

ANALYTICAL REPORT

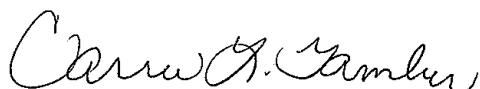
Job Number: 180-71131-1

Job Description: Harley Davidson

For:

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

Attention: Christopher O'Neil



Approved for release.
Carrie L Gamber
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10/19/2017 3:59 PM

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10/19/2017

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Definitions/Glossary

Client: Groundwater Sciences Corporation
Project/Site: Harley Davidson

TestAmerica Job ID: 180-71131-1

Qualifiers

GC/MS VOA

| Qualifier | Qualifier Description |
|-----------|--|
| U | Indicates the analyte was analyzed for but not detected. |
| ^c | CCV Recovery is outside acceptance limits. |
| J | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |

Glossary

| Abbreviation | These commonly used abbreviations may or may not be present in this report. |
|----------------|---|
| ¤ | 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 | Detection Limit (DoD/DOE) |
| DL, RA, RE, IN | Indicates a Dilution, Re-analysis, Re-extraction, or additional Initial metals/anion analysis of the sample |
| DLC | Decision Level Concentration (Radiochemistry) |
| EDL | Estimated Detection Limit (Dioxin) |
| LOD | Limit of Detection (DoD/DOE) |
| LOQ | Limit of Quantitation (DoD/DOE) |
| MDA | Minimum Detectable Activity (Radiochemistry) |
| MDC | Minimum Detectable Concentration (Radiochemistry) |
| 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 (Radiochemistry) |
| 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: 180-71131-1

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 10/10/2017; the samples arrived in good condition, properly preserved and on ice. The temperature of the coolers at receipt was 1.8 C.

The following Trip Blank sample was received with headspace in 2 of 2 vials: HD-QC6-0/1-2 (180-71131-3). The client was emailed. If the analyst determines the headspace is greater than 6 millimeters and a headspace vial is required to be used. An NCM will be written.

The Chain-of-Custody (COC) was incomplete as received: it was not relinquished by the client. The client emailed a picture of the pink copy that was signed. This copy has been added to the log in and reported as part of the final report.

VOLATILES

The continuing calibration verification (CCV) analyzed in batch 180-226148 was outside the method criteria for the following analytes: 4-Methyl-2-pentanone (MIBK), Acetone, Acrylonitrile and Carbon tetrachloride. A CCV standard at or below the reporting limit (RL) was analyzed with the affected samples and found to be acceptable. As indicated in the reference method, sample analysis may proceed; however, any detection for the affected analytes is considered estimated.

Detection Summary

Client: Groundwater Sciences Corporation
Project/Site: Harley Davidson

TestAmerica Job ID: 180-71131-1

Client Sample ID: HD-SPBA-CW-23-0/1-0

Lab Sample ID: 180-71131-1

| Analyte | Result | Qualifier | RL | MDL | Unit | Dil Fac | D | Method | Prep Type |
|-------------------|--------|-----------|-----|------|------|---------|---|--------|-----------|
| Trichloroethene | 4.1 | | 1.0 | 0.69 | ug/L | 1 | | 8260C | Total/NA |
| Tetrachloroethene | 42 | | 1.0 | 0.47 | ug/L | 1 | | 8260C | Total/NA |

Client Sample ID: HD-CW-23-0/1-0

Lab Sample ID: 180-71131-2

| Analyte | Result | Qualifier | RL | MDL | Unit | Dil Fac | D | Method | Prep Type |
|-------------------|--------|-----------|-----|------|------|---------|---|--------|-----------|
| Trichloroethene | 4.4 | | 1.0 | 0.69 | ug/L | 1 | | 8260C | Total/NA |
| Tetrachloroethene | 43 | | 1.0 | 0.47 | ug/L | 1 | | 8260C | Total/NA |

Client Sample ID: HD-QC6-0/1-2

Lab Sample ID: 180-71131-3

| Analyte | Result | Qualifier | RL | MDL | Unit | Dil Fac | D | Method | Prep Type |
|--------------------|--------|-----------|-----|------|------|---------|---|--------|-----------|
| Methylene Chloride | 1.0 | | 1.0 | 0.36 | ug/L | 1 | | 8260C | Total/NA |

This Detection Summary does not include radiochemical test results.

TestAmerica Pittsburgh

Client Sample Results

Client: Groundwater Sciences Corporation
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-71131-1

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

Client Sample ID: HD-SPBA-CW-23-0/1-0

Date Collected: 10/05/17 09:50

Date Received: 10/10/17 09:00

Lab Sample ID: 180-71131-1

Matrix: Water

| Analyte | Result | Qualifier | RL | MDL | Unit | D | Prepared | Analyzed | Dil Fac |
|------------------------------|------------------|------------------|---------------|------|------|-----------------|-----------------|----------------|---------|
| Chloromethane | 1.0 | U | 1.0 | 0.90 | ug/L | | | 10/18/17 06:33 | 1 |
| Vinyl chloride | 1.0 | U | 1.0 | 0.88 | ug/L | | | 10/18/17 06:33 | 1 |
| Bromomethane | 1.0 | U | 1.0 | 0.89 | ug/L | | | 10/18/17 06:33 | 1 |
| Chloroethane | 1.0 | U | 1.0 | 0.90 | ug/L | | | 10/18/17 06:33 | 1 |
| 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.55 | ug/L | | | 10/18/17 06:33 | 1 |
| Acetone | 5.0 | U ^c | 5.0 | 3.4 | ug/L | | | 10/18/17 06:33 | 1 |
| Carbon disulfide | 1.0 | U | 1.0 | 0.88 | ug/L | | | 10/18/17 06:33 | 1 |
| Methylene Chloride | 1.0 | U | 1.0 | 0.36 | ug/L | | | 10/18/17 06:33 | 1 |
| trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.67 | ug/L | | | 10/18/17 06:33 | 1 |
| Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.59 | ug/L | | | 10/18/17 06:33 | 1 |
| 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.63 | ug/L | | | 10/18/17 06:33 | 1 |
| cis-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.71 | ug/L | | | 10/18/17 06:33 | 1 |
| Bromoform | 1.0 | U | 1.0 | 0.63 | ug/L | | | 10/18/17 06:33 | 1 |
| 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 | ug/L | | | 10/18/17 06:33 | 1 |
| Chloroform | 1.0 | U | 1.0 | 0.60 | ug/L | | | 10/18/17 06:33 | 1 |
| 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.60 | ug/L | | | 10/18/17 06:33 | 1 |
| Carbon tetrachloride | 1.0 | U ^c | 1.0 | 0.88 | ug/L | | | 10/18/17 06:33 | 1 |
| Benzene | 1.0 | U | 1.0 | 0.60 | ug/L | | | 10/18/17 06:33 | 1 |
| 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.57 | ug/L | | | 10/18/17 06:33 | 1 |
| Trