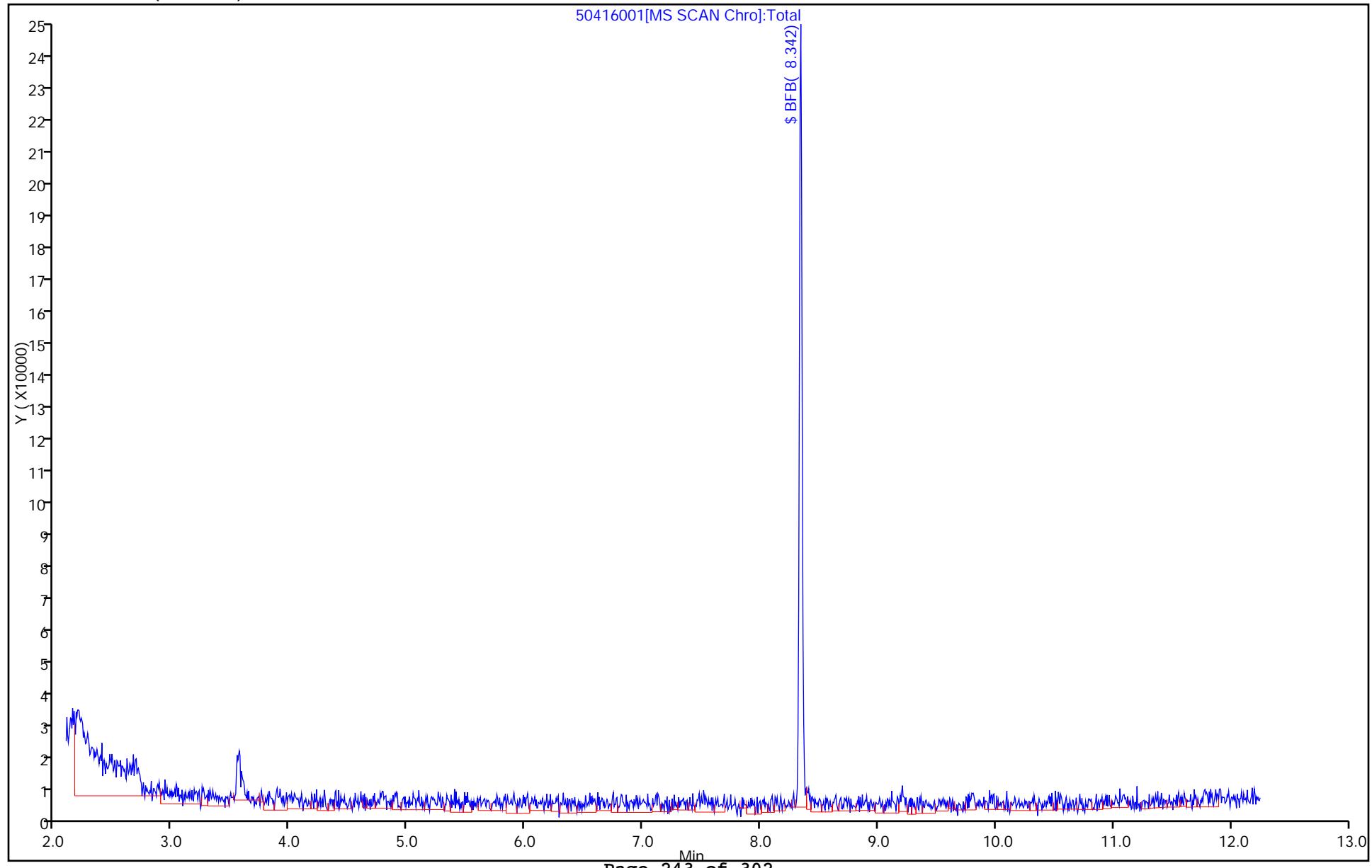
Report Date: 16-Apr-2015 12:37:09

Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Pittsburgh

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150416-6494.b\\50416001.D  
Injection Date: 16-Apr-2015 09:31:30 Instrument ID: CHHP5  
Lims ID: BFB Operator ID: 001562  
Client ID:  
Injection Vol: 5.0 mL Dil. Factor: 1.0000 ALS Bottle#: 1  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)

Worklist Smp#: 1



FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: \_\_\_\_\_ Lab Sample ID: MB 180-138583/5  
Matrix: Water Lab File ID: 50415005.D  
Analysis Method: 8260C Date Collected: \_\_\_\_\_  
Sample wt/vol: 5 (mL) Date Analyzed: 04/15/2015 14:33  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138583 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
74-87-3	Chloromethane	1.0	U	1.0	0.28
75-01-4	Vinyl chloride	1.0	U	1.0	0.23
74-83-9	Bromomethane	1.0	U	1.0	0.31
75-00-3	Chloroethane	1.0	U	1.0	0.21
75-35-4	1,1-Dichloroethene	1.0	U	1.0	0.30
67-64-1	Acetone	5.0	U	5.0	2.5
75-15-0	Carbon disulfide	1.0	U	1.0	0.21
75-09-2	Methylene Chloride	1.0	U	1.0	0.13
156-60-5	trans-1,2-Dichloroethene	1.0	U	1.0	0.17
1634-04-4	Methyl tert-butyl ether	1.0	U	1.0	0.18
75-34-3	1,1-Dichloroethane	1.0	U	1.0	0.12
156-59-2	cis-1,2-Dichloroethene	1.0	U	1.0	0.24
74-97-5	Bromochloromethane	1.0	U	1.0	0.18
78-93-3	2-Butanone (MEK)	5.0	U	5.0	0.55
67-66-3	Chloroform	1.0	U	1.0	0.17
71-55-6	1,1,1-Trichloroethane	1.0	U	1.0	0.29
56-23-5	Carbon tetrachloride	1.0	U	1.0	0.14
71-43-2	Benzene	1.0	U	1.0	0.11
107-06-2	1,2-Dichloroethane	1.0	U	1.0	0.21
79-01-6	Trichloroethene	1.0	U	1.0	0.14
78-87-5	1,2-Dichloropropane	1.0	U	1.0	0.095
75-27-4	Bromodichloromethane	1.0	U	1.0	0.13
10061-01-5	cis-1,3-Dichloropropene	1.0	U	1.0	0.19
108-10-1	4-Methyl-2-pentanone (MIBK)	5.0	U	5.0	0.53
108-88-3	Toluene	1.0	U	1.0	0.15
10061-02-6	trans-1,3-Dichloropropene	1.0	U	1.0	0.15
79-00-5	1,1,2-Trichloroethane	1.0	U	1.0	0.20
127-18-4	Tetrachloroethene	1.0	U	1.0	0.15
591-78-6	2-Hexanone	5.0	U	5.0	0.16
124-48-1	Dibromochloromethane	1.0	U	1.0	0.14
106-93-4	1,2-Dibromoethane (EDB)	1.0	U	1.0	0.18
108-90-7	Chlorobenzene	1.0	U	1.0	0.14
630-20-6	1,1,1,2-Tetrachloroethane	1.0	U	1.0	0.28
100-41-4	Ethylbenzene	1.0	U	1.0	0.23
1330-20-7	Xylenes, Total	3.0	U	3.0	0.49
100-42-5	Styrene	1.0	U	1.0	0.097

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: \_\_\_\_\_ Lab Sample ID: MB 180-138583/5  
Matrix: Water Lab File ID: 50415005.D  
Analysis Method: 8260C Date Collected: \_\_\_\_\_  
Sample wt/vol: 5 (mL) Date Analyzed: 04/15/2015 14:33  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138583 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-25-2	Bromoform	1.0	U	1.0	0.19
79-34-5	1,1,2,2-Tetrachloroethane	1.0	U	1.0	0.20
107-13-1	Acrylonitrile	20	U	20	0.55
123-91-1	1,4-Dioxane	200	U	200	34

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	113		64-135
2037-26-5	Toluene-d8 (Surr)	101		71-118
460-00-4	4-Bromofluorobenzene (Surr)	93		70-118
1868-53-7	Dibromofluoromethane (Surr)	107		70-128

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415005.D  
 Lims ID: MB  
 Client ID:  
 Sample Type: MB  
 Inject. Date: 15-Apr-2015 14:33:30 ALS Bottle#: 5 Worklist Smp#: 5  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: MB  
 Misc. Info.: 180-0006480-005  
 Operator ID: 001562 Instrument ID: CHHP5  
 Method: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 07:30:04 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICAL File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 16-Apr-2015 07:30:04

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.305	4.302	0.003	0	174312	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.274	7.271	0.003	99	460612	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.358	10.361	-0.003	89	107539	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.682	12.679	0.003	97	150393	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.537	6.528	0.009	92	111889	50.0	53.4	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.902	6.899	0.003	0	156630	50.0	56.7	
\$ 7 Toluene-d8 (Surr)	98	8.922	8.919	0.003	93	434905	50.0	50.7	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.532	11.529	0.003	88	144274	50.0	46.7	
11 Dichlorodifluoromethane	85		1.619					ND	
12 Chloromethane	50		1.789					ND	
13 Vinyl chloride	62		1.917					ND	
14 Butadiene	39		1.959					ND	
15 Bromomethane	94		2.264					ND	
16 Chloroethane	64		2.416					ND	
17 Dichlorofluoromethane	67		2.671					ND	
18 Trichlorofluoromethane	101		2.726					ND	
19 Ethanol	45		3.018					ND	
20 Ethyl ether	59		3.091					ND	
21 Acrolein	56		3.261					ND	
22 1,1-Dichloroethene	96		3.395					ND	
23 1,1,2-Trichloro-1,2,2-trif	101		3.450					ND	
24 Acetone	43	3.526	3.499	0.027	69	1778		1.88	
25 Iodomethane	142		3.626					ND	
26 Carbon disulfide	76		3.675					ND	
27 Isopropyl alcohol	45		3.785					ND	
29 Acetonitrile	40		3.931					ND	
28 3-Chloro-1-propene	76		3.943					ND	
30 Methyl acetate	43		4.022					ND	
31 Methylene Chloride	84		4.143					ND	
32 2-Methyl-2-propanol	59		4.441					ND	
33 Acrylonitrile	53		4.551					ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
34 trans-1,2-Dichloroethene	96		4.563					ND	
35 Methyl tert-butyl ether	73		4.600					ND	
36 Hexane	57		4.983					ND	
37 1,1-Dichloroethane	63		5.165					ND	
38 Vinyl acetate	43		5.299					ND	
39 2-Chloro-1,3-butadiene	53		5.312					ND	
41 Isopropyl ether	45		5.324					ND	
40 Isopropyl ether TIC	45		5.409					ND	
42 Tert-butyl ethyl ether	59		5.799					ND	
44 2,2-Dichloropropane	77		5.926					ND	
45 cis-1,2-Dichloroethene	96		5.938					ND	
43 Tert-butyl ethyl ether (TI)	59		5.961					ND	
46 2-Butanone (MEK)	43		5.987					ND	
47 Propionitrile	54		6.060					ND	
48 Ethyl acetate	43		6.084					ND	
49 Chlorobromomethane	128		6.224					ND	
50 Methacrylonitrile	41		6.237					ND	
51 Tetrahydrofuran	42		6.279					ND	
52 Chloroform	83	6.343	6.339	0.004	1	961		0.2157	
53 1,1,1-Trichloroethane	97		6.528					ND	
54 Cyclohexane	56		6.583					ND	
56 Carbon tetrachloride	117		6.723					ND	
55 1,1-Dichloropropene	75		6.723					ND	
57 Isobutyl alcohol	41		6.942					ND	
58 Benzene	78		6.954					ND	
59 1,2-Dichloroethane	62		6.984					ND	
61 Tert-amyl methyl ether	73		7.113					ND	
60 Tert-amyl methyl ether (TI)	73		7.262					ND	
62 n-Heptane	43		7.276					ND	
63 n-Butanol	56		7.654					ND	
64 Trichloroethene	130		7.666					ND	
65 Ethyl acrylate	55		7.818					ND	
66 Methylcyclohexane	83		7.860					ND	
67 1,2-Dichloropropane	63		7.897					ND	
68 Dibromomethane	93		8.025					ND	
70 1,4-Dioxane	88		8.049					ND	
69 Methyl methacrylate	69		8.049					ND	
71 Dichlorobromomethane	83		8.195					ND	
72 2-Nitropropane	41		8.439					ND	
73 2-Chloroethyl vinyl ether	63		8.517					ND	
74 cis-1,3-Dichloropropene	75		8.651					ND	
75 4-Methyl-2-pentanone (MIBK)	43		8.822					ND	
76 Toluene	91		8.986					ND	
77 trans-1,3-Dichloropropene	75		9.217					ND	
78 Ethyl methacrylate	69		9.314					ND	
79 1,1,2-Trichloroethane	97		9.399					ND	
80 Tetrachloroethene	164		9.533					ND	
81 1,3-Dichloropropane	76		9.558					ND	
82 2-Hexanone	43		9.655					ND	
83 n-Butyl acetate	43		9.783					ND	
84 Chlorodibromomethane	129		9.789					ND	
85 Ethylene Dibromide	107		9.898					ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
86 3-Chlorobenzotrifluoride	180		10.373					ND	
87 Chlorobenzene	112		10.391					ND	
88 4-Chlorobenzotrifluoride	180		10.428					ND	
89 1,1,1,2-Tetrachloroethane	131		10.470					ND	
90 Ethylbenzene	106		10.501					ND	
91 m-Xylene & p-Xylene	106		10.616					ND	
92 o-Xylene	106		11.012					ND	
93 Styrene	104		11.024					ND	
94 Bromoform	173		11.212					ND	
95 Cyclohexanol	57		11.231					ND	
96 2-Chlorobenzotrifluoride	180		11.273					ND	
97 Isopropylbenzene	105		11.377					ND	
98 Cyclohexanone	55		11.481					ND	
99 1,1,2,2-Tetrachloroethane	83		11.675					ND	
100 Bromobenzene	156		11.681					ND	
101 1,2,3-Trichloropropane	110		11.717					ND	
102 trans-1,4-Dichloro-2-buten	53		11.729					ND	
103 N-Propylbenzene	120		11.784					ND	
104 2-Chlorotoluene	126		11.869					ND	
105 3-Chlorotoluene	126		11.930					ND	
106 1,3,5-Trimethylbenzene	105		11.961					ND	
107 4-Chlorotoluene	126		11.985					ND	
108 tert-Butylbenzene	119		12.283					ND	
109 Pentachloroethane	167		12.302					ND	
110 1,2,4-Trimethylbenzene	105		12.332					ND	
111 1,2-dichloro-4-(trifluoromethyl)	214		12.399					ND	
112 sec-Butylbenzene	105		12.502					ND	
113 1,3-Dichlorobenzene	146		12.618					ND	
114 4-Isopropyltoluene	119		12.648					ND	
115 1,4-Dichlorobenzene	146		12.703					ND	
116 2,4-Dichloro-1-(trifluoromethyl)	214		12.758					ND	
117 1,2,3-Trimethylbenzene	105		12.758					ND	
118 2,5-Dichlorobenzotrifluoride	214		12.806					ND	
119 Benzyl chloride	91		12.843					ND	
120 n-Butylbenzene	91		13.062					ND	
121 1,2-Dichlorobenzene	146		13.080					ND	
122 1,2-Dibromo-3-Chloropropan	75		13.853					ND	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125		14.005					ND	
124 1,3,5-Trichlorobenzene	180		14.072					ND	
125 2,3- & 3,4- Dichlorotoluene	125		14.424					ND	
126 1,2,4-Trichlorobenzene	180		14.686					ND	
127 Hexachlorobutadiene	225		14.856					ND	
128 Naphthalene	128		14.942					ND	
129 1,2,3-Trichlorobenzene	180		15.185					ND	
131 2,4,5-Trichlorotoluene	159		15.964					ND	
130 2,3,6-Trichlorotoluene	159		16.061					ND	
132 2-Methylnaphthalene	142		16.074					ND	
146 2,5-Dichlorotoluene	1		0.000					ND	
150 2,6-Dichlorotoluene	1		0.000					ND	
151 Isooctane	57		0.000					ND	
149 3,4-Dichlorotoluene	1		0.000					ND	
152 Formaldehyde TIC	1		0.000					ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
148 2,3-Dichlorotoluene	1		0.000						ND
147 2,4-Dichlorotoluene	1		0.000						ND
S 133 Xylenes, Total	106		1.000						ND
S 134 1,2-Dichloroethene, Total	96		1.000						ND
S 135 1,3-Dichloropropene, Total	1		0.000						ND
T 153 1,2 Epoxybutane TIC	42		0.000						ND
T 136 Mesityl oxide TIC	83		0.000						ND
T 137 Tetrahydrofuran TIC	42		0.000						ND
T 138 Methyl n-amyl ketone TIC	43		0.000						ND

**Reagents:**

VOA8260INT\_00031

Amount Added: 2.00

Units: uL

Run Reagent

VOA8260SURR\_00033

Amount Added: 2.00

Units: uL

Run Reagent

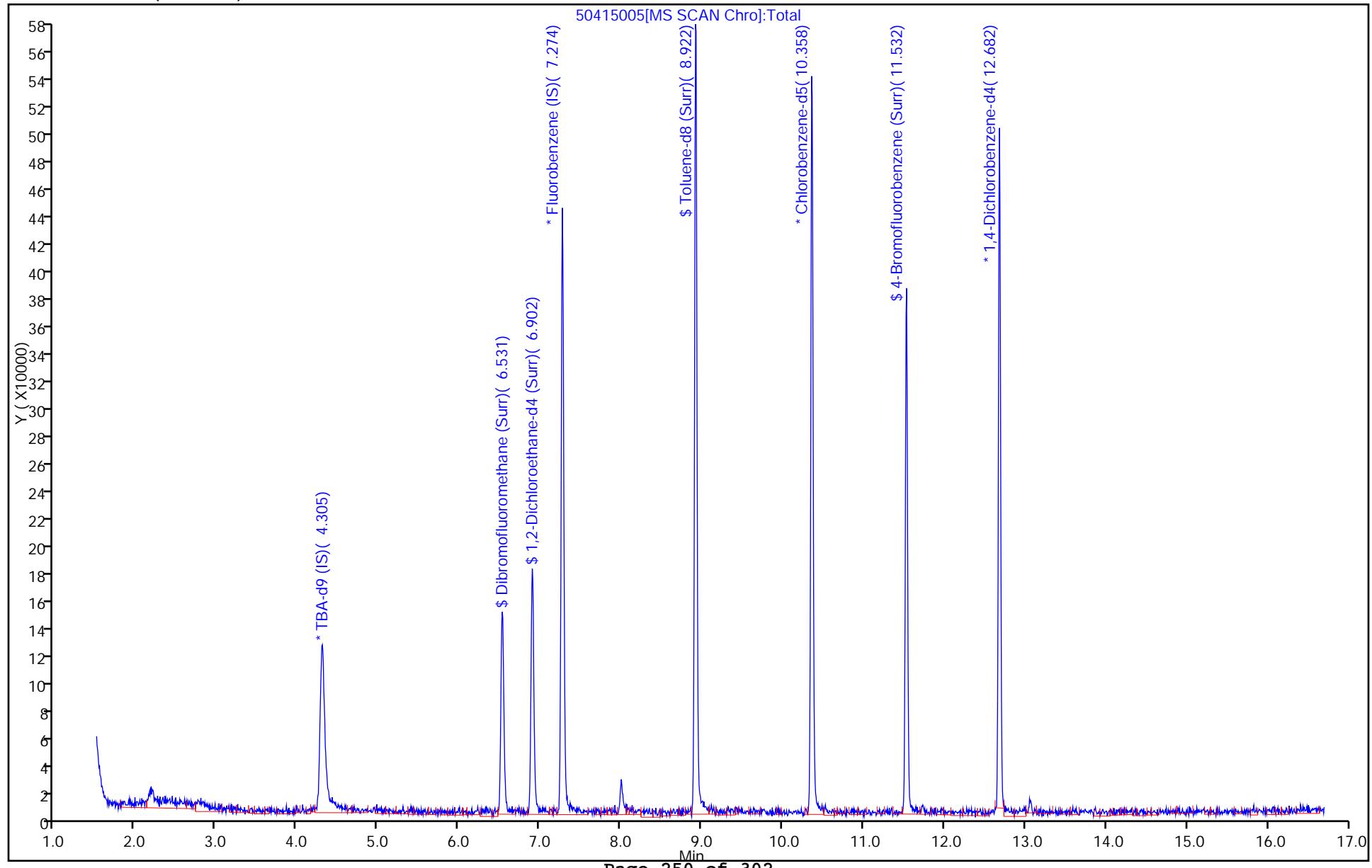
Report Date: 16-Apr-2015 07:30:05

Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Pittsburgh

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150415-6480.b\\50415005.D  
Injection Date: 15-Apr-2015 14:33:30 Instrument ID: CHHP5  
Lims ID: MB Operator ID: 001562  
Client ID:  
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 5  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)

Worklist Smp#: 5



FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: \_\_\_\_\_ Lab Sample ID: MB 180-138685/5  
Matrix: Water Lab File ID: 50416005.D  
Analysis Method: 8260C Date Collected: \_\_\_\_\_  
Sample wt/vol: 5 (mL) Date Analyzed: 04/16/2015 11:32  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138685 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
74-87-3	Chloromethane	1.0	U	1.0	0.28
75-01-4	Vinyl chloride	1.0	U	1.0	0.23
74-83-9	Bromomethane	1.0	U	1.0	0.31
75-00-3	Chloroethane	1.0	U	1.0	0.21
75-35-4	1,1-Dichloroethene	1.0	U	1.0	0.30
67-64-1	Acetone	5.0	U	5.0	2.5
75-15-0	Carbon disulfide	1.0	U	1.0	0.21
75-09-2	Methylene Chloride	0.162	J	1.0	0.13
156-60-5	trans-1,2-Dichloroethene	1.0	U	1.0	0.17
1634-04-4	Methyl tert-butyl ether	1.0	U	1.0	0.18
75-34-3	1,1-Dichloroethane	1.0	U	1.0	0.12
156-59-2	cis-1,2-Dichloroethene	1.0	U	1.0	0.24
74-97-5	Bromochloromethane	1.0	U	1.0	0.18
78-93-3	2-Butanone (MEK)	5.0	U	5.0	0.55
67-66-3	Chloroform	1.0	U	1.0	0.17
71-55-6	1,1,1-Trichloroethane	1.0	U	1.0	0.29
56-23-5	Carbon tetrachloride	1.0	U	1.0	0.14
71-43-2	Benzene	1.0	U	1.0	0.11
107-06-2	1,2-Dichloroethane	1.0	U	1.0	0.21
79-01-6	Trichloroethene	1.0	U	1.0	0.14
78-87-5	1,2-Dichloropropane	1.0	U	1.0	0.095
75-27-4	Bromodichloromethane	1.0	U	1.0	0.13
10061-01-5	cis-1,3-Dichloropropene	1.0	U	1.0	0.19
108-10-1	4-Methyl-2-pentanone (MIBK)	5.0	U	5.0	0.53
108-88-3	Toluene	1.0	U	1.0	0.15
10061-02-6	trans-1,3-Dichloropropene	1.0	U	1.0	0.15
79-00-5	1,1,2-Trichloroethane	1.0	U	1.0	0.20
127-18-4	Tetrachloroethene	1.0	U	1.0	0.15
591-78-6	2-Hexanone	5.0	U	5.0	0.16
124-48-1	Dibromochloromethane	1.0	U	1.0	0.14
106-93-4	1,2-Dibromoethane (EDB)	1.0	U	1.0	0.18
108-90-7	Chlorobenzene	1.0	U	1.0	0.14
630-20-6	1,1,1,2-Tetrachloroethane	1.0	U	1.0	0.28
100-41-4	Ethylbenzene	1.0	U	1.0	0.23
1330-20-7	Xylenes, Total	3.0	U	3.0	0.49
100-42-5	Styrene	1.0	U	1.0	0.097

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: \_\_\_\_\_ Lab Sample ID: MB 180-138685/5  
Matrix: Water Lab File ID: 50416005.D  
Analysis Method: 8260C Date Collected: \_\_\_\_\_  
Sample wt/vol: 5 (mL) Date Analyzed: 04/16/2015 11:32  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138685 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-25-2	Bromoform	1.0	U	1.0	0.19
79-34-5	1,1,2,2-Tetrachloroethane	1.0	U	1.0	0.20
107-13-1	Acrylonitrile	20	U	20	0.55
123-91-1	1,4-Dioxane	200	U	200	34

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	116		64-135
2037-26-5	Toluene-d8 (Surr)	100		71-118
460-00-4	4-Bromofluorobenzene (Surr)	97		70-118
1868-53-7	Dibromofluoromethane (Surr)	111		70-128

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\50416005.D  
 Lims ID: MB  
 Client ID:  
 Sample Type: MB  
 Inject. Date: 16-Apr-2015 11:32:30 ALS Bottle#: 5 Worklist Smp#: 5  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: MB  
 Misc. Info.: 180-0006494-005  
 Operator ID: 001562 Instrument ID: CHHP5  
 Method: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 12:37:09 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICAL File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 16-Apr-2015 11:51:47

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.303	4.317	-0.014	0	194715	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.278	7.268	0.010	98	448391	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.362	10.358	0.004	89	108906	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.686	12.682	0.004	97	157942	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.529	6.525	0.004	92	113002	50.0	55.4	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.907	6.897	0.010	0	156007	50.0	58.0	
\$ 7 Toluene-d8 (Surr)	98	8.926	8.916	0.010	95	434094	50.0	50.0	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.530	11.532	-0.002	87	151494	50.0	48.5	
11 Dichlorodifluoromethane	85		1.628					ND	
12 Chloromethane	50		1.786					ND	
13 Vinyl chloride	62		1.914					ND	
14 Butadiene	39	1.979	1.957	0.022	19	315		0.0930	
15 Bromomethane	94		2.267					ND	
16 Chloroethane	64		2.413					ND	
17 Dichlorofluoromethane	67		2.669					ND	
18 Trichlorofluoromethane	101		2.735					ND	
19 Ethanol	45		3.018					ND	
20 Ethyl ether	59		3.094					ND	
21 Acrolein	56		3.265					ND	
22 1,1-Dichloroethene	96		3.386					ND	
23 1,1,2-Trichloro-1,2,2-trif	101		3.441					ND	
24 Acetone	43		3.496					ND	
25 Iodomethane	142		3.618					ND	
26 Carbon disulfide	76		3.691					ND	
27 Isopropyl alcohol	45		3.785					ND	
29 Acetonitrile	40		3.931					ND	
28 3-Chloro-1-propene	76		3.946					ND	
30 Methyl acetate	43		4.025					ND	
31 Methylene Chloride	84	4.175	4.147	0.028	1	2426		0.8113	M
32 2-Methyl-2-propanol	59		4.439					ND	
33 Acrylonitrile	53		4.548					ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
34 trans-1,2-Dichloroethene	96		4.561					ND	
35 Methyl tert-butyl ether	73		4.591					ND	
36 Hexane	57		4.986					ND	
37 1,1-Dichloroethane	63		5.175					ND	
38 Vinyl acetate	43		5.291					ND	
39 2-Chloro-1,3-butadiene	53		5.312					ND	
41 Isopropyl ether	45		5.324					ND	
40 Isopropyl ether TIC	45		5.409					ND	
42 Tert-butyl ethyl ether	59		5.799					ND	
44 2,2-Dichloropropane	77		5.929					ND	
45 cis-1,2-Dichloroethene	96		5.935					ND	
43 Tert-butyl ethyl ether (TI)	59		5.961					ND	
46 2-Butanone (MEK)	43		5.990					ND	
47 Propionitrile	54		6.060					ND	
48 Ethyl acetate	43	6.037	6.084	-0.047	1	122		0.0894	
49 Chlorobromomethane	128		6.227					ND	
50 Methacrylonitrile	41		6.237					ND	
51 Tetrahydrofuran	42		6.282					ND	
52 Chloroform	83		6.337					ND	
53 1,1,1-Trichloroethane	97		6.525					ND	
54 Cyclohexane	56		6.580					ND	
56 Carbon tetrachloride	117		6.720					ND	
55 1,1-Dichloropropene	75		6.720					ND	
57 Isobutyl alcohol	41		6.945					ND	
58 Benzene	78		6.951					ND	
59 1,2-Dichloroethane	62		6.982					ND	
61 Tert-amyl methyl ether	73		7.113					ND	
60 Tert-amyl methyl ether (TI)	73		7.262					ND	
62 n-Heptane	43	7.241	7.280	-0.039	1	78		0.0214	
63 n-Butanol	56		7.654					ND	
64 Trichloroethene	130		7.669					ND	
65 Ethyl acrylate	55	7.801	7.818	-0.017	1	161		0.0575	
66 Methylcyclohexane	83		7.858					ND	
67 1,2-Dichloropropane	63		7.900					ND	
68 Dibromomethane	93		8.022					ND	
69 Methyl methacrylate	69		8.049					ND	
70 1,4-Dioxane	88		8.052					ND	
71 Dichlorobromomethane	83		8.192					ND	
72 2-Nitropropane	41		8.439					ND	
73 2-Chloroethyl vinyl ether	63		8.515					ND	
74 cis-1,3-Dichloropropene	75		8.655					ND	
75 4-Methyl-2-pentanone (MIBK)	43		8.819					ND	
76 Toluene	91		8.989					ND	
77 trans-1,3-Dichloropropene	75		9.220					ND	
78 Ethyl methacrylate	69		9.318					ND	
79 1,1,2-Trichloroethane	97		9.397					ND	
80 Tetrachloroethene	164		9.531					ND	
81 1,3-Dichloropropane	76		9.561					ND	
82 2-Hexanone	43		9.652					ND	
83 n-Butyl acetate	43		9.783					ND	
84 Chlorodibromomethane	129		9.786					ND	
85 Ethylene Dibromide	107		9.902					ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
86 3-Chlorobenzotrifluoride	180		10.370					ND	
87 Chlorobenzene	112		10.388					ND	
88 4-Chlorobenzotrifluoride	180		10.425					ND	
89 1,1,1,2-Tetrachloroethane	131		10.474					ND	
90 Ethylbenzene	106		10.498					ND	
91 m-Xylene & p-Xylene	106		10.614					ND	
92 o-Xylene	106		11.009					ND	
93 Styrene	104		11.021					ND	
94 Bromoform	173		11.204					ND	
95 Cyclohexanol	57		11.231					ND	
96 2-Chlorobenzotrifluoride	180		11.271					ND	
97 Isopropylbenzene	105		11.380					ND	
98 Cyclohexanone	55		11.481					ND	
99 1,1,2,2-Tetrachloroethane	83		11.672					ND	
100 Bromobenzene	156		11.678					ND	
101 1,2,3-Trichloropropane	110		11.721					ND	
102 trans-1,4-Dichloro-2-buten	53		11.727					ND	
103 N-Propylbenzene	120		11.788					ND	
104 2-Chlorotoluene	126		11.873					ND	
105 3-Chlorotoluene	126		11.934					ND	
106 1,3,5-Trimethylbenzene	105		11.958					ND	
107 4-Chlorotoluene	126		11.976					ND	
108 tert-Butylbenzene	119		12.287					ND	
109 Pentachloroethane	167		12.302					ND	
110 1,2,4-Trimethylbenzene	105		12.335					ND	
111 1,2-dichloro-4-(trifluoromethyl)	214		12.402					ND	
112 sec-Butylbenzene	105		12.506					ND	
113 1,3-Dichlorobenzene	146		12.615					ND	
114 4-Isopropyltoluene	119		12.645					ND	
115 1,4-Dichlorobenzene	146		12.706					ND	
116 2,4-Dichloro-1-(trifluoromethyl)	214		12.755					ND	
117 1,2,3-Trimethylbenzene	105		12.758					ND	
118 2,5-Dichlorobenzotrifluoride	214		12.804					ND	
119 Benzyl chloride	91		12.843					ND	
120 n-Butylbenzene	91		13.059					ND	
121 1,2-Dichlorobenzene	146		13.077					ND	
122 1,2-Dibromo-3-Chloropropan	75		13.862					ND	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125		14.008					ND	
124 1,3,5-Trichlorobenzene	180		14.069					ND	
125 2,3- & 3,4- Dichlorotoluene	125		14.422					ND	
126 1,2,4-Trichlorobenzene	180		14.690					ND	
127 Hexachlorobutadiene	225		14.860					ND	
128 Naphthalene	128		14.939					ND	
129 1,2,3-Trichlorobenzene	180		15.182					ND	
131 2,4,5-Trichlorotoluene	159		15.961					ND	
130 2,3,6-Trichlorotoluene	159		16.058					ND	
132 2-Methylnaphthalene	142		16.074					ND	
148 2,3-Dichlorotoluene	1		0.000					ND	
147 2,4-Dichlorotoluene	1		0.000					ND	
146 2,5-Dichlorotoluene	1		0.000					ND	
150 2,6-Dichlorotoluene	1		0.000					ND	
152 Formaldehyde TIC	1		0.000					ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
151 Isooctane	57		0.000						ND
149 3,4-Dichlorotoluene	1		0.000						ND
S 133 Xylenes, Total	106		1.000						ND
S 134 1,2-Dichloroethene, Total	96		1.000						ND
S 135 1,3-Dichloropropene, Total	1		0.000						ND
T 153 1,2 Epoxybutane TIC	42		0.000						ND
T 136 Mesityl oxide TIC	83		0.000						ND
T 137 Tetrahydrofuran TIC	42		0.000						ND
T 138 Methyl n-amyl ketone TIC	43		0.000						ND

**QC Flag Legend**

Review Flags

M - Manually Integrated

**Reagents:**

VOA8260INT\_00031  
 VOA8260SURR\_00033

Amount Added: 2.00 Units: uL Run Reagent  
 Amount Added: 2.00 Units: uL Run Reagent

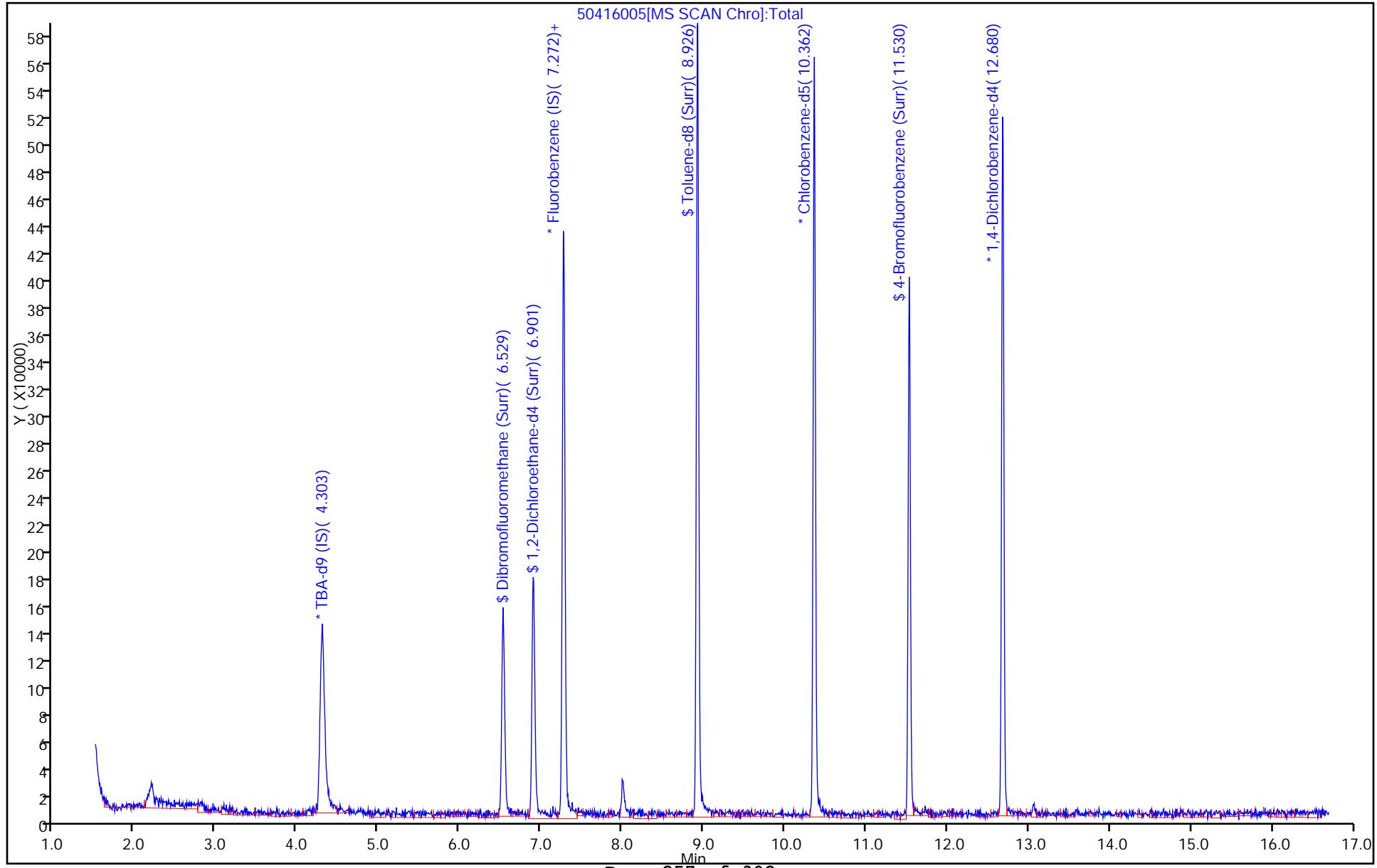
Report Date: 16-Apr-2015 12:37:14

Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Pittsburgh

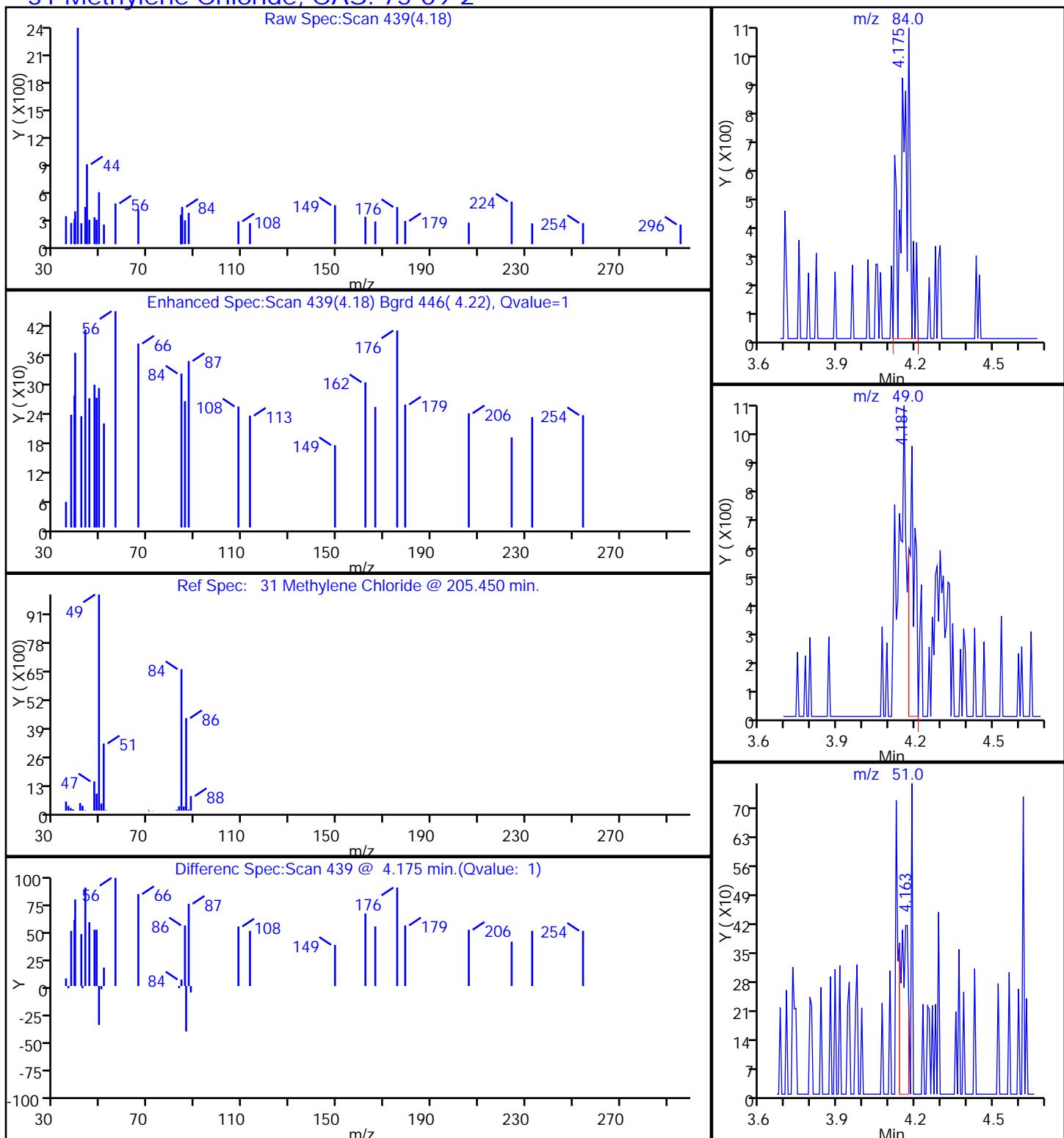
Data File: \\PITCHROM\\ChromData\\CHHP5\\20150416-6494.b\\50416005.D  
Injection Date: 16-Apr-2015 11:32:30 Instrument ID: CHHP5  
Lims ID: MB Operator ID: 001562  
Client ID:  
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 5  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)

Worklist Smp#: 5



TestAmerica Pittsburgh  
 Data File: \PITCHROM\ChromData\CHHP5\20150416-6494.b\50416005.D  
 Injection Date: 16-Apr-2015 11:32:30 Instrument ID: CHHP5  
 Lims ID: MB  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 5 Worklist Smp#: 5  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 ( 0.18 mm) Detector: MS SCAN

### 31 Methylene Chloride, CAS: 75-09-2



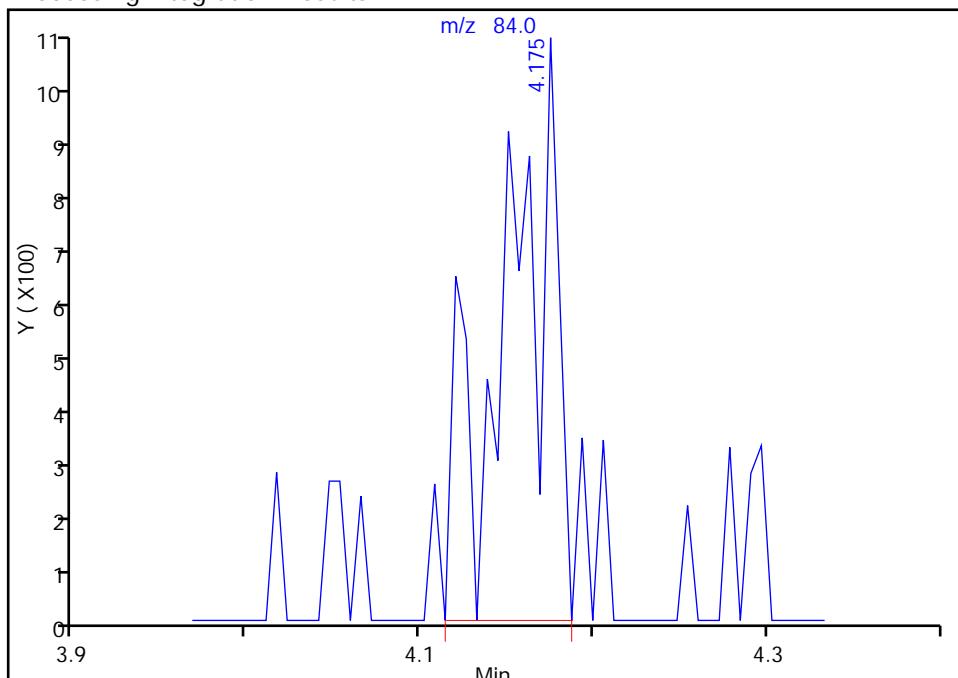
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\50416005.D  
 Injection Date: 16-Apr-2015 11:32:30 Instrument ID: CHHP5  
 Lims ID: MB  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 5 Worklist Smp#: 5  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 31 Methylene Chloride, CAS: 75-09-2

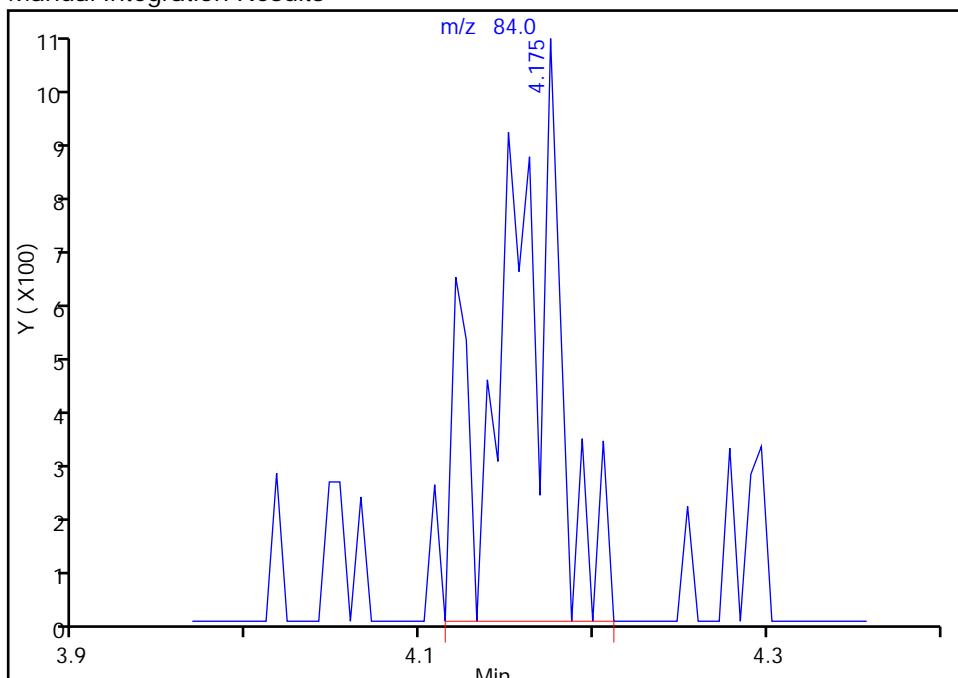
RT: 4.18  
 Area: 2187  
 Amount: 0.731357  
 Amount Units: ng

## Processing Integration Results



RT: 4.18  
 Area: 2426  
 Amount: 0.811281  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 16-Apr-2015 11:51:47

Audit Action: Manually Integrated

Audit Reason: Split Peak

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: \_\_\_\_\_ Lab Sample ID: LCS 180-138583/8  
Matrix: Water Lab File ID: 50415008.D  
Analysis Method: 8260C Date Collected: \_\_\_\_\_  
Sample wt/vol: 5 (mL) Date Analyzed: 04/15/2015 15:57  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138583 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
74-87-3	Chloromethane	10.1		1.0	0.28
75-01-4	Vinyl chloride	10.0		1.0	0.23
74-83-9	Bromomethane	8.98		1.0	0.31
75-00-3	Chloroethane	10.5		1.0	0.21
75-35-4	1,1-Dichloroethene	7.04		1.0	0.30
67-64-1	Acetone	19.0		5.0	2.5
75-15-0	Carbon disulfide	3.87		1.0	0.21
75-09-2	Methylene Chloride	7.62		1.0	0.13
156-60-5	trans-1,2-Dichloroethene	7.74		1.0	0.17
1634-04-4	Methyl tert-butyl ether	7.94		1.0	0.18
75-34-3	1,1-Dichloroethane	8.31		1.0	0.12
156-59-2	cis-1,2-Dichloroethene	7.88		1.0	0.24
74-97-5	Bromochloromethane	8.32		1.0	0.18
78-93-3	2-Butanone (MEK)	16.7		5.0	0.55
67-66-3	Chloroform	9.11		1.0	0.17
71-55-6	1,1,1-Trichloroethane	8.87		1.0	0.29
56-23-5	Carbon tetrachloride	9.33		1.0	0.14
71-43-2	Benzene	8.86		1.0	0.11
107-06-2	1,2-Dichloroethane	8.84		1.0	0.21
79-01-6	Trichloroethene	8.40		1.0	0.14
78-87-5	1,2-Dichloroproppane	9.27		1.0	0.095
75-27-4	Bromodichloromethane	8.85		1.0	0.13
10061-01-5	cis-1,3-Dichloropropene	9.60		1.0	0.19
108-10-1	4-Methyl-2-pentanone (MIBK)	16.8		5.0	0.53
108-88-3	Toluene	9.81		1.0	0.15
10061-02-6	trans-1,3-Dichloropropene	11.6		1.0	0.15
79-00-5	1,1,2-Trichloroethane	11.2		1.0	0.20
127-18-4	Tetrachloroethene	9.33		1.0	0.15
591-78-6	2-Hexanone	15.3		5.0	0.16
124-48-1	Dibromochloromethane	10.1		1.0	0.14
106-93-4	1,2-Dibromoethane (EDB)	9.99		1.0	0.18
108-90-7	Chlorobenzene	10.2		1.0	0.14
630-20-6	1,1,1,2-Tetrachloroethane	11.8		1.0	0.28
100-41-4	Ethylbenzene	9.72		1.0	0.23
1330-20-7	Xylenes, Total	19.0		3.0	0.49
100-42-5	Styrene	9.70		1.0	0.097

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: \_\_\_\_\_ Lab Sample ID: LCS 180-138583/8  
Matrix: Water Lab File ID: 50415008.D  
Analysis Method: 8260C Date Collected: \_\_\_\_\_  
Sample wt/vol: 5 (mL) Date Analyzed: 04/15/2015 15:57  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138583 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-25-2	Bromoform	9.35		1.0	0.19
79-34-5	1,1,2,2-Tetrachloroethane	10.9		1.0	0.20
107-13-1	Acrylonitrile	87.6		20	0.55
123-91-1	1,4-Dioxane	179	J	200	34

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	95		64-135
2037-26-5	Toluene-d8 (Surr)	99		71-118
460-00-4	4-Bromofluorobenzene (Surr)	94		70-118
1868-53-7	Dibromofluoromethane (Surr)	91		70-128

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415008.D  
 Lims ID: LCS  
 Client ID:  
 Sample Type: LCS  
 Inject. Date: 15-Apr-2015 15:57:30 ALS Bottle#: 8 Worklist Smp#: 8  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: LCS  
 Misc. Info.: 180-0006480-008  
 Operator ID: 001562 Instrument ID: CHHP5  
 Method: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 07:56:19 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICAL File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 16-Apr-2015 07:34:37

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.317	4.302	0.015	0	163019	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.273	7.271	0.002	98	552275	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.358	10.361	-0.003	88	123275	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.682	12.679	0.003	96	185676	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.525	6.528	-0.003	87	114892	50.0	45.7	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.896	6.899	-0.003	0	156966	50.0	47.4	
\$ 7 Toluene-d8 (Surr)	98	8.922	8.919	0.003	94	484341	50.0	49.3	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.526	11.529	-0.003	90	166121	50.0	46.9	
11 Dichlorodifluoromethane	85	1.616	1.619	-0.003	99	151519	50.0	64.0	
12 Chloromethane	50	1.786	1.789	-0.003	98	165254	50.0	50.6	
13 Vinyl chloride	62	1.908	1.917	-0.009	99	182699	50.0	50.0	
14 Butadiene	39	1.963	1.959	0.004	96	185184	50.0	44.4	
15 Bromomethane	94	2.267	2.264	0.003	90	89652	50.0	44.9	
16 Chloroethane	64	2.394	2.416	-0.022	99	132184	50.0	52.3	
17 Dichlorofluoromethane	67	2.662	2.671	-0.009	97	294079	50.0	51.0	
18 Trichlorofluoromethane	101	2.699	2.726	-0.027	90	196082	50.0	44.8	
20 Ethyl ether	59	3.094	3.091	0.003	93	115394	50.0	39.9	
21 Acrolein	56	3.264	3.261	0.003	99	47654	150.0	135.8	
22 1,1-Dichloroethene	96	3.380	3.395	-0.015	97	112036	50.0	35.2	
23 1,1,2-Trichloro-1,2,2-trif	101	3.453	3.450	0.003	92	121517	50.0	37.7	
24 Acetone	43	3.502	3.499	0.003	100	107684	100.0	95.2	
25 Iodomethane	142	3.611	3.626	-0.015	98	157366	50.0	35.6	
26 Carbon disulfide	76	3.672	3.675	-0.003	100	150609	50.0	19.3	
28 3-Chloro-1-propene	76	3.940	3.943	-0.003	87	52088	50.0	30.9	
30 Methyl acetate	43	4.025	4.022	0.003	98	593539	250.0	224.2	
31 Methylene Chloride	84	4.153	4.143	0.010	93	140387	50.0	38.1	
32 2-Methyl-2-propanol	59	4.445	4.441	0.004	84	102153	500.0	532.0	
33 Acrylonitrile	53	4.554	4.551	0.003	97	596596	500.0	438.2	
34 trans-1,2-Dichloroethene	96	4.560	4.563	-0.003	98	127451	50.0	38.7	
35 Methyl tert-butyl ether	73	4.597	4.600	-0.003	95	289287	50.0	39.7	
36 Hexane	57	4.986	4.983	0.003	93	163398	50.0	31.0	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
37 1,1-Dichloroethane	63	5.175	5.165	0.010	97	244205	50.0	41.5	
38 Vinyl acetate	43	5.290	5.299	-0.009	97	200662	50.0	48.1	
44 2,2-Dichloropropane	77	5.923	5.926	-0.003	60	73522	50.0	50.0	
45 cis-1,2-Dichloroethene	96	5.941	5.938	0.003	84	136676	50.0	39.4	
46 2-Butanone (MEK)	43	5.990	5.987	0.003	100	150748	100.0	83.3	
49 Chlorobromomethane	128	6.221	6.224	-0.003	96	62457	50.0	41.6	
51 Tetrahydrofuran	42	6.294	6.279	0.015	88	87660	100.0	77.4	
52 Chloroform	83	6.343	6.339	0.004	93	243265	50.0	45.5	
53 1,1,1-Trichloroethane	97	6.525	6.528	-0.003	96	151325	50.0	44.4	
54 Cyclohexane	56	6.592	6.583	0.009	93	213523	50.0	32.6	
56 Carbon tetrachloride	117	6.720	6.723	-0.003	70	127699	50.0	46.7	
55 1,1-Dichloropropene	75	6.726	6.723	0.003	93	179419	50.0	40.5	
57 Isobutyl alcohol	41	6.945	6.942	0.003	43	113809	1250.0	1543.9	
58 Benzene	78	6.957	6.954	0.003	97	579590	50.0	44.3	
59 1,2-Dichloroethane	62	6.987	6.984	0.003	97	189332	50.0	44.2	
62 n-Heptane	43	7.279	7.276	0.003	88	152111	50.0	33.8	
64 Trichloroethene	130	7.663	7.666	-0.003	96	137761	50.0	42.0	
66 Methylcyclohexane	83	7.863	7.860	0.003	91	205457	50.0	35.1	
67 1,2-Dichloropropane	63	7.900	7.897	0.003	95	150044	50.0	46.3	
68 Dibromomethane	93	8.022	8.025	-0.003	96	78270	50.0	44.9	
70 1,4-Dioxane	88	8.052	8.049	0.003	95	30531	1000.0	895.7	
71 Dichlorobromomethane	83	8.198	8.195	0.003	98	157314	50.0	44.2	
73 2-Chloroethyl vinyl ether	63	8.514	8.517	-0.003	93	160898	100.0	88.2	
74 cis-1,3-Dichloropropene	75	8.654	8.651	0.003	93	164775	50.0	48.0	
75 4-Methyl-2-pentanone (MIBK)	43	8.819	8.822	-0.003	98	280386	100.0	84.1	
76 Toluene	91	8.989	8.986	0.003	98	619625	50.0	49.0	
77 trans-1,3-Dichloropropene	75	9.220	9.217	0.003	97	131982	50.0	57.8	
78 Ethyl methacrylate	69	9.311	9.314	-0.003	91	144170	50.0	48.4	
79 1,1,2-Trichloroethane	97	9.403	9.399	0.004	93	132086	50.0	55.8	
80 Tetrachloroethene	164	9.536	9.533	0.003	96	115233	50.0	46.6	
81 1,3-Dichloropropane	76	9.561	9.558	0.003	94	233746	50.0	53.1	
82 2-Hexanone	43	9.652	9.655	-0.003	98	194993	100.0	76.5	
84 Chlorodibromomethane	129	9.786	9.789	-0.003	89	95576	50.0	50.5	
85 Ethylene Dibromide	107	9.895	9.898	-0.003	98	112949	50.0	50.0	
86 3-Chlorobenzotrifluoride	180	10.370	10.373	-0.003	94	233331	50.0	48.4	
87 Chlorobenzene	112	10.388	10.391	-0.003	94	409960	50.0	51.2	
88 4-Chlorobenzotrifluoride	180	10.425	10.428	-0.003	94	221035	50.0	47.4	
89 1,1,1,2-Tetrachloroethane	131	10.473	10.470	0.003	91	121921	50.0	59.0	
90 Ethylbenzene	106	10.498	10.501	-0.003	99	223153	50.0	48.6	
91 m-Xylene & p-Xylene	106	10.619	10.616	0.003	0	270181	50.0	48.1	
92 o-Xylene	106	11.009	11.012	-0.003	96	256327	50.0	46.7	
93 Styrene	104	11.021	11.024	-0.003	94	429342	50.0	48.5	
94 Bromoform	173	11.209	11.212	-0.003	96	54595	50.0	46.7	
96 2-Chlorobenzotrifluoride	180	11.270	11.273	-0.003	97	224424	50.0	46.6	
97 Isopropylbenzene	105	11.380	11.377	0.003	97	647124	50.0	47.2	
99 1,1,2,2-Tetrachloroethane	83	11.672	11.675	-0.003	95	185129	50.0	54.5	
100 Bromobenzene	156	11.678	11.681	-0.003	94	165756	50.0	48.2	
101 1,2,3-Trichloropropane	110	11.714	11.717	-0.003	86	58741	50.0	52.0	
102 trans-1,4-Dichloro-2-butene	53	11.726	11.729	-0.003	65	46297	50.0	49.3	
103 N-Propylbenzene	120	11.787	11.784	0.003	99	191069	50.0	45.1	
104 2-Chlorotoluene	126	11.872	11.869	0.003	96	162769	50.0	45.7	
105 3-Chlorotoluene	126	11.933	11.930	0.003	98	179619	50.0	45.1	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
106 1,3,5-Trimethylbenzene	105	11.958	11.961	-0.003	94	556049	50.0	47.0	
107 4-Chlorotoluene	126	11.982	11.985	-0.003	98	184731	50.0	47.9	
108 tert-Butylbenzene	119	12.286	12.283	0.003	93	428757	50.0	41.9	
110 1,2,4-Trimethylbenzene	105	12.335	12.332	0.003	97	547006	50.0	45.1	
111 1,2-dichloro-4-(trifluoromethyl)	214	12.402	12.399	0.003	99	167924	50.0	43.8	
112 sec-Butylbenzene	105	12.505	12.502	0.003	95	642917	50.0	44.6	
113 1,3-Dichlorobenzene	146	12.615	12.618	-0.003	97	298086	50.0	47.1	
114 4-Isopropyltoluene	119	12.651	12.648	0.003	96	523673	50.0	44.0	
115 1,4-Dichlorobenzene	146	12.706	12.703	0.003	94	316006	50.0	48.9	
116 2,4-Dichloro-1-(trifluoromethyl)	214	12.755	12.758	-0.003	99	152027	50.0	42.3	
118 2,5-Dichlorobenzotrifluoride	214	12.803	12.806	-0.003	0	173559	50.0	43.2	
120 n-Butylbenzene	91	13.059	13.062	-0.003	98	454579	50.0	42.0	
121 1,2-Dichlorobenzene	146	13.077	13.080	-0.003	96	279566	50.0	47.7	
122 1,2-Dibromo-3-Chloropropan	75	13.856	13.853	0.003	80	20632	50.0	43.0	
123 2,4- & 2,5- & 2,6- Dichloro	125	14.008	14.005	0.003	0	501013	150.0	113.0	
124 1,3,5-Trichlorobenzene	180	14.069	14.072	-0.003	97	165012	50.0	46.4	
125 2,3- & 3,4- Dichlorotoluene	125	14.421	14.424	-0.003	0	312659	100.0	72.5	
126 1,2,4-Trichlorobenzene	180	14.689	14.686	0.003	94	116847	50.0	38.3	
127 Hexachlorobutadiene	225	14.859	14.856	0.003	96	61409	50.0	42.0	
128 Naphthalene	128	14.939	14.942	-0.003	97	267486	50.0	33.4	
129 1,2,3-Trichlorobenzene	180	15.182	15.185	-0.003	94	91405	50.0	36.5	
131 2,4,5-Trichlorotoluene	159	15.954	15.964	-0.010	0	31131	50.0	23.1	
130 2,3,6-Trichlorotoluene	159	16.058	16.061	-0.003	93	29773	50.0	24.5	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
S 133 Xylenes, Total	106				0		100.0	94.8	
S 134 1,2-Dichloroethene, Total	96				0		100.0	78.1	
S 135 1,3-Dichloropropene, Total	1				0		100.0	105.9	

**QC Flag Legend**

Processing Flags

ND - Not Detected or Marked ND

**Reagents:**

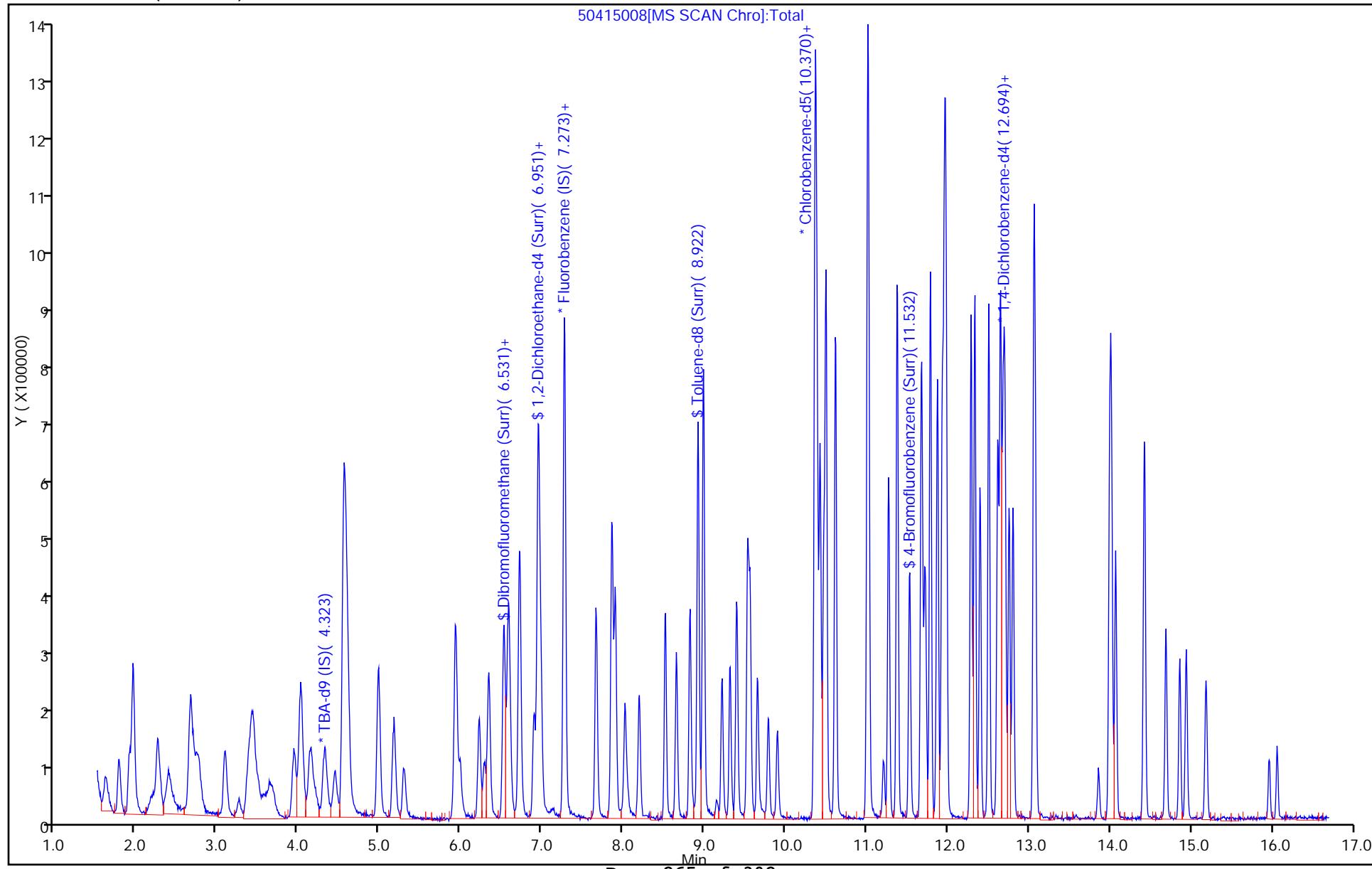
VOA8260VOA2ND_00111	Amount Added: 2.00	Units: uL	
voaWKet2 Rest_00002	Amount Added: 2.00	Units: uL	
voaW ee2nd_00001	Amount Added: 2.00	Units: uL	
voaWVA2nd Res_00006	Amount Added: 2.00	Units: uL	
voaW 135tcb a_00001	Amount Added: 2.00	Units: uL	
voaW2-cl 2ndR_00002	Amount Added: 2.00	Units: uL	
VOAACRO2ND_00007	Amount Added: 6.00	Units: uL	
VOA8260INT_00031	Amount Added: 2.00	Units: uL	Run Reagent
VOA8260SURR_00033	Amount Added: 2.00	Units: uL	Run Reagent

Report Date: 16-Apr-2015 08:48:32

Chrom Revision: 2.2 13-Mar-2015 11:20:44

## TestAmerica Pittsburgh

Data File: \PITCHROM\ChromData\CHHP5\20150415-6480.b\50415008.D  
Injection Date: 15-Apr-2015 15:57:30 Instrument ID: CHHP5  
Lims ID: LCS Operator ID: 001562  
Client ID:  
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 8  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)



FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: \_\_\_\_\_ Lab Sample ID: LCS 180-138685/8  
Matrix: Water Lab File ID: 50416008.D  
Analysis Method: 8260C Date Collected: \_\_\_\_\_  
Sample wt/vol: 5 (mL) Date Analyzed: 04/16/2015 13:05  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138685 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
74-87-3	Chloromethane	10.0		1.0	0.28
75-01-4	Vinyl chloride	10.5		1.0	0.23
74-83-9	Bromomethane	9.71		1.0	0.31
75-00-3	Chloroethane	11.0		1.0	0.21
75-35-4	1,1-Dichloroethene	9.09		1.0	0.30
67-64-1	Acetone	21.0		5.0	2.5
75-15-0	Carbon disulfide	6.75		1.0	0.21
75-09-2	Methylene Chloride	9.92		1.0	0.13
156-60-5	trans-1,2-Dichloroethene	9.42		1.0	0.17
1634-04-4	Methyl tert-butyl ether	9.22		1.0	0.18
75-34-3	1,1-Dichloroethane	9.66		1.0	0.12
156-59-2	cis-1,2-Dichloroethene	9.13		1.0	0.24
74-97-5	Bromochloromethane	9.67		1.0	0.18
78-93-3	2-Butanone (MEK)	17.2		5.0	0.55
67-66-3	Chloroform	10.1		1.0	0.17
71-55-6	1,1,1-Trichloroethane	10.6		1.0	0.29
56-23-5	Carbon tetrachloride	11.0		1.0	0.14
71-43-2	Benzene	10.2		1.0	0.11
107-06-2	1,2-Dichloroethane	10.0		1.0	0.21
79-01-6	Trichloroethene	9.34		1.0	0.14
78-87-5	1,2-Dichloroproppane	9.73		1.0	0.095
75-27-4	Bromodichloromethane	9.17		1.0	0.13
10061-01-5	cis-1,3-Dichloropropene	9.80		1.0	0.19
108-10-1	4-Methyl-2-pentanone (MIBK)	16.2		5.0	0.53
108-88-3	Toluene	10.5		1.0	0.15
10061-02-6	trans-1,3-Dichloropropene	10.8		1.0	0.15
79-00-5	1,1,2-Trichloroethane	10.7		1.0	0.20
127-18-4	Tetrachloroethene	10.2		1.0	0.15
591-78-6	2-Hexanone	16.8		5.0	0.16
124-48-1	Dibromochloromethane	9.51		1.0	0.14
106-93-4	1,2-Dibromoethane (EDB)	10.2		1.0	0.18
108-90-7	Chlorobenzene	10.3		1.0	0.14
630-20-6	1,1,1,2-Tetrachloroethane	11.1		1.0	0.28
100-41-4	Ethylbenzene	9.73		1.0	0.23
1330-20-7	Xylenes, Total	19.8		3.0	0.49
100-42-5	Styrene	9.71		1.0	0.097

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: \_\_\_\_\_ Lab Sample ID: LCS 180-138685/8  
Matrix: Water Lab File ID: 50416008.D  
Analysis Method: 8260C Date Collected: \_\_\_\_\_  
Sample wt/vol: 5 (mL) Date Analyzed: 04/16/2015 13:05  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138685 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-25-2	Bromoform	9.19		1.0	0.19
79-34-5	1,1,2,2-Tetrachloroethane	10.0		1.0	0.20
107-13-1	Acrylonitrile	95.8		20	0.55
123-91-1	1,4-Dioxane	167	J	200	34

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	93		64-135
2037-26-5	Toluene-d8 (Surr)	101		71-118
460-00-4	4-Bromofluorobenzene (Surr)	94		70-118
1868-53-7	Dibromofluoromethane (Surr)	93		70-128

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\50416008.D  
 Lims ID: LCS  
 Client ID:  
 Sample Type: LCS  
 Inject. Date: 16-Apr-2015 13:05:30 ALS Bottle#: 8 Worklist Smp#: 8  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: LCS  
 Misc. Info.: 180-0006494-008  
 Operator ID: 001562 Instrument ID: CHHP5  
 Method: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 13:57:55 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICAL File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 16-Apr-2015 13:52:52

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.316	4.317	-0.001	0	158617	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.273	7.268	0.005	98	554542	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.363	10.358	0.005	88	125698	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.681	12.682	-0.001	96	186690	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.531	6.525	0.006	80	117647	50.0	46.6	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.902	6.897	0.005	0	154924	50.0	46.6	
\$ 7 Toluene-d8 (Surr)	98	8.921	8.916	0.005	93	507014	50.0	50.6	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.531	11.532	-0.001	89	170279	50.0	47.2	
11 Dichlorodifluoromethane	85	1.615	1.628	-0.013	98	142573	50.0	60.0	
12 Chloromethane	50	1.779	1.786	-0.007	99	164484	50.0	50.1	
13 Vinyl chloride	62	1.913	1.914	-0.001	99	192109	50.0	52.4	
14 Butadiene	39	1.956	1.957	-0.001	96	207211	50.0	49.5	
15 Bromomethane	94	2.254	2.267	-0.013	92	96674	50.0	48.6	
16 Chloroethane	64	2.394	2.413	-0.019	99	139998	50.0	55.2	
17 Dichlorofluoromethane	67	2.662	2.669	-0.007	98	309688	50.0	53.5	
18 Trichlorofluoromethane	101	2.692	2.735	-0.043	97	206762	50.0	47.0	
20 Ethyl ether	59	3.087	3.094	-0.007	92	144771	50.0	49.9	
21 Acrolein	56	3.264	3.265	-0.001	96	43469	150.0	123.4	
22 1,1-Dichloroethene	96	3.386	3.386	0.000	97	145280	50.0	45.4	
23 1,1,2-Trichloro-1,2,2-trif	101	3.440	3.441	-0.001	94	163569	50.0	50.6	
24 Acetone	43	3.495	3.496	-0.001	98	119501	100.0	105.2	
25 Iodomethane	142	3.611	3.618	-0.007	99	203913	50.0	45.9	
26 Carbon disulfide	76	3.684	3.691	-0.007	99	264084	50.0	33.8	
28 3-Chloro-1-propene	76	3.945	3.946	-0.001	88	71568	50.0	42.3	
30 Methyl acetate	43	4.018	4.025	-0.007	98	635208	250.0	239.0	
31 Methylene Chloride	84	4.158	4.147	0.011	95	183520	50.0	49.6	
32 2-Methyl-2-propanol	59	4.444	4.439	0.005	85	91254	500.0	488.4	
33 Acrylonitrile	53	4.554	4.548	0.006	99	655138	500.0	479.2	
34 trans-1,2-Dichloroethene	96	4.566	4.561	0.005	62	155816	50.0	47.1	
35 Methyl tert-butyl ether	73	4.596	4.591	0.005	95	337215	50.0	46.1	
36 Hexane	57	4.985	4.986	-0.001	95	218992	50.0	41.4	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
37 1,1-Dichloroethane	63	5.168	5.175	-0.007	96	285135	50.0	48.3	
38 Vinyl acetate	43	5.296	5.291	0.005	97	164953	50.0	39.4	
44 2,2-Dichloropropane	77	5.922	5.929	-0.007	68	91592	50.0	62.1	
45 cis-1,2-Dichloroethene	96	5.934	5.935	-0.001	82	159048	50.0	45.6	
46 2-Butanone (MEK)	43	5.983	5.990	-0.007	100	156296	100.0	86.1	
49 Chlorobromomethane	128	6.226	6.227	-0.001	97	72941	50.0	48.4	
51 Tetrahydrofuran	42	6.287	6.282	0.005	88	89016	100.0	78.3	
52 Chloroform	83	6.342	6.337	0.005	95	270001	50.0	50.3	
53 1,1,1-Trichloroethane	97	6.531	6.525	0.006	96	180759	50.0	52.8	
54 Cyclohexane	56	6.585	6.580	0.005	91	290837	50.0	44.2	
55 1,1-Dichloropropene	75	6.725	6.720	0.005	92	199451	50.0	44.8	
56 Carbon tetrachloride	117	6.719	6.720	-0.001	74	151677	50.0	55.2	
57 Isobutyl alcohol	41	6.944	6.945	-0.001	45	103185	1250.0	1394.0	
58 Benzene	78	6.950	6.951	-0.001	97	671875	50.0	51.1	
59 1,2-Dichloroethane	62	6.987	6.982	0.005	97	215447	50.0	50.1	
62 n-Heptane	43	7.279	7.280	-0.001	89	195888	50.0	43.4	
64 Trichloroethene	130	7.668	7.669	-0.001	97	153772	50.0	46.7	
66 Methylcyclohexane	83	7.863	7.858	0.005	91	257954	50.0	43.9	
67 1,2-Dichloropropane	63	7.899	7.900	-0.001	93	158198	50.0	48.7	
68 Dibromomethane	93	8.021	8.022	-0.001	95	84112	50.0	48.1	
70 1,4-Dioxane	88	8.058	8.052	0.006	97	28593	1000.0	835.4	M
71 Dichlorobromomethane	83	8.198	8.192	0.006	98	163813	50.0	45.9	
73 2-Chloroethyl vinyl ether	63	8.520	8.515	0.005	93	159404	100.0	87.0	
74 cis-1,3-Dichloropropene	75	8.654	8.655	-0.001	92	168788	50.0	49.0	
75 4-Methyl-2-pentanone (MIBK)	43	8.824	8.819	0.005	98	275399	100.0	81.0	
76 Toluene	91	8.988	8.989	-0.001	98	678535	50.0	52.7	
77 trans-1,3-Dichloropropene	75	9.220	9.220	0.000	97	126022	50.0	54.2	
78 Ethyl methacrylate	69	9.317	9.318	-0.001	91	141853	50.0	46.7	
79 1,1,2-Trichloroethane	97	9.396	9.397	-0.001	93	128975	50.0	53.4	
80 Tetrachloroethene	164	9.536	9.531	0.005	97	128550	50.0	51.0	
81 1,3-Dichloropropane	76	9.560	9.561	-0.001	93	228777	50.0	51.0	
82 2-Hexanone	43	9.658	9.652	0.006	97	218428	100.0	84.0	
84 Chlorodibromomethane	129	9.791	9.786	0.005	93	91655	50.0	47.5	
85 Ethylene Dibromide	107	9.901	9.902	-0.001	96	117737	50.0	51.1	
86 3-Chlorobenzotrifluoride	180	10.369	10.370	-0.001	95	234802	50.0	47.8	
87 Chlorobenzene	112	10.388	10.388	0.000	94	420095	50.0	51.5	
88 4-Chlorobenzotrifluoride	180	10.424	10.425	-0.001	94	223092	50.0	46.9	
89 1,1,2-Tetrachloroethane	131	10.473	10.474	-0.001	91	116973	50.0	55.5	
90 Ethylbenzene	106	10.503	10.498	0.005	99	227864	50.0	48.7	
91 m-Xylene & p-Xylene	106	10.619	10.614	0.005	0	288585	50.0	50.4	
92 o-Xylene	106	11.008	11.009	-0.001	98	271589	50.0	48.5	
93 Styrene	104	11.020	11.021	-0.001	95	438285	50.0	48.6	
94 Bromoform	173	11.215	11.204	0.011	96	54705	50.0	45.9	
96 2-Chlorobenzotrifluoride	180	11.270	11.271	-0.001	97	224797	50.0	45.8	
97 Isopropylbenzene	105	11.379	11.380	-0.001	97	675852	50.0	48.4	
99 1,1,2,2-Tetrachloroethane	83	11.671	11.672	-0.001	96	173934	50.0	50.2	
100 Bromobenzene	156	11.683	11.678	0.005	95	161314	50.0	46.7	
101 1,2,3-Trichloropropane	110	11.714	11.721	-0.007	87	54241	50.0	47.8	
102 trans-1,4-Dichloro-2-butene	53	11.732	11.727	0.005	69	40659	50.0	43.1	
103 N-Propylbenzene	120	11.787	11.788	-0.001	99	194718	50.0	45.7	
104 2-Chlorotoluene	126	11.872	11.873	-0.001	96	171139	50.0	47.8	
105 3-Chlorotoluene	126	11.933	11.934	-0.001	96	177881	50.0	44.4	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
106 1,3,5-Trimethylbenzene	105	11.963	11.958	0.005	94	575788	50.0	48.4	
107 4-Chlorotoluene	126	11.981	11.976	0.005	98	184803	50.0	47.7	
108 tert-Butylbenzene	119	12.286	12.287	-0.001	93	438763	50.0	42.6	
110 1,2,4-Trimethylbenzene	105	12.334	12.335	-0.001	97	560239	50.0	45.9	
111 1,2-dichloro-4-(trifluoromethyl)	214	12.401	12.402	-0.001	98	167982	50.0	43.6	
112 sec-Butylbenzene	105	12.505	12.506	-0.001	95	648621	50.0	44.8	
113 1,3-Dichlorobenzene	146	12.614	12.615	-0.001	97	301423	50.0	47.4	
114 4-Isopropyltoluene	119	12.651	12.645	0.006	96	532768	50.0	44.5	
115 1,4-Dichlorobenzene	146	12.705	12.706	-0.001	94	311180	50.0	47.9	
116 2,4-Dichloro-1-(trifluoromethyl)	214	12.754	12.755	-0.001	98	150663	50.0	41.7	
118 2,5-Dichlorobenzotrifluoride	214	12.803	12.804	-0.001	0	180027	50.0	44.6	
120 n-Butylbenzene	91	13.058	13.059	-0.001	98	471441	50.0	43.3	
121 1,2-Dichlorobenzene	146	13.076	13.077	-0.001	95	280617	50.0	47.6	
122 1,2-Dibromo-3-Chloropropan	75	13.855	13.862	-0.007	76	19013	50.0	39.4	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125	14.007	14.008	-0.001	0	515822	150.0	115.8	
124 1,3,5-Trichlorobenzene	180	14.068	14.069	-0.001	96	170685	50.0	47.7	
125 2,3- & 3,4- Dichlorotoluene	125	14.421	14.422	-0.001	0	320296	100.0	73.9	
126 1,2,4-Trichlorobenzene	180	14.689	14.690	-0.001	94	118081	50.0	38.5	
127 Hexachlorobutadiene	225	14.859	14.860	-0.001	96	62498	50.0	42.5	
128 Naphthalene	128	14.938	14.939	-0.001	97	264517	50.0	32.8	
129 1,2,3-Trichlorobenzene	180	15.181	15.182	-0.001	94	97570	50.0	38.8	
131 2,4,5-Trichlorotoluene	159	15.960	15.961	-0.001	0	30053	50.0	22.2	
130 2,3,6-Trichlorotoluene	159	16.057	16.058	-0.001	95	30235	50.0	24.7	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
S 134 1,2-Dichloroethene, Total	96				0		100.0	92.8	
S 133 Xylenes, Total	106				0		100.0	98.9	
S 135 1,3-Dichloropropene, Total	1				0		100.0	103.2	

**QC Flag Legend**

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

**Reagents:**

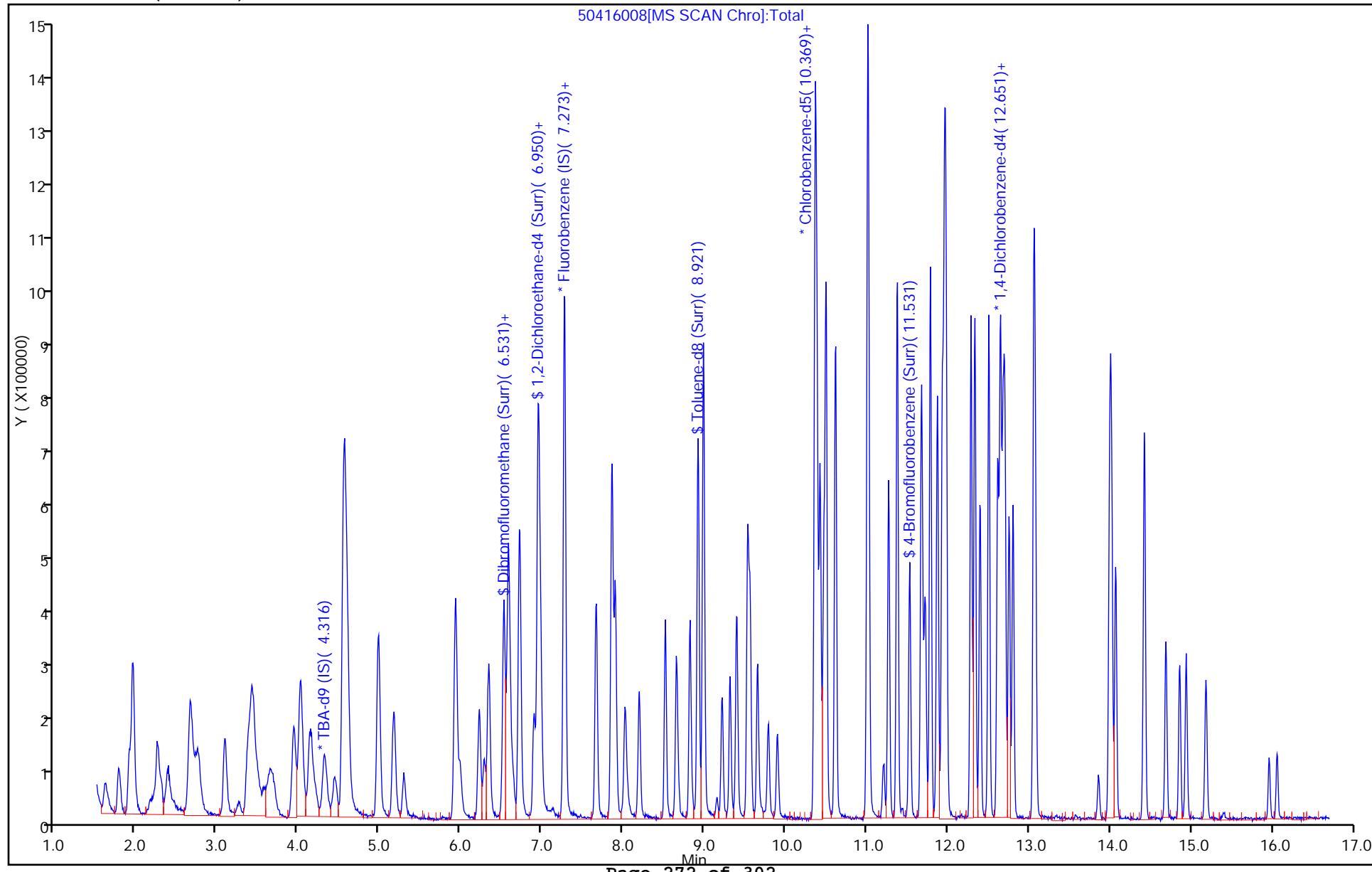
VOA8260VOA2ND_00111	Amount Added: 2.00	Units: uL	
voaWKet2 Rest_00002	Amount Added: 2.00	Units: uL	
voaW ee2nd_00001	Amount Added: 2.00	Units: uL	
voaWVA2nd Res_00006	Amount Added: 2.00	Units: uL	
voaW2-cl 2ndR_00002	Amount Added: 2.00	Units: uL	
voaW 135tcb a_00001	Amount Added: 2.00	Units: uL	
VOAACRO2ND_00007	Amount Added: 6.00	Units: uL	
VOA8260INT_00031	Amount Added: 2.00	Units: uL	Run Reagent
VOA8260SURR_00033	Amount Added: 2.00	Units: uL	Run Reagent

Report Date: 16-Apr-2015 13:57:55

Chrom Revision: 2.2 13-Mar-2015 11:20:44

## TestAmerica Pittsburgh

Data File: \PITCHROM\ChromData\CHHP5\20150416-6494.b\50416008.D  
Injection Date: 16-Apr-2015 13:05:30 Instrument ID: CHHP5  
Lims ID: LCS Operator ID: 001562  
Client ID:  
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 8  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)



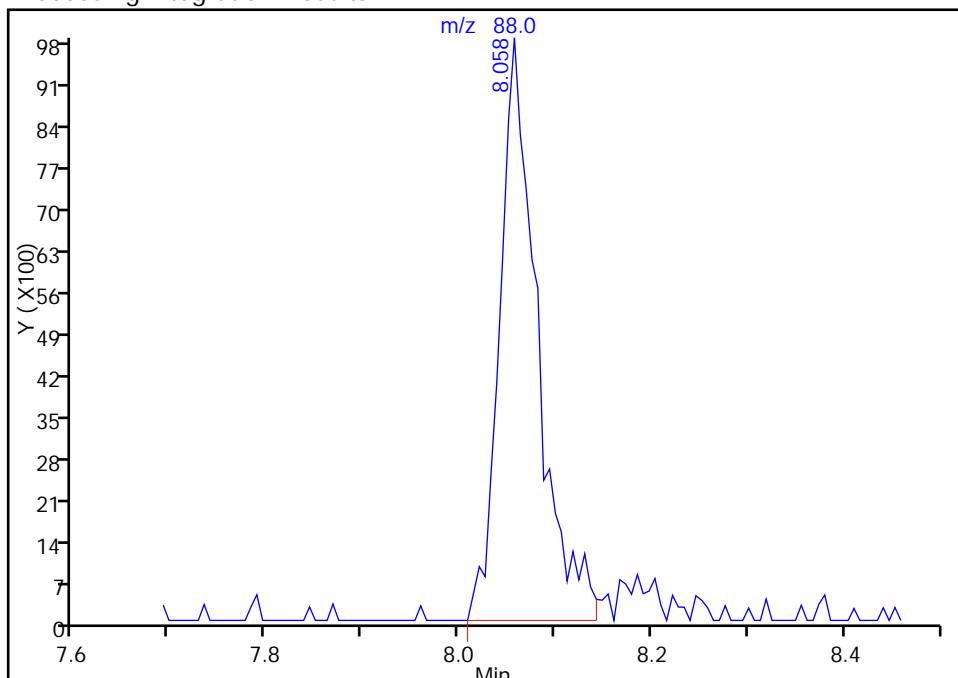
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\50416008.D  
 Injection Date: 16-Apr-2015 13:05:30 Instrument ID: CHHP5  
 Lims ID: LCS  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 8 Worklist Smp#: 8  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 70 1,4-Dioxane, CAS: 123-91-1

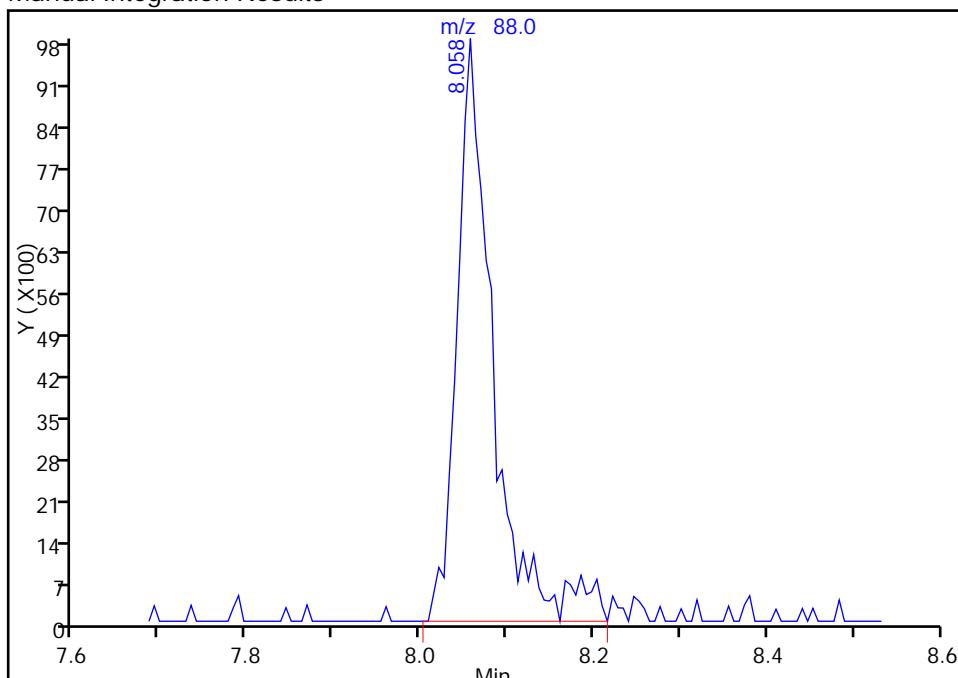
RT: 8.06  
 Area: 26688  
 Amount: 779.7828  
 Amount Units: ng

## Processing Integration Results



RT: 8.06  
 Area: 28593  
 Amount: 835.4440  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 16-Apr-2015 13:37:43

Audit Action: Manually Integrated

Audit Reason: Peak Tail

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: \_\_\_\_\_ Lab Sample ID: LCSD 180-138685/10  
Matrix: Water Lab File ID: 50416010.D  
Analysis Method: 8260C Date Collected: \_\_\_\_\_  
Sample wt/vol: 5 (mL) Date Analyzed: 04/16/2015 13:53  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138685 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
74-87-3	Chloromethane	10.8		1.0	0.28
75-01-4	Vinyl chloride	10.3		1.0	0.23
74-83-9	Bromomethane	8.47		1.0	0.31
75-00-3	Chloroethane	10.7		1.0	0.21
75-35-4	1,1-Dichloroethene	9.04		1.0	0.30
67-64-1	Acetone	20.2		5.0	2.5
75-15-0	Carbon disulfide	6.12		1.0	0.21
75-09-2	Methylene Chloride	10.2		1.0	0.13
156-60-5	trans-1,2-Dichloroethene	9.27		1.0	0.17
1634-04-4	Methyl tert-butyl ether	9.56		1.0	0.18
75-34-3	1,1-Dichloroethane	9.47		1.0	0.12
156-59-2	cis-1,2-Dichloroethene	9.45		1.0	0.24
74-97-5	Bromochloromethane	9.04		1.0	0.18
78-93-3	2-Butanone (MEK)	17.7		5.0	0.55
67-66-3	Chloroform	9.80		1.0	0.17
71-55-6	1,1,1-Trichloroethane	10.2		1.0	0.29
56-23-5	Carbon tetrachloride	10.3		1.0	0.14
71-43-2	Benzene	10.1		1.0	0.11
107-06-2	1,2-Dichloroethane	9.65		1.0	0.21
79-01-6	Trichloroethene	9.24		1.0	0.14
78-87-5	1,2-Dichloroproppane	9.60		1.0	0.095
75-27-4	Bromodichloromethane	8.86		1.0	0.13
10061-01-5	cis-1,3-Dichloropropene	9.88		1.0	0.19
108-10-1	4-Methyl-2-pentanone (MIBK)	16.9		5.0	0.53
108-88-3	Toluene	10.4		1.0	0.15
10061-02-6	trans-1,3-Dichloropropene	10.3		1.0	0.15
79-00-5	1,1,2-Trichloroethane	10.4		1.0	0.20
127-18-4	Tetrachloroethene	10.0		1.0	0.15
591-78-6	2-Hexanone	17.3		5.0	0.16
124-48-1	Dibromochloromethane	9.55		1.0	0.14
106-93-4	1,2-Dibromoethane (EDB)	10.1		1.0	0.18
108-90-7	Chlorobenzene	10.2		1.0	0.14
630-20-6	1,1,1,2-Tetrachloroethane	10.7		1.0	0.28
100-41-4	Ethylbenzene	9.71		1.0	0.23
1330-20-7	Xylenes, Total	19.2		3.0	0.49
100-42-5	Styrene	9.79		1.0	0.097

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: \_\_\_\_\_ Lab Sample ID: LCSD 180-138685/10  
Matrix: Water Lab File ID: 50416010.D  
Analysis Method: 8260C Date Collected: \_\_\_\_\_  
Sample wt/vol: 5 (mL) Date Analyzed: 04/16/2015 13:53  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138685 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-25-2	Bromoform	8.77		1.0	0.19
79-34-5	1,1,2,2-Tetrachloroethane	9.87		1.0	0.20
107-13-1	Acrylonitrile	97.2		20	0.55
123-91-1	1,4-Dioxane	193	J	200	34

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	92		64-135
2037-26-5	Toluene-d8 (Surr)	97		71-118
460-00-4	4-Bromofluorobenzene (Surr)	92		70-118
1868-53-7	Dibromofluoromethane (Surr)	89		70-128

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\50416010.D  
 Lims ID: LCSD  
 Client ID:  
 Sample Type: LCSD  
 Inject. Date: 16-Apr-2015 13:53:30 ALS Bottle#: 10 Worklist Smp#: 10  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: LCSD  
 Misc. Info.: 180-0006494-010  
 Operator ID: 001562 Instrument ID: CHHP5  
 Method: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 15:44:05 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICAL File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 16-Apr-2015 14:11:22

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.321	4.317	0.004	0	173426	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.271	7.268	0.003	98	571009	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.361	10.358	0.003	89	129134	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.679	12.682	-0.003	92	189386	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.529	6.525	0.004	79	115162	50.0	44.3	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.900	6.897	0.003	0	157245	50.0	45.9	
\$ 7 Toluene-d8 (Surr)	98	8.920	8.916	0.004	93	499124	50.0	48.5	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.529	11.532	-0.003	90	171017	50.0	46.1	
11 Dichlorodifluoromethane	85	1.619	1.628	-0.009	98	142318	50.0	58.1	
12 Chloromethane	50	1.790	1.786	0.004	99	183081	50.0	54.2	
13 Vinyl chloride	62	1.924	1.914	0.010	98	195171	50.0	51.7	
14 Butadiene	39	1.966	1.957	0.009	98	206820	50.0	48.0	
15 Bromomethane	94	2.270	2.267	0.003	92	87984	50.0	42.4	
16 Chloroethane	64	2.404	2.413	-0.009	99	139652	50.0	53.5	
17 Dichlorofluoromethane	67	2.660	2.669	-0.009	98	307885	50.0	51.6	
18 Trichlorofluoromethane	101	2.727	2.735	-0.008	95	203865	50.0	45.0	
20 Ethyl ether	59	3.098	3.094	0.004	93	149623	50.0	50.1	
21 Acrolein	56	3.256	3.265	-0.009	98	50605	150.0	139.5	
22 1,1-Dichloroethene	96	3.408	3.386	0.022	95	148765	50.0	45.2	
23 1,1,2-Trichloro-1,2,2-trif	101	3.438	3.441	-0.003	93	171706	50.0	51.6	
24 Acetone	43	3.505	3.496	0.009	100	118276	100.0	101.1	
25 Iodomethane	142	3.627	3.618	0.009	98	207493	50.0	45.4	
26 Carbon disulfide	76	3.676	3.691	-0.015	100	246283	50.0	30.6	
28 3-Chloro-1-propene	76	3.937	3.946	-0.009	89	75693	50.0	43.5	
30 Methyl acetate	43	4.029	4.025	0.004	98	685496	250.0	250.5	
31 Methylene Chloride	84	4.150	4.147	0.003	96	194019	50.0	50.9	
32 2-Methyl-2-propanol	59	4.442	4.439	0.003	86	95983	500.0	469.9	
33 Acrylonitrile	53	4.552	4.548	0.004	99	684240	500.0	486.1	
34 trans-1,2-Dichloroethene	96	4.564	4.561	0.003	59	157792	50.0	46.3	
35 Methyl tert-butyl ether	73	4.594	4.591	0.003	95	359785	50.0	47.8	
36 Hexane	57	4.984	4.986	-0.002	95	233302	50.0	42.9	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
37 1,1-Dichloroethane	63	5.172	5.175	-0.003	96	287698	50.0	47.3	
38 Vinyl acetate	43	5.300	5.291	0.009	97	177892	50.0	41.3	
44 2,2-Dichloropropane	77	5.933	5.929	0.004	55	92464	50.0	60.8	
45 cis-1,2-Dichloroethene	96	5.939	5.935	0.004	82	169553	50.0	47.3	
46 2-Butanone (MEK)	43	5.994	5.990	0.004	100	165698	100.0	88.6	
49 Chlorobromomethane	128	6.225	6.227	-0.002	96	70174	50.0	45.2	
51 Tetrahydrofuran	42	6.292	6.282	0.010	89	101178	100.0	86.4	
52 Chloroform	83	6.340	6.337	0.003	97	270545	50.0	49.0	
53 1,1,1-Trichloroethane	97	6.529	6.525	0.004	95	180178	50.0	51.1	
54 Cyclohexane	56	6.584	6.580	0.004	92	296857	50.0	43.8	
55 1,1-Dichloropropene	75	6.724	6.720	0.004	95	209495	50.0	45.7	
56 Carbon tetrachloride	117	6.717	6.720	-0.003	95	145806	50.0	51.5	
57 Isobutyl alcohol	41	6.943	6.945	-0.002	46	109028	1250.0	1430.5	
58 Benzene	78	6.955	6.951	0.004	97	682565	50.0	50.4	
59 1,2-Dichloroethane	62	6.985	6.982	0.003	97	213741	50.0	48.2	
62 n-Heptane	43	7.283	7.280	0.003	93	198183	50.0	42.6	
64 Trichloroethene	130	7.673	7.669	0.004	97	156546	50.0	46.2	
66 Methylcyclohexane	83	7.861	7.858	0.003	92	269360	50.0	44.5	
67 1,2-Dichloropropane	63	7.904	7.900	0.004	94	160598	50.0	48.0	
68 Dibromomethane	93	8.019	8.022	-0.003	96	83927	50.0	46.6	
70 1,4-Dioxane	88	8.056	8.052	0.004	96	33929	1000.0	962.8	M
71 Dichlorobromomethane	83	8.202	8.192	0.010	98	162930	50.0	44.3	
73 2-Chloroethyl vinyl ether	63	8.518	8.515	0.003	92	171254	100.0	90.8	
74 cis-1,3-Dichloropropene	75	8.658	8.655	0.003	93	175274	50.0	49.4	
75 4-Methyl-2-pentanone (MIBK)	43	8.822	8.819	0.003	98	295118	100.0	84.5	
76 Toluene	91	8.987	8.989	-0.002	98	686403	50.0	51.9	
77 trans-1,3-Dichloropropene	75	9.218	9.220	-0.002	98	122900	50.0	51.4	
78 Ethyl methacrylate	69	9.315	9.318	-0.003	89	155565	50.0	49.9	
79 1,1,2-Trichloroethane	97	9.400	9.397	0.003	92	129395	50.0	52.1	
80 Tetrachloroethene	164	9.540	9.531	0.009	96	129748	50.0	50.1	
81 1,3-Dichloropropane	76	9.558	9.561	-0.003	93	239198	50.0	51.9	
82 2-Hexanone	43	9.656	9.652	0.004	98	230670	100.0	86.4	
84 Chlorodibromomethane	129	9.790	9.786	0.004	91	94565	50.0	47.7	
85 Ethylene Dibromide	107	9.899	9.902	-0.003	98	119861	50.0	50.6	
86 3-Chlorobenzotrifluoride	180	10.368	10.370	-0.002	95	252337	50.0	50.0	
87 Chlorobenzene	112	10.392	10.388	0.004	95	425910	50.0	50.8	
88 4-Chlorobenzotrifluoride	180	10.428	10.425	0.003	95	244743	50.0	50.1	
89 1,1,2-Tetrachloroethane	131	10.471	10.474	-0.003	91	115829	50.0	53.5	
90 Ethylbenzene	106	10.495	10.498	-0.003	99	233596	50.0	48.6	
91 m-Xylene & p-Xylene	106	10.617	10.614	0.003	0	288984	50.0	49.1	
92 o-Xylene	106	11.006	11.009	-0.003	97	271203	50.0	47.1	
93 Styrene	104	11.025	11.021	0.004	95	454073	50.0	49.0	
94 Bromoform	173	11.207	11.204	0.003	95	53615	50.0	43.8	
96 2-Chlorobenzotrifluoride	180	11.274	11.271	0.003	97	241288	50.0	47.9	
97 Isopropylbenzene	105	11.377	11.380	-0.003	97	689338	50.0	48.0	
99 1,1,2,2-Tetrachloroethane	83	11.675	11.672	0.003	95	175634	50.0	49.4	
100 Bromobenzene	156	11.675	11.678	-0.003	95	166892	50.0	47.6	
101 1,2,3-Trichloropropane	110	11.718	11.721	-0.003	88	54354	50.0	47.2	
102 trans-1,4-Dichloro-2-butene	53	11.730	11.727	0.003	66	43276	50.0	45.2	
103 N-Propylbenzene	120	11.785	11.788	-0.003	99	197860	50.0	45.8	
104 2-Chlorotoluene	126	11.870	11.873	-0.003	96	170723	50.0	47.0	
105 3-Chlorotoluene	126	11.931	11.934	-0.003	94	188679	50.0	46.5	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
106 1,3,5-Trimethylbenzene	105	11.961	11.958	0.003	94	574325	50.0	47.6	
107 4-Chlorotoluene	126	11.980	11.976	0.004	98	194023	50.0	49.3	
108 tert-Butylbenzene	119	12.284	12.287	-0.003	93	450766	50.0	43.2	
110 1,2,4-Trimethylbenzene	105	12.332	12.335	-0.003	97	572040	50.0	46.2	
111 1,2-dichloro-4-(trifluoromethyl)	214	12.399	12.402	-0.003	99	178041	50.0	45.5	
112 sec-Butylbenzene	105	12.509	12.506	0.003	95	660628	50.0	44.9	
113 1,3-Dichlorobenzene	146	12.618	12.615	0.003	97	297636	50.0	46.1	
114 4-Isopropyltoluene	119	12.649	12.645	0.004	96	533428	50.0	44.0	
115 1,4-Dichlorobenzene	146	12.704	12.706	-0.002	95	314806	50.0	47.7	
116 2,4-Dichloro-1-(trifluoromethyl)	214	12.758	12.755	0.003	98	157249	50.0	42.9	
118 2,5-Dichlorobenzotrifluoride	214	12.801	12.804	-0.003	0	191342	50.0	46.7	
120 n-Butylbenzene	91	13.056	13.059	-0.003	98	479546	50.0	43.4	
121 1,2-Dichlorobenzene	146	13.075	13.077	-0.002	96	287547	50.0	48.1	
122 1,2-Dibromo-3-Chloropropan	75	13.859	13.862	-0.003	72	19245	50.0	39.3	
123 2,4- & 2,5- & 2,6- Dichlorobenzene	125	14.005	14.008	-0.003	0	550998	150.0	121.9	
124 1,3,5-Trichlorobenzene	180	14.066	14.069	-0.003	96	169934	50.0	46.8	
125 2,3- & 3,4- Dichlorotoluene	125	14.425	14.422	0.003	0	336638	100.0	76.6	
126 1,2,4-Trichlorobenzene	180	14.693	14.690	0.003	94	118499	50.0	38.1	
127 Hexachlorobutadiene	225	14.857	14.860	-0.003	95	61898	50.0	41.5	
128 Naphthalene	128	14.936	14.939	-0.003	97	285958	50.0	35.0	
129 1,2,3-Trichlorobenzene	180	15.180	15.182	-0.002	95	99066	50.0	38.8	
131 2,4,5-Trichlorotoluene	159	15.958	15.961	-0.003	0	34573	50.0	25.2	
130 2,3,6-Trichlorotoluene	159	16.062	16.058	0.004	93	32762	50.0	26.4	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
S 134 1,2-Dichloroethene, Total	96				0		100.0	93.6	
S 133 Xylenes, Total	106				0		100.0	96.2	
S 135 1,3-Dichloropropene, Total	1				0		100.0	100.8	

**QC Flag Legend**

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

**Reagents:**

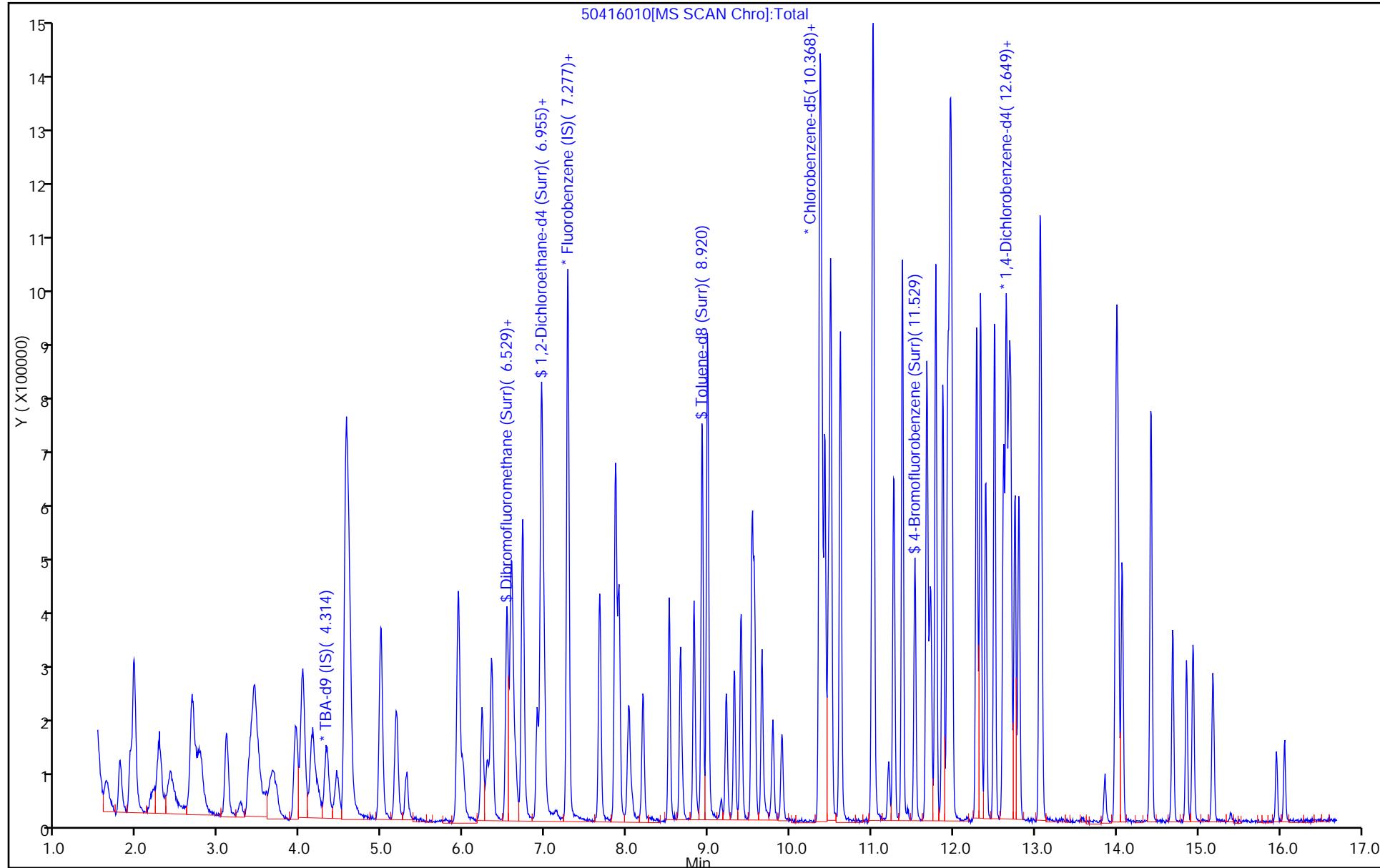
voaW 135tcb a_00001	Amount Added: 2.00	Units: uL
voaW2-cl 2ndR_00002	Amount Added: 2.00	Units: uL
VOA8260VOA2ND_00111	Amount Added: 2.00	Units: uL
voaWKet2 Rest_00002	Amount Added: 2.00	Units: uL
voaW ee2nd_00001	Amount Added: 2.00	Units: uL
voaWVA2nd Res_00006	Amount Added: 2.00	Units: uL
VOAACRO2ND_00007	Amount Added: 6.00	Units: uL
VOA8260INT_00031	Amount Added: 2.00	Units: uL Run Reagent
VOA8260SURR_00033	Amount Added: 2.00	Units: uL Run Reagent

Report Date: 16-Apr-2015 15:44:05

Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Pittsburgh  
Data File: \PITCHROM\ChromData\CHHP5\20150416-6494.b\50416010.D  
Injection Date: 16-Apr-2015 13:53:30 Instrument ID: CHHP5  
Lims ID: LCSD Operator ID: 001562  
Client ID:  
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 10  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)

Worklist Smp#: 10



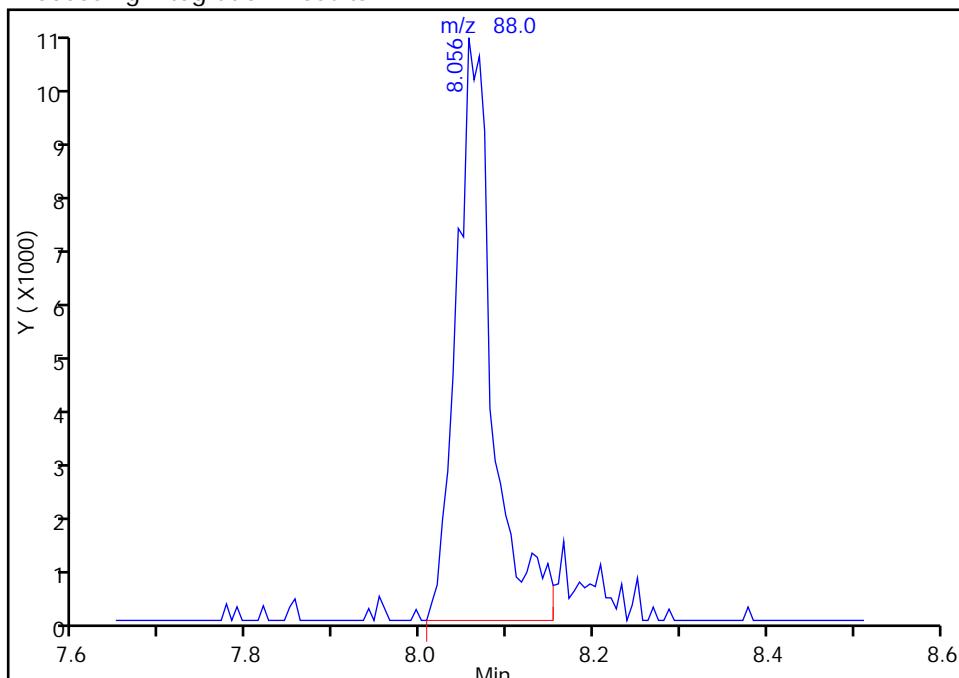
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150416-6494.b\50416010.D  
 Injection Date: 16-Apr-2015 13:53:30 Instrument ID: CHHP5  
 Lims ID: LCSD  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 10 Worklist Smp#: 10  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 70 1,4-Dioxane, CAS: 123-91-1

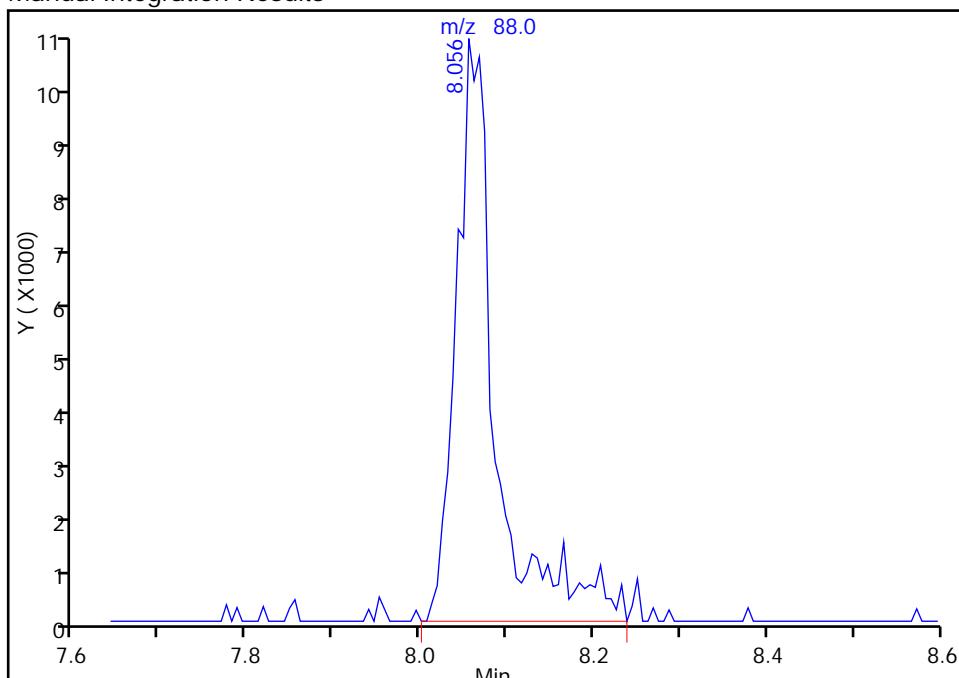
RT: 8.06  
 Area: 30856  
 Amount: 875.5657  
 Amount Units: ng

## Processing Integration Results



RT: 8.06  
 Area: 33929  
 Amount: 962.7647  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 16-Apr-2015 14:11:22

Audit Action: Manually Integrated

Audit Reason: Peak Tail

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Client Sample ID: HD-MW-165-0/1-0 MS Lab Sample ID: 180-42975-1 MS

Matrix: Water Lab File ID: 50415009.D

Analysis Method: 8260C Date Collected: 04/10/2015 08:40

Sample wt/vol: 5 (mL) Date Analyzed: 04/15/2015 16:21

Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1

Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)

% Moisture: \_\_\_\_\_ Level: (low/med) Low

Analysis Batch No.: 138583 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
74-87-3	Chloromethane	10.5		1.0	0.28
75-01-4	Vinyl chloride	9.56		1.0	0.23
74-83-9	Bromomethane	9.20		1.0	0.31
75-00-3	Chloroethane	10.3		1.0	0.21
75-35-4	1,1-Dichloroethene	6.75		1.0	0.30
67-64-1	Acetone	20.2		5.0	2.5
75-15-0	Carbon disulfide	3.81		1.0	0.21
75-09-2	Methylene Chloride	7.73		1.0	0.13
156-60-5	trans-1,2-Dichloroethene	7.54		1.0	0.17
1634-04-4	Methyl tert-butyl ether	8.13		1.0	0.18
75-34-3	1,1-Dichloroethane	8.29		1.0	0.12
156-59-2	cis-1,2-Dichloroethene	8.26		1.0	0.24
74-97-5	Bromochloromethane	7.70		1.0	0.18
78-93-3	2-Butanone (MEK)	16.6		5.0	0.55
67-66-3	Chloroform	9.04		1.0	0.17
71-55-6	1,1,1-Trichloroethane	8.69		1.0	0.29
56-23-5	Carbon tetrachloride	8.96		1.0	0.14
71-43-2	Benzene	8.62		1.0	0.11
107-06-2	1,2-Dichloroethane	8.54		1.0	0.21
79-01-6	Trichloroethene	19.5		1.0	0.14
78-87-5	1,2-Dichloroproppane	8.78		1.0	0.095
75-27-4	Bromodichloromethane	8.29		1.0	0.13
10061-01-5	cis-1,3-Dichloropropene	8.80		1.0	0.19
108-10-1	4-Methyl-2-pentanone (MIBK)	16.8		5.0	0.53
108-88-3	Toluene	9.59		1.0	0.15
10061-02-6	trans-1,3-Dichloropropene	11.0		1.0	0.15
79-00-5	1,1,2-Trichloroethane	10.7		1.0	0.20
127-18-4	Tetrachloroethene	15.2		1.0	0.15
591-78-6	2-Hexanone	15.4		5.0	0.16
124-48-1	Dibromochloromethane	9.09		1.0	0.14
106-93-4	1,2-Dibromoethane (EDB)	9.76		1.0	0.18
108-90-7	Chlorobenzene	9.83		1.0	0.14
630-20-6	1,1,1,2-Tetrachloroethane	10.7		1.0	0.28
100-41-4	Ethylbenzene	9.18		1.0	0.23
1330-20-7	Xylenes, Total	18.5		3.0	0.49
100-42-5	Styrene	9.29		1.0	0.097

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: HD-MW-165-0/1-0 MS Lab Sample ID: 180-42975-1 MS  
Matrix: Water Lab File ID: 50415009.D  
Analysis Method: 8260C Date Collected: 04/10/2015 08:40  
Sample wt/vol: 5 (mL) Date Analyzed: 04/15/2015 16:21  
Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1  
Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)  
% Moisture: \_\_\_\_\_ Level: (low/med) Low  
Analysis Batch No.: 138583 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-25-2	Bromoform	8.75		1.0	0.19
79-34-5	1,1,2,2-Tetrachloroethane	10.1		1.0	0.20
107-13-1	Acrylonitrile	87.1		20	0.55
123-91-1	1,4-Dioxane	175	J	200	34

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	90		64-135
2037-26-5	Toluene-d8 (Surr)	99		71-118
460-00-4	4-Bromofluorobenzene (Surr)	89		70-118
1868-53-7	Dibromofluoromethane (Surr)	92		70-128

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415009.D  
 Lims ID: 180-42975-B-1 MS  
 Client ID:  
 Sample Type: MS  
 Inject. Date: 15-Apr-2015 16:21:30 ALS Bottle#: 9 Worklist Smp#: 9  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: 180-42975-B-1 MS  
 Misc. Info.: 180-0006480-009  
 Operator ID: 001562 Instrument ID: CHHP5  
 Method: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 07:36:15 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICAL File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 16-Apr-2015 07:36:15

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.317	4.302	0.015	0	171044	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.274	7.271	0.003	99	564634	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.358	10.361	-0.003	88	124370	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.682	12.679	0.003	93	179202	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.531	6.528	0.003	86	118675	50.0	46.2	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.903	6.899	0.004	0	152297	50.0	45.0	
\$ 7 Toluene-d8 (Surr)	98	8.922	8.919	0.003	94	491669	50.0	49.6	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.532	11.529	0.003	88	159670	50.0	44.7	
11 Dichlorodifluoromethane	85	1.622	1.619	0.003	99	144468	50.0	59.7	
12 Chloromethane	50	1.780	1.789	-0.009	99	175230	50.0	52.5	
13 Vinyl chloride	62	1.914	1.917	-0.003	98	178485	50.0	47.8	
14 Butadiene	39	1.957	1.959	-0.002	96	176635	50.0	41.4	
15 Bromomethane	94	2.261	2.264	-0.003	92	93739	50.0	46.0	
16 Chloroethane	64	2.389	2.416	-0.027	99	133274	50.0	51.6	
17 Dichlorofluoromethane	67	2.656	2.671	-0.015	98	297243	50.0	50.4	
18 Trichlorofluoromethane	101	2.729	2.726	0.003	96	206967	50.0	46.2	
20 Ethyl ether	59	3.100	3.091	0.009	93	119047	50.0	40.3	
21 Acrolein	56	3.259	3.261	-0.002	99	48740	150.0	135.9	
22 1,1-Dichloroethene	96	3.392	3.395	-0.003	98	109858	50.0	33.7	
23 1,1,2-Trichloro-1,2,2-trif	101	3.429	3.450	-0.021	96	121320	50.0	36.8	
24 Acetone	43	3.502	3.499	0.003	98	116551	100.0	100.8	
25 Iodomethane	142	3.617	3.626	-0.009	97	159233	50.0	35.2	
26 Carbon disulfide	76	3.678	3.675	0.003	99	151798	50.0	19.1	
28 3-Chloro-1-propene	76	3.934	3.943	-0.009	88	52535	50.0	30.5	
30 Methyl acetate	43	4.025	4.022	0.003	98	576784	250.0	213.1	
31 Methylene Chloride	84	4.141	4.143	-0.002	95	145585	50.0	38.7	
32 2-Methyl-2-propanol	59	4.439	4.441	-0.002	86	98826	500.0	490.5	
33 Acrylonitrile	53	4.554	4.551	0.003	98	606250	500.0	435.5	
34 trans-1,2-Dichloroethene	96	4.560	4.563	-0.003	50	126905	50.0	37.7	
35 Methyl tert-butyl ether	73	4.597	4.600	-0.003	95	302608	50.0	40.6	
36 Hexane	57	4.986	4.983	0.003	94	167066	50.0	31.1	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
37 1,1-Dichloroethane	63	5.175	5.165	0.010	97	249275	50.0	41.5	
38 Vinyl acetate	43	5.297	5.299	-0.002	97	206999	50.0	48.5	
44 2,2-Dichloropropane	77	5.935	5.926	0.009	52	73672	50.0	49.0	
45 cis-1,2-Dichloroethene	96	5.935	5.938	-0.003	82	146542	50.0	41.3	
46 2-Butanone (MEK)	43	5.996	5.987	0.009	100	153467	100.0	83.0	
49 Chlorobromomethane	128	6.227	6.224	0.003	96	59145	50.0	38.5	
51 Tetrahydrofuran	42	6.288	6.279	0.009	88	86779	100.0	74.9	
52 Chloroform	83	6.343	6.339	0.004	94	246732	50.0	45.2	
53 1,1,1-Trichloroethane	97	6.531	6.528	0.003	97	151504	50.0	43.4	
54 Cyclohexane	56	6.586	6.583	0.003	94	215770	50.0	32.2	
56 Carbon tetrachloride	117	6.714	6.723	-0.009	73	125303	50.0	44.8	
55 1,1-Dichloropropene	75	6.720	6.723	-0.003	93	174743	50.0	38.6	
57 Isobutyl alcohol	41	6.945	6.942	0.003	47	99161	1250.0	1315.7	
58 Benzene	78	6.957	6.954	0.003	97	576872	50.0	43.1	
59 1,2-Dichloroethane	62	6.988	6.984	0.004	97	187086	50.0	42.7	
62 n-Heptane	43	7.280	7.276	0.004	87	158992	50.0	34.6	
64 Trichloroethene	130	7.669	7.666	0.003	97	327201	50.0	97.6	
66 Methylcyclohexane	83	7.864	7.860	0.004	92	201471	50.0	33.7	
67 1,2-Dichloropropane	63	7.900	7.897	0.003	95	145307	50.0	43.9	
68 Dibromomethane	93	8.022	8.025	-0.003	95	80514	50.0	45.2	
70 1,4-Dioxane	88	8.058	8.049	0.009	95	30486	1000.0	874.8	M
71 Dichlorobromomethane	83	8.198	8.195	0.003	98	150639	50.0	41.4	
74 cis-1,3-Dichloropropene	75	8.655	8.651	0.004	93	154327	50.0	44.0	
75 4-Methyl-2-pentanone (MIBK)	43	8.825	8.822	0.003	98	282226	100.0	83.9	
76 Toluene	91	8.989	8.986	0.003	97	611329	50.0	48.0	
77 trans-1,3-Dichloropropene	75	9.214	9.217	-0.003	97	127015	50.0	55.2	
78 Ethyl methacrylate	69	9.312	9.314	-0.002	91	139359	50.0	46.4	
79 1,1,2-Trichloroethane	97	9.397	9.399	-0.002	92	127652	50.0	53.4	
80 Tetrachloroethene	164	9.537	9.533	0.004	96	189361	50.0	76.0	
81 1,3-Dichloropropane	76	9.561	9.558	0.003	94	219497	50.0	49.4	
82 2-Hexanone	43	9.658	9.655	0.003	98	198453	100.0	77.2	
84 Chlorodibromomethane	129	9.792	9.789	0.003	90	86740	50.0	45.5	
85 Ethylene Dibromide	107	9.896	9.898	-0.002	100	111275	50.0	48.8	
86 3-Chlorobenzotrifluoride	180	10.370	10.373	-0.003	96	234834	50.0	48.3	
87 Chlorobenzene	112	10.388	10.391	-0.003	94	397047	50.0	49.2	
88 4-Chlorobenzotrifluoride	180	10.425	10.428	-0.003	96	217606	50.0	46.3	
89 1,1,1,2-Tetrachloroethane	131	10.467	10.470	-0.003	91	111834	50.0	53.6	
90 Ethylbenzene	106	10.498	10.501	-0.003	99	212688	50.0	45.9	
91 m-Xylene & p-Xylene	106	10.613	10.616	-0.003	0	265021	50.0	46.8	
92 o-Xylene	106	11.009	11.012	-0.003	96	253539	50.0	45.7	
93 Styrene	104	11.021	11.024	-0.003	91	415059	50.0	46.5	
94 Bromoform	173	11.210	11.212	-0.002	95	51537	50.0	43.7	
96 2-Chlorobenzotrifluoride	180	11.270	11.273	-0.003	97	222158	50.0	45.7	
97 Isopropylbenzene	105	11.374	11.377	-0.003	97	637754	50.0	46.1	
99 1,1,2,2-Tetrachloroethane	83	11.672	11.675	-0.003	95	173381	50.0	50.6	
100 Bromobenzene	156	11.684	11.681	0.003	96	154530	50.0	46.6	
101 1,2,3-Trichloropropane	110	11.721	11.717	0.004	87	55090	50.0	50.5	
102 trans-1,4-Dichloro-2-butene	53	11.733	11.729	0.004	66	43700	50.0	48.2	
103 N-Propylbenzene	120	11.788	11.784	0.004	99	181947	50.0	44.5	
104 2-Chlorotoluene	126	11.873	11.869	0.004	95	156147	50.0	45.4	
105 3-Chlorotoluene	126	11.934	11.930	0.004	95	175084	50.0	45.6	
106 1,3,5-Trimethylbenzene	105	11.958	11.961	-0.003	94	527957	50.0	46.3	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
107 4-Chlorotoluene	126	11.976	11.985	-0.009	98	173598	50.0	46.7	
108 tert-Butylbenzene	119	12.286	12.283	0.003	92	420758	50.0	42.6	
110 1,2,4-Trimethylbenzene	105	12.335	12.332	0.003	95	520451	50.0	44.5	
111 1,2-dichloro-4-(trifluorom	214	12.396	12.399	-0.003	98	162348	50.0	43.9	
112 sec-Butylbenzene	105	12.505	12.502	0.003	95	620992	50.0	44.6	
113 1,3-Dichlorobenzene	146	12.621	12.618	0.003	97	281351	50.0	46.0	
114 4-Isopropyltoluene	119	12.651	12.648	0.003	97	512873	50.0	44.7	
115 1,4-Dichlorobenzene	146	12.706	12.703	0.003	93	291448	50.0	46.7	
116 2,4-Dichloro-1-(trifluorom	214	12.755	12.758	-0.003	98	151697	50.0	43.8	
118 2,5-Dichlorobenzotrifluori	214	12.804	12.806	-0.002	0	171573	50.0	44.3	
120 n-Butylbenzene	91	13.059	13.062	-0.003	98	439422	50.0	42.0	
121 1,2-Dichlorobenzene	146	13.077	13.080	-0.003	96	267602	50.0	47.3	
122 1,2-Dibromo-3-Chloropropan	75	13.856	13.853	0.003	75	19436	50.0	42.0	
123 2,4- & 2,5- & 2,6- Dichlor	125	14.002	14.005	-0.003	0	503999	150.0	117.8	
125 2,3- & 3,4- Dichlorotoluen	125	14.422	14.424	-0.002	0	305938	100.0	73.5	
126 1,2,4-Trichlorobenzene	180	14.689	14.686	0.003	92	117212	50.0	39.8	
127 Hexachlorobutadiene	225	14.860	14.856	0.004	97	62143	50.0	44.0	
128 Naphthalene	128	14.939	14.942	-0.003	97	273191	50.0	35.3	
129 1,2,3-Trichlorobenzene	180	15.182	15.185	-0.003	95	92525	50.0	38.3	
131 2,4,5-Trichlorotoluene	159	15.961	15.964	-0.003	0	33586	50.0	25.9	
130 2,3,6-Trichlorotoluene	159	16.058	16.061	-0.003	96	34777	50.0	29.6	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
S 133 Xylenes, Total	106				0		100.0	92.5	
S 134 1,2-Dichloroethene, Total	96				0		100.0	79.0	
S 135 1,3-Dichloropropene, Total	1				0		100.0	99.2	

### QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

### Reagents:

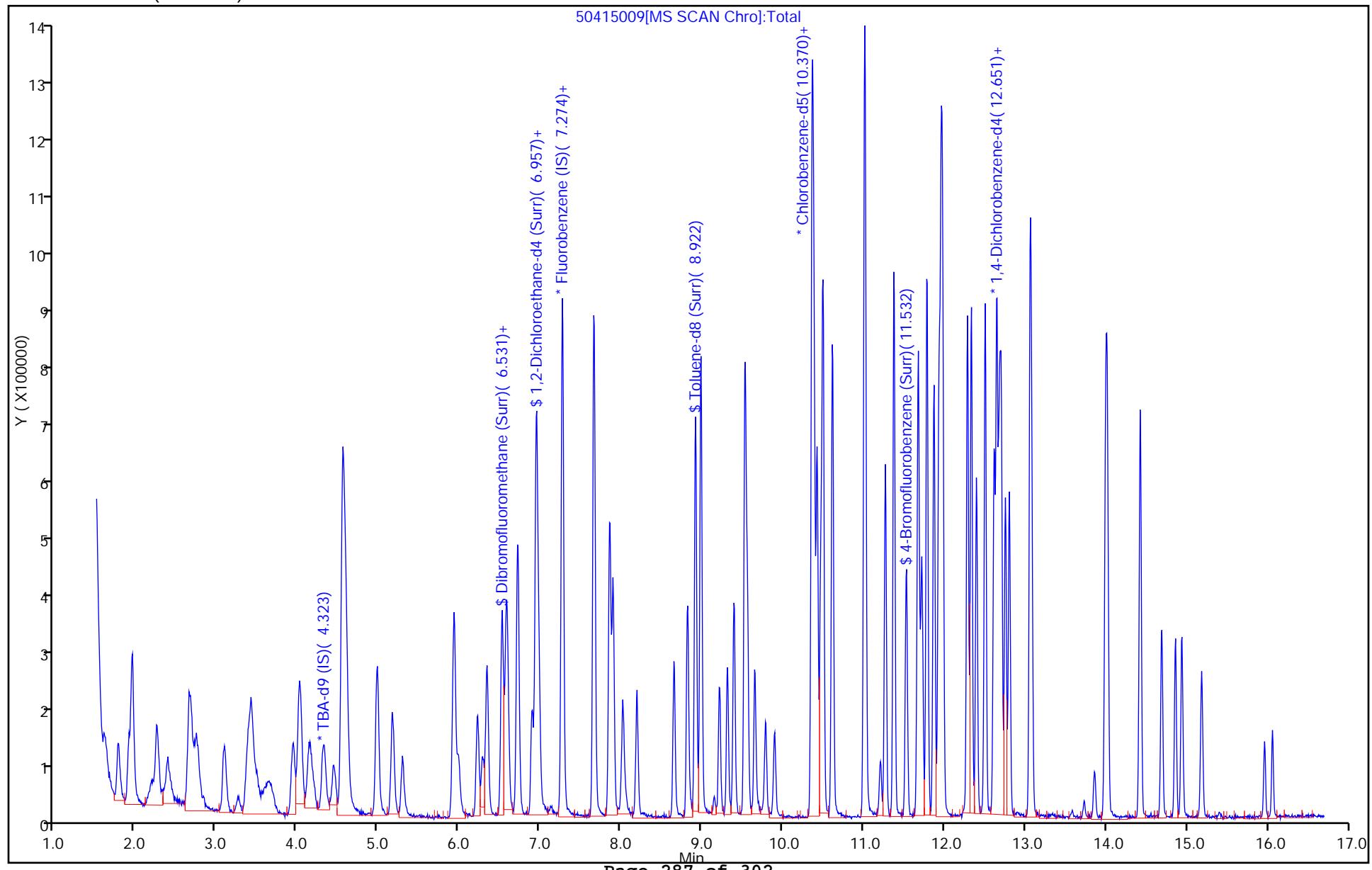
VOA8260VOA2ND_00111	Amount Added: 2.00	Units: uL	
voaWKet2 Rest_00002	Amount Added: 2.00	Units: uL	
voaW ee2nd_00001	Amount Added: 2.00	Units: uL	
voaWVA2nd Res_00006	Amount Added: 2.00	Units: uL	
VOAACRO2ND_00007	Amount Added: 6.00	Units: uL	
VOA8260INT_00031	Amount Added: 2.00	Units: uL	Run Reagent
VOA8260SURR_00033	Amount Added: 2.00	Units: uL	Run Reagent

Report Date: 16-Apr-2015 07:36:16

Chrom Revision: 2.2 13-Mar-2015 11:20:44

## TestAmerica Pittsburgh

Data File: \\PITCHROM\\ChromData\\CHHP5\\20150415-6480.b\\50415009.D  
Injection Date: 15-Apr-2015 16:21:30 Instrument ID: CHHP5  
Lims ID: 180-42975-B-1 MS  
Client ID:  
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 9  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)

Operator ID: 001562  
Worklist Smp#: 9

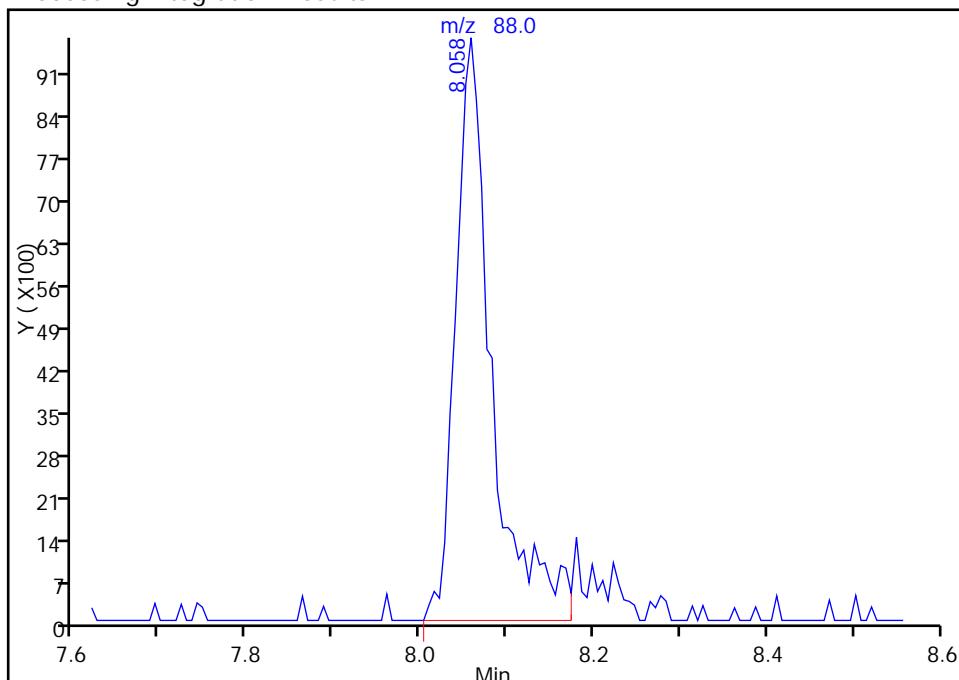
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415009.D  
 Injection Date: 15-Apr-2015 16:21:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-B-1 MS  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 9 Worklist Smp#: 9  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 70 1,4-Dioxane, CAS: 123-91-1

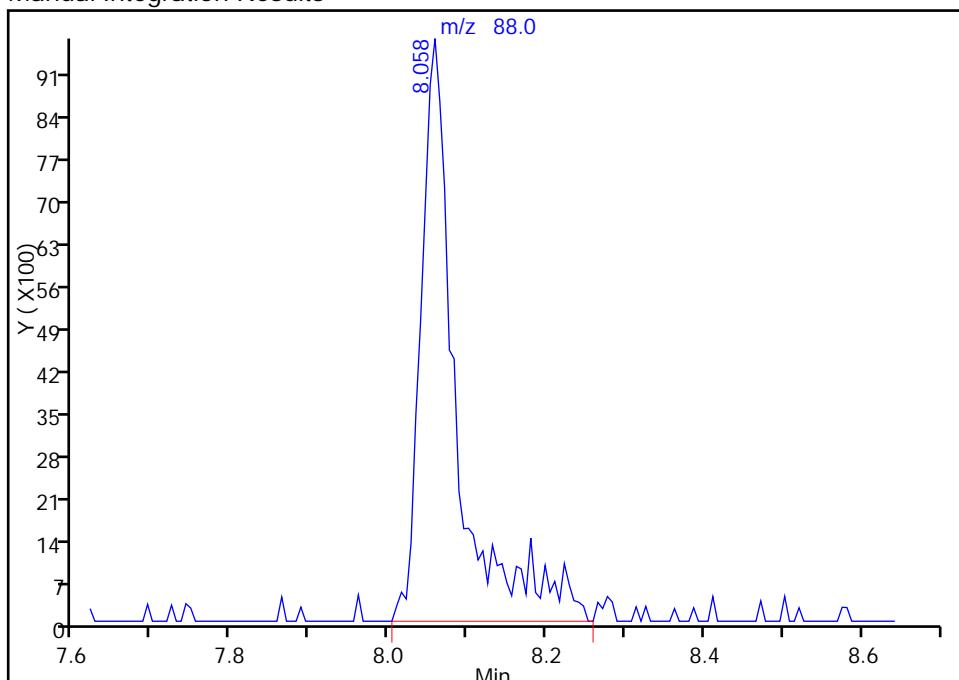
RT: 8.06  
 Area: 27909  
 Amount: 800.8834  
 Amount Units: ng

## Processing Integration Results



RT: 8.06  
 Area: 30486  
 Amount: 874.8337  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 16-Apr-2015 07:36:15

Audit Action: Manually Integrated

Audit Reason: Peak Tail

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Client Sample ID: HD-MW-165-0/1-0 MSD Lab Sample ID: 180-42975-1 MSD

Matrix: Water Lab File ID: 50415010.D

Analysis Method: 8260C Date Collected: 04/10/2015 08:40

Sample wt/vol: 5 (mL) Date Analyzed: 04/15/2015 16:46

Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1

Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)

% Moisture: \_\_\_\_\_ Level: (low/med) Low

Analysis Batch No.: 138583 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
74-87-3	Chloromethane	10.8		1.0	0.28
75-01-4	Vinyl chloride	9.37		1.0	0.23
74-83-9	Bromomethane	10.1		1.0	0.31
75-00-3	Chloroethane	10.4		1.0	0.21
75-35-4	1,1-Dichloroethene	6.67		1.0	0.30
67-64-1	Acetone	19.1		5.0	2.5
75-15-0	Carbon disulfide	3.88		1.0	0.21
75-09-2	Methylene Chloride	8.19		1.0	0.13
156-60-5	trans-1,2-Dichloroethene	7.67		1.0	0.17
1634-04-4	Methyl tert-butyl ether	8.73		1.0	0.18
75-34-3	1,1-Dichloroethane	8.39		1.0	0.12
156-59-2	cis-1,2-Dichloroethene	8.61		1.0	0.24
74-97-5	Bromochloromethane	9.01		1.0	0.18
78-93-3	2-Butanone (MEK)	18.7		5.0	0.55
67-66-3	Chloroform	9.15		1.0	0.17
71-55-6	1,1,1-Trichloroethane	8.67		1.0	0.29
56-23-5	Carbon tetrachloride	8.65		1.0	0.14
71-43-2	Benzene	8.79		1.0	0.11
107-06-2	1,2-Dichloroethane	9.38		1.0	0.21
79-01-6	Trichloroethene	19.7		1.0	0.14
78-87-5	1,2-Dichloroproppane	9.80		1.0	0.095
75-27-4	Bromodichloromethane	9.01		1.0	0.13
10061-01-5	cis-1,3-Dichloropropene	9.60		1.0	0.19
108-10-1	4-Methyl-2-pentanone (MIBK)	18.0		5.0	0.53
108-88-3	Toluene	9.38		1.0	0.15
10061-02-6	trans-1,3-Dichloropropene	11.3		1.0	0.15
79-00-5	1,1,2-Trichloroethane	10.4		1.0	0.20
127-18-4	Tetrachloroethene	14.2		1.0	0.15
591-78-6	2-Hexanone	16.2		5.0	0.16
124-48-1	Dibromochloromethane	9.15		1.0	0.14
106-93-4	1,2-Dibromoethane (EDB)	10.2		1.0	0.18
108-90-7	Chlorobenzene	9.90		1.0	0.14
630-20-6	1,1,1,2-Tetrachloroethane	10.8		1.0	0.28
100-41-4	Ethylbenzene	9.33		1.0	0.23
1330-20-7	Xylenes, Total	18.3		3.0	0.49
100-42-5	Styrene	9.54		1.0	0.097

FORM I  
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Client Sample ID: HD-MW-165-0/1-0 MSD Lab Sample ID: 180-42975-1 MSD

Matrix: Water Lab File ID: 50415010.D

Analysis Method: 8260C Date Collected: 04/10/2015 08:40

Sample wt/vol: 5 (mL) Date Analyzed: 04/15/2015 16:46

Soil Aliquot Vol: \_\_\_\_\_ Dilution Factor: 1

Soil Extract Vol.: \_\_\_\_\_ GC Column: DB-624 ID: 0.18 (mm)

% Moisture: \_\_\_\_\_ Level: (low/med) Low

Analysis Batch No.: 138583 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-25-2	Bromoform	8.52		1.0	0.19
79-34-5	1,1,2,2-Tetrachloroethane	10.5		1.0	0.20
107-13-1	Acrylonitrile	94.6		20	0.55
123-91-1	1,4-Dioxane	191	J	200	34

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	96		64-135
2037-26-5	Toluene-d8 (Surr)	96		71-118
460-00-4	4-Bromofluorobenzene (Surr)	93		70-118
1868-53-7	Dibromofluoromethane (Surr)	92		70-128

TestAmerica Pittsburgh  
Target Compound Quantitation Report

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415010.D  
 Lims ID: 180-42975-C-1 MSD  
 Client ID:  
 Sample Type: MSD  
 Inject. Date: 15-Apr-2015 16:46:30 ALS Bottle#: 10 Worklist Smp#: 10  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Sample Info: 180-42975-C-1 MSD  
 Misc. Info.: 180-0006480-010  
 Operator ID: 001562 Instrument ID: CHHP5  
 Method: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\MSVOA\_LL\_CHHP5.m  
 Limit Group: VOA 8260C ICAL  
 Last Update: 16-Apr-2015 07:37:38 Calib Date: 14-Apr-2015 13:44:30  
 Integrator: RTE ID Type: Deconvolution ID  
 Quant Method: Internal Standard Quant By: Initial Calibration  
 Last ICAL File: \\PITCHROM\ChromData\CHHP5\20150414-6459.b\50414005.D  
 Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN  
 Process Host: XAWRK025

First Level Reviewer: fergusond Date: 16-Apr-2015 07:37:38

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
* 1 TBA-d9 (IS)	65	4.311	4.302	0.009	0	208690	1000.0	1000.0	
* 2 Fluorobenzene (IS)	96	7.274	7.271	0.003	99	561630	50.0	50.0	
* 3 Chlorobenzene-d5	119	10.358	10.361	-0.003	87	131185	50.0	50.0	
* 4 1,4-Dichlorobenzene-d4	152	12.676	12.679	-0.003	95	190234	50.0	50.0	
\$ 5 Dibromofluoromethane (Surr)	113	6.532	6.528	0.004	87	117065	50.0	45.8	
\$ 6 1,2-Dichloroethane-d4 (Sur)	65	6.903	6.899	0.004	0	161730	50.0	48.0	
\$ 7 Toluene-d8 (Surr)	98	8.923	8.919	0.004	93	500912	50.0	47.9	
\$ 8 4-Bromofluorobenzene (Surr)	95	11.526	11.529	-0.003	88	175867	50.0	46.7	
11 Dichlorodifluoromethane	85	1.616	1.619	-0.003	99	138723	50.0	57.6	
12 Chloromethane	50	1.787	1.789	-0.002	99	179928	50.0	54.2	
13 Vinyl chloride	62	1.915	1.917	-0.002	99	173934	50.0	46.8	
14 Butadiene	39	1.951	1.959	-0.008	97	163700	50.0	38.6	
15 Bromomethane	94	2.255	2.264	-0.009	91	101574	50.0	50.6	
16 Chloroethane	64	2.389	2.416	-0.027	98	133013	50.0	51.8	
17 Dichlorofluoromethane	67	2.663	2.671	-0.008	98	294651	50.0	50.2	
18 Trichlorofluoromethane	101	2.736	2.726	0.010	94	196877	50.0	44.2	
20 Ethyl ether	59	3.095	3.091	0.004	94	120922	50.0	41.2	
21 Acrolein	56	3.265	3.261	0.004	99	48965	150.0	137.2	
22 1,1-Dichloroethene	96	3.381	3.395	-0.014	97	107958	50.0	33.3	
23 1,1,2-Trichloro-1,2,2-trif	101	3.435	3.450	-0.015	95	117202	50.0	35.8	
24 Acetone	43	3.496	3.499	-0.003	98	110022	100.0	95.6	
25 Iodomethane	142	3.588	3.626	-0.038	98	162152	50.0	36.0	
26 Carbon disulfide	76	3.685	3.675	0.010	99	153606	50.0	19.4	
28 3-Chloro-1-propene	76	3.946	3.943	0.003	91	54865	50.0	32.1	
30 Methyl acetate	43	4.032	4.022	0.010	98	645732	250.0	239.9	
31 Methylene Chloride	84	4.141	4.143	-0.002	92	153291	50.0	40.9	
32 2-Methyl-2-propanol	59	4.439	4.441	-0.002	84	108108	500.0	439.8	
33 Acrylonitrile	53	4.555	4.551	0.004	99	654835	500.0	473.0	
34 trans-1,2-Dichloroethene	96	4.561	4.563	-0.002	49	128548	50.0	38.4	
35 Methyl tert-butyl ether	73	4.597	4.600	-0.003	95	323417	50.0	43.7	
36 Hexane	57	4.981	4.983	-0.002	95	164136	50.0	30.7	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
37 1,1-Dichloroethane	63	5.175	5.165	0.010	97	250874	50.0	42.0	
38 Vinyl acetate	43	5.297	5.299	-0.002	98	220123	50.0	51.9	
44 2,2-Dichloropropane	77	5.930	5.926	0.004	52	75362	50.0	50.4	
45 cis-1,2-Dichloroethene	96	5.942	5.938	0.004	82	152018	50.0	43.1	
46 2-Butanone (MEK)	43	5.991	5.987	0.003	100	172190	100.0	93.6	
49 Chlorobromomethane	128	6.228	6.224	0.004	94	68776	50.0	45.0	
51 Tetrahydrofuran	42	6.289	6.279	0.010	93	90790	100.0	78.8	
52 Chloroform	83	6.343	6.339	0.004	95	248630	50.0	45.8	
53 1,1,1-Trichloroethane	97	6.532	6.528	0.004	97	150357	50.0	43.4	
54 Cyclohexane	56	6.587	6.583	0.004	93	214122	50.0	32.2	
56 Carbon tetrachloride	117	6.721	6.723	-0.002	93	120323	50.0	43.2	
55 1,1-Dichloropropene	75	6.721	6.723	-0.002	94	174363	50.0	38.7	
57 Isobutyl alcohol	41	6.940	6.942	-0.002	52	120399	1250.0	1606.1	
58 Benzene	78	6.952	6.954	-0.002	97	584745	50.0	43.9	
59 1,2-Dichloroethane	62	6.982	6.984	-0.002	97	204477	50.0	46.9	
62 n-Heptane	43	7.274	7.276	-0.002	92	151871	50.0	33.2	
64 Trichloroethene	130	7.663	7.666	-0.003	98	328224	50.0	98.4	
66 Methylcyclohexane	83	7.858	7.860	-0.002	90	202898	50.0	34.1	
67 1,2-Dichloropropane	63	7.901	7.897	0.004	95	161282	50.0	49.0	
68 Dibromomethane	93	8.022	8.025	-0.003	96	81893	50.0	46.2	
70 1,4-Dioxane	88	8.053	8.049	0.004	97	33017	1000.0	952.5	M
71 Dichlorobromomethane	83	8.199	8.195	0.004	99	162905	50.0	45.0	
74 cis-1,3-Dichloropropene	75	8.655	8.651	0.004	93	167450	50.0	48.0	
75 4-Methyl-2-pentanone (MIBK)	43	8.819	8.822	-0.003	98	319991	100.0	90.1	
76 Toluene	91	8.990	8.986	0.004	98	630641	50.0	46.9	
77 trans-1,3-Dichloropropene	75	9.221	9.217	0.004	97	137652	50.0	56.7	
78 Ethyl methacrylate	69	9.318	9.314	0.004	90	157466	50.0	49.7	
79 1,1,2-Trichloroethane	97	9.397	9.399	-0.002	93	130513	50.0	51.8	
80 Tetrachloroethene	164	9.531	9.533	-0.002	96	186782	50.0	71.0	
81 1,3-Dichloropropane	76	9.562	9.558	0.004	94	243904	50.0	52.1	
82 2-Hexanone	43	9.659	9.655	0.004	99	220123	100.0	81.1	
84 Chlorodibromomethane	129	9.787	9.789	-0.002	91	92056	50.0	45.7	
85 Ethylene Dibromide	107	9.902	9.898	0.004	99	122121	50.0	50.8	
86 3-Chlorobenzotrifluoride	180	10.371	10.373	-0.002	94	226719	50.0	44.2	
87 Chlorobenzene	112	10.389	10.391	-0.002	94	421555	50.0	49.5	
88 4-Chlorobenzotrifluoride	180	10.425	10.428	-0.003	95	208661	50.0	42.1	
89 1,1,1,2-Tetrachloroethane	131	10.468	10.470	-0.002	91	118449	50.0	53.9	
90 Ethylbenzene	106	10.498	10.501	-0.003	99	227932	50.0	46.6	
91 m-Xylene & p-Xylene	106	10.614	10.616	-0.002	0	275364	50.0	46.1	
92 o-Xylene	106	11.009	11.012	-0.003	95	264354	50.0	45.2	
93 Styrene	104	11.022	11.024	-0.002	91	449546	50.0	47.7	
94 Bromoform	173	11.210	11.212	-0.002	95	52937	50.0	42.6	
96 2-Chlorobenzotrifluoride	180	11.271	11.273	-0.002	97	221120	50.0	43.2	
97 Isopropylbenzene	105	11.374	11.377	-0.003	97	653479	50.0	44.8	
99 1,1,2,2-Tetrachloroethane	83	11.672	11.675	-0.003	96	190352	50.0	52.7	
100 Bromobenzene	156	11.679	11.681	-0.002	96	167748	50.0	47.6	
101 1,2,3-Trichloropropane	110	11.715	11.717	-0.002	88	61037	50.0	52.8	
102 trans-1,4-Dichloro-2-butene	53	11.727	11.729	-0.002	67	47629	50.0	49.5	
103 N-Propylbenzene	120	11.782	11.784	-0.002	99	191488	50.0	44.1	
104 2-Chlorotoluene	126	11.873	11.869	0.004	96	164035	50.0	45.0	
105 3-Chlorotoluene	126	11.934	11.930	0.004	94	175084	50.0	42.9	
106 1,3,5-Trimethylbenzene	105	11.958	11.961	-0.003	94	561466	50.0	46.4	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ng	OnCol Amt ng	Flags
107 4-Chlorotoluene	126	11.983	11.985	-0.002	98	183736	50.0	46.5	
108 tert-Butylbenzene	119	12.287	12.283	0.004	93	440939	50.0	42.0	
110 1,2,4-Trimethylbenzene	105	12.336	12.332	0.004	98	561097	50.0	45.1	
111 1,2-dichloro-4-(trifluorom	214	12.402	12.399	0.003	99	160795	50.0	40.9	
112 sec-Butylbenzene	105	12.506	12.502	0.004	95	648966	50.0	43.9	
113 1,3-Dichlorobenzene	146	12.615	12.618	-0.003	98	308813	50.0	47.6	
114 4-Isopropyltoluene	119	12.646	12.648	-0.002	96	522628	50.0	42.9	
115 1,4-Dichlorobenzene	146	12.707	12.703	0.004	94	325063	50.0	49.1	
116 2,4-Dichloro-1-(trifluorom	214	12.755	12.758	-0.003	98	155324	50.0	42.2	
118 2,5-Dichlorobenzotrifluori	214	12.804	12.806	-0.002	0	167755	50.0	40.8	
120 n-Butylbenzene	91	13.059	13.062	-0.003	98	442985	50.0	39.9	
121 1,2-Dichlorobenzene	146	13.078	13.080	-0.002	96	287180	50.0	47.8	
122 1,2-Dibromo-3-Chloropropan	75	13.856	13.853	0.003	76	21004	50.0	42.7	
123 2,4- & 2,5- & 2,6- Dichlor	125	14.002	14.005	-0.003	0	517247	150.0	113.9	
125 2,3- & 3,4- Dichlorotoluen	125	14.422	14.424	-0.002	0	315127	100.0	71.4	
126 1,2,4-Trichlorobenzene	180	14.690	14.686	0.004	94	119977	50.0	38.4	
127 Hexachlorobutadiene	225	14.860	14.856	0.004	96	55585	50.0	37.1	
128 Naphthalene	128	14.939	14.942	-0.003	97	304354	50.0	37.1	
129 1,2,3-Trichlorobenzene	180	15.183	15.185	-0.002	94	98273	50.0	38.3	
131 2,4,5-Trichlorotoluene	159	15.961	15.964	-0.003	0	33493	50.0	24.3	
130 2,3,6-Trichlorotoluene	159	16.059	16.061	-0.002	94	31782	50.0	25.5	
149 3,4-Dichlorotoluene	1	0.000					ND	ND	
148 2,3-Dichlorotoluene	1	0.000					ND	ND	
147 2,4-Dichlorotoluene	1	0.000					ND	ND	
146 2,5-Dichlorotoluene	1	0.000					ND	ND	
150 2,6-Dichlorotoluene	1	0.000					ND	ND	
S 133 Xylenes, Total	106				0		100.0	91.3	
S 134 1,2-Dichloroethene, Total	96				0		100.0	81.4	
S 135 1,3-Dichloropropene, Total	1				0		100.0	104.7	

### QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

### Reagents:

VOAACRO2ND_00007	Amount Added: 6.00	Units: uL	
voaWVA2nd Res_00006	Amount Added: 2.00	Units: uL	
voaW ee2nd_00001	Amount Added: 2.00	Units: uL	
voaWKet2 Rest_00002	Amount Added: 2.00	Units: uL	
VOA8260VOA2ND_00111	Amount Added: 2.00	Units: uL	
VOA8260INT_00031	Amount Added: 2.00	Units: uL	Run Reagent
VOA8260SURR_00033	Amount Added: 2.00	Units: uL	Run Reagent

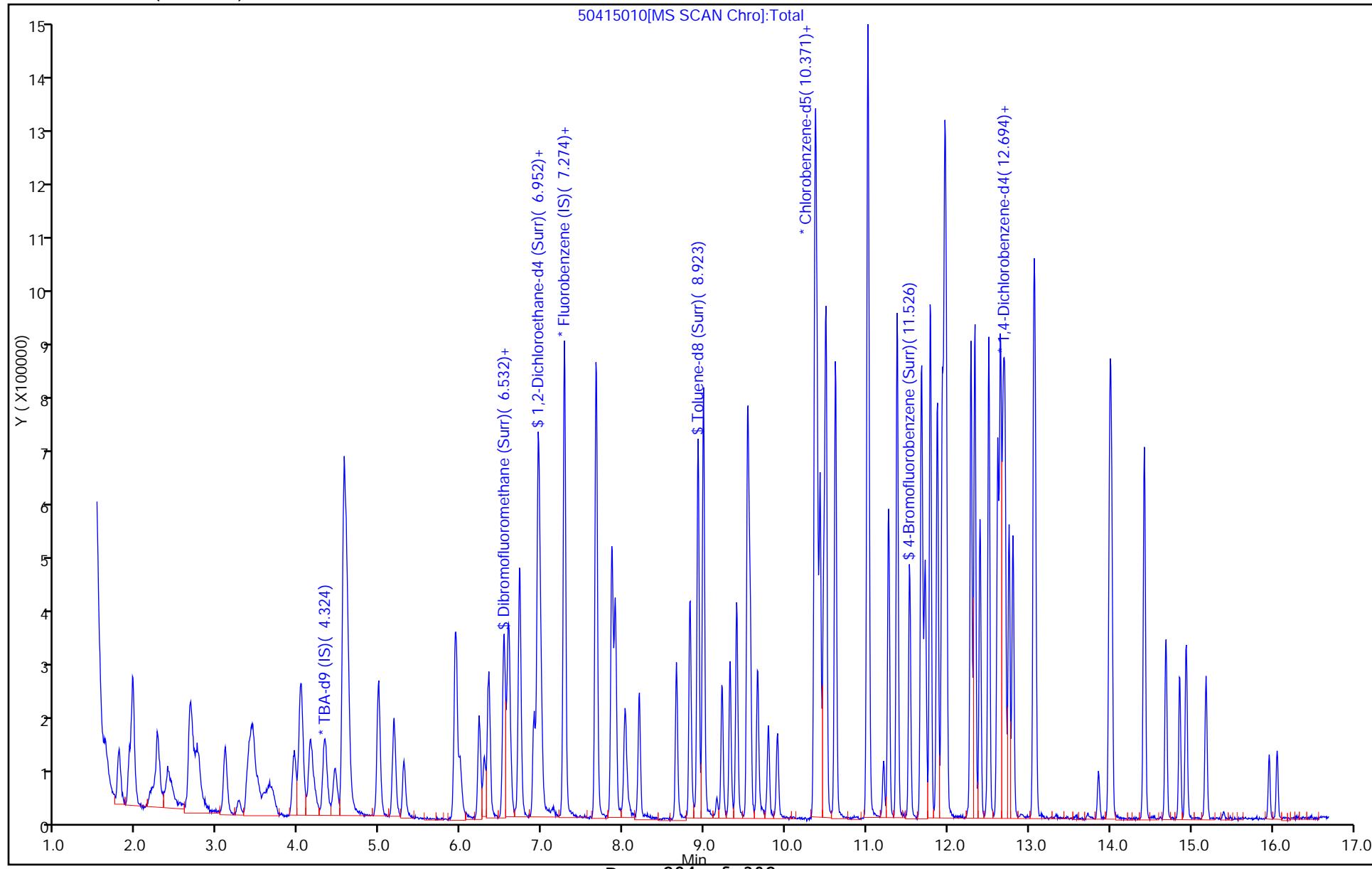
Report Date: 16-Apr-2015 07:37:38

Chrom Revision: 2.2 13-Mar-2015 11:20:44

## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415010.D  
Injection Date: 15-Apr-2015 16:46:30 Instrument ID: CHHP5  
Lims ID: 180-42975-C-1 MSD Operator ID: 001562  
Client ID:  
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 10  
Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
Column: DB-624 ( 0.18 mm)

Worklist Smp#: 10



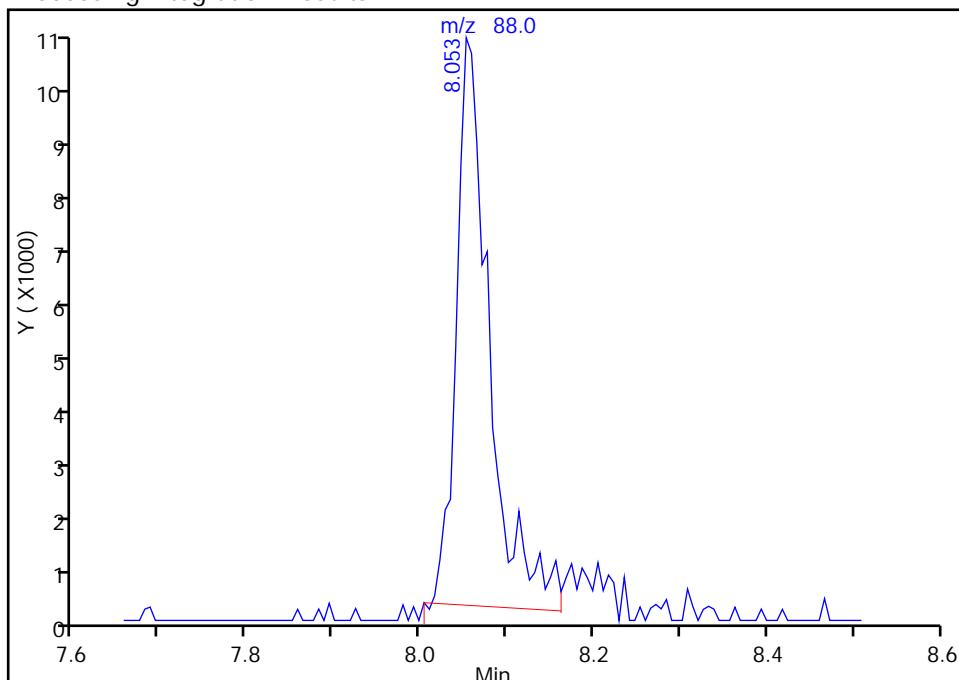
## TestAmerica Pittsburgh

Data File: \\PITCHROM\ChromData\CHHP5\20150415-6480.b\50415010.D  
 Injection Date: 15-Apr-2015 16:46:30 Instrument ID: CHHP5  
 Lims ID: 180-42975-C-1 MSD  
 Client ID:  
 Operator ID: 001562 ALS Bottle#: 10 Worklist Smp#: 10  
 Purge Vol: 5.000 mL Dil. Factor: 1.0000  
 Method: MSVOA\_LL\_CHHP5 Limit Group: VOA 8260C ICAL  
 Column: DB-624 (0.18 mm) Detector: MS SCAN

## 70 1,4-Dioxane, CAS: 123-91-1

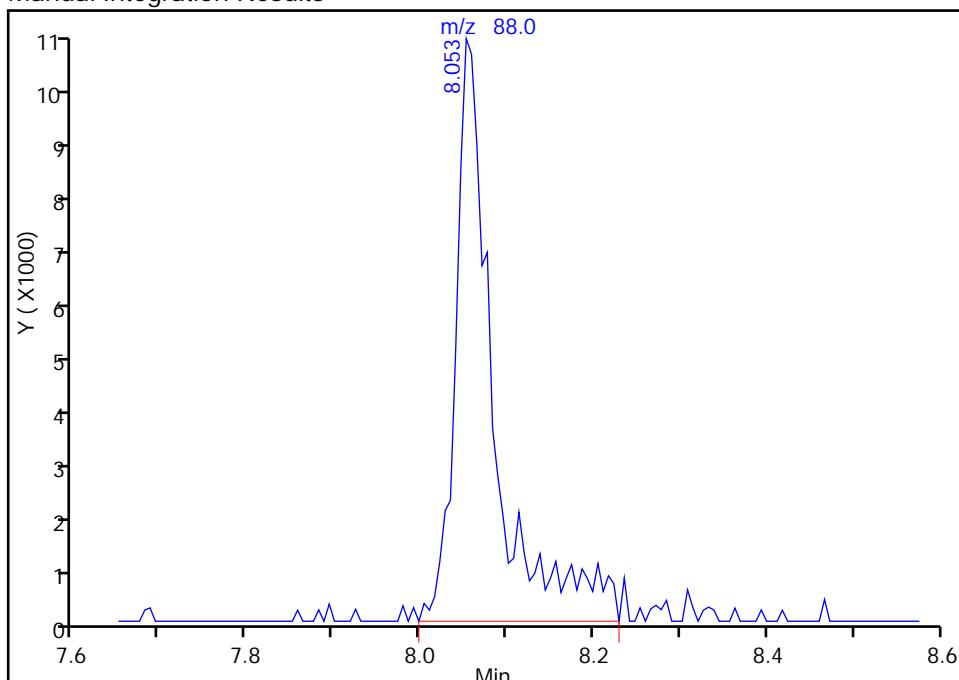
RT: 8.05  
 Area: 27677  
 Amount: 798.4740  
 Amount Units: ng

## Processing Integration Results



RT: 8.05  
 Area: 33017  
 Amount: 952.5316  
 Amount Units: ng

## Manual Integration Results



Reviewer: fergusond, 16-Apr-2015 07:37:38

Audit Action: Manually Integrated

Audit Reason: Peak Tail

## GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Instrument ID: CHHP5 Start Date: 03/16/2015 10:49Analysis Batch Number: 135593 End Date: 03/16/2015 17:05

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
BFB 180-135593/1		03/16/2015 10:49	1	50316001.D	DB-624 0.18 (mm)
IC 180-135593/4		03/16/2015 12:41	1	50316004.D	DB-624 0.18 (mm)
ICIS 180-135593/5		03/16/2015 13:05	1	50316005.D	DB-624 0.18 (mm)
IC 180-135593/6		03/16/2015 13:29	1	50316006.D	DB-624 0.18 (mm)
IC 180-135593/7		03/16/2015 13:53	1	50316007.D	DB-624 0.18 (mm)
IC 180-135593/8		03/16/2015 14:17	1	50316008.D	DB-624 0.18 (mm)
IC 180-135593/9		03/16/2015 14:41	1	50316009.D	DB-624 0.18 (mm)
IC 180-135593/10		03/16/2015 15:05	1	50316010.D	DB-624 0.18 (mm)
IC 180-135593/13		03/16/2015 16:17	1	50316013.D	DB-624 0.18 (mm)
ICV 180-135593/15		03/16/2015 17:05	1		DB-624 0.18 (mm)

## GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica PittsburghJob No.: 180-42975-1

SDG No.:

Instrument ID: CHHP5Start Date: 04/15/2015 12:36Analysis Batch Number: 138583End Date: 04/15/2015 20:47

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
BFB 180-138583/1		04/15/2015 12:36	1	50415001.D	DB-624 0.18 (mm)
CCVIS 180-138583/2		04/15/2015 13:21	1	50415002.D	DB-624 0.18 (mm)
ZZZZZ		04/15/2015 13:21	1		DB-624 0.18 (mm)
CCV 180-138583/3		04/15/2015 13:45	1	50415003.D	DB-624 0.18 (mm)
MB 180-138583/5		04/15/2015 14:33	1	50415005.D	DB-624 0.18 (mm)
180-42975-1	HD-MW-165-0/1-0	04/15/2015 15:09	1	50415006.D	DB-624 0.18 (mm)
180-42975-4	HD-QC2-0/1-2	04/15/2015 15:33	1	50415007.D	DB-624 0.18 (mm)
LCS 180-138583/8		04/15/2015 15:57	1	50415008.D	DB-624 0.18 (mm)
180-42975-1 MS	HD-MW-165-0/1-0 MS	04/15/2015 16:21	1	50415009.D	DB-624 0.18 (mm)
180-42975-1 MSD	HD-MW-165-0/1-0 MSD	04/15/2015 16:46	1	50415010.D	DB-624 0.18 (mm)
ZZZZZ		04/15/2015 17:34	4		DB-624 0.18 (mm)
ZZZZZ		04/15/2015 17:58	1		DB-624 0.18 (mm)
ZZZZZ		04/15/2015 18:46	1		DB-624 0.18 (mm)
ZZZZZ		04/15/2015 19:10	12.5		DB-624 0.18 (mm)
ZZZZZ		04/15/2015 19:35	1		DB-624 0.18 (mm)
ZZZZZ		04/15/2015 19:59	1		DB-624 0.18 (mm)
ZZZZZ		04/15/2015 20:23	1		DB-624 0.18 (mm)
180-42975-2	HD-MW-162-0/1-0	04/15/2015 20:47	4	50415020.D	DB-624 0.18 (mm)

## GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Pittsburgh Job No.: 180-42975-1

SDG No.: \_\_\_\_\_

Instrument ID: CHHP5 Start Date: 04/16/2015 09:31Analysis Batch Number: 138685 End Date: 04/16/2015 16:53

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
BFB 180-138685/1		04/16/2015 09:31	1	50416001.D	DB-624 0.18 (mm)
CCVIS 180-138685/2		04/16/2015 10:19	1	50416002.D	DB-624 0.18 (mm)
ZZZZZ		04/16/2015 10:19	1		DB-624 0.18 (mm)
MB 180-138685/5		04/16/2015 11:32	1	50416005.D	DB-624 0.18 (mm)
LCS 180-138685/8		04/16/2015 13:05	1	50416008.D	DB-624 0.18 (mm)
LCSD 180-138685/10		04/16/2015 13:53	1	50416010.D	DB-624 0.18 (mm)
ZZZZZ		04/16/2015 16:04	2		DB-624 0.18 (mm)
180-42975-2 DL	HD-MW-162-0/1-0 DL	04/16/2015 16:29	40	50416016.D	DB-624 0.18 (mm)
180-42975-3	HD-MW-169-0/1-0	04/16/2015 16:53	1	50416017.D	DB-624 0.18 (mm)

# **Shipping and Receiving Documents**

ORIGIN ID:KPDA (610) 337-9992  
SAMPLE RECEIPT  
TEST AMERICA  
1008 WEST 9TH AVE

SHIP DATE: 10APR15  
ACTWGT: 21.0 LB  
CAD: 8490299/INET3610

KING OF PRUSSIA, PA 19406  
UNITED STATES US

BILL RECIPIENT

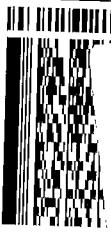
TO **SAMPLE RECEIPT**  
**TEST AMERICA - PITTSBURGH**  
**301 ALPHA DR**

**PITTSBURGH PA 15238**

(412) 963-7058  
INV:  
PO:

REF:

DEPT:



Uncorrected temp  
Thermometer ID

CF C Initials NS

PT-WI-SR-001 effective 7/26/13

2.9 °C

FedEx  
Express



J15172150223011V

TRK#  
0201 7733 4533 7703

**SATURDAY 12:00P**  
**PRIORITY OVERNIGHT**

**XO AGCA**

**15238**  
PA-US R/T

Part # 156297-456 RJT2 02/15

180-42975 Chain of Custody

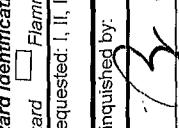


**TestAmerica Pittsburgh**  
301 Alpha Drive RIDC Park  
Pittsburgh, PA 15238  
Phone (412) 963-7058 Fax (412) 963-2468

**TestAmerica**  
THE LEADERS IN ENVIRONMENTAL TESTING

## Chain of Custody Record

**Client Information**

Client Contact Alisa Miller	Jennifer Reese	Sampler: <b>R. Ulrich</b>	Lab PI#: Gamber, Carrie L.	Carrier Tracking No(s): COC No: <b>TAP 20150808</b> Phone: 480-23296-4180
Company: Groundwater Sciences Corporation	Phone: Email: <b>jreese@groundwatersciences.com</b>	Job #: <b>10012-26</b>	Page: <b>1 of 1</b>	
Address: 2501 Market Place Street, Suite 310	TAT Requested (days): <b>5 days</b>	Analysis Requested		
City: Harrisburg	PO #:			
State, Zip: PA, 17110-9307	Purchase Order not required			
Phone: 724-941-4544	WO #:			
Email: amiller@groundwatersciences.com	Project #: 18010144			
Project Name: Harter-Davison SPBA Groundwater Sampling	SSOW #:			
Site: Harley-Davidson, York PA	Sample Date	Sample Time	Sample Type (C=comp, G=grab)	Matrix (W=water, S=solid, O=oil, B=bioassay, A=air)
				Preservation Codes: A - HCl B - NaOH C - Zn Acetate D - Nitric Acid E - NaHSO4 F - MeOH G - Amchlor H - Ascorbic Acid I - Ice J - DI Water K - EDTA L - EDA Other:
				Total Number of Containers: Z - other (specify)
				Special Instructions/Note:
				Performs MSDS (Yes or No)
				Field Filled Sample (Yes or No)
				Designation Code: <b>A</b>
<b>HD-MW-165-01-0</b>	<b>4/10/15</b>	<b>0840</b>	<b>G</b>	<b>W</b>
<b>HD-MW-162-01-0</b>	<b>4/10/15</b>	<b>0950</b>	<b>G</b>	<b>W</b>
<b>HD-MW-169-01-0</b>	<b>4/10/15</b>	<b>1232</b>	<b>G</b>	<b>W</b>
<b>HD-QC2-01-2</b>	<b>4/10/15</b>	<b>1200</b>	<b>X</b>	<b>W</b>
<b>VOCs (8260C)</b>				
<input checked="" type="checkbox"/> Possible Hazard Identification <input type="checkbox"/> Non-Hazard <input type="checkbox"/> Flammable <input type="checkbox"/> Skin Irritant <input type="checkbox"/> Poison B <input type="checkbox"/> Unknown <input type="checkbox"/> Radiological				
Deliverable Requested: I, II, III, IV, Other (specify)				
Empty Kit Relinquished by:   180-42975 Waybill				
Date:	Time:	Company	Received By:	Method of Shipment:
Relinquished by:  	Date/Time: <b>4/10/15 14:20</b>	<b>GSC</b>	Received by:  	Date/Time:
Relinquished by:  	Date/Time: <b>4/10/15 1642</b>	<b>TA</b>	Received by:  	Date/Time: <b>4/10/15 930</b>
Cooler Temperature(s) °C and Other Remarks: △ Yes ▲ No				
Special Instructions/QC Requirements: <b>CLP Like Deliverables</b>				
Sample Disposal / A fee may be assessed if samples are retained longer than 1 month <input type="checkbox"/> Return To Client <input checked="" type="checkbox"/> Disposal By Lab <input type="checkbox"/> Archive For Months				

## Login Sample Receipt Checklist

Client: Groundwater Sciences Corporation

Job Number: 180-42975-1

**Login Number: 42975**

**List Source: TestAmerica Pittsburgh**

**List Number: 1**

**Creator: Lonzo, Michael A**

Question	Answer	Comment
Radioactivity wasn't checked or is </= background as measured by a survey meter.	True	
The cooler's custody seal, if present, is intact.	True	
Sample custody seals, if present, are intact.	True	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
Is the Field Sampler's name present on COC?	True	
There are no discrepancies between the containers received and the COC.	True	
Samples are received within Holding Time.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
Sample Preservation Verified.	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
Containers requiring zero headspace have no headspace or bubble is <6mm (1/4").	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	
Residual Chlorine Checked.	N/A	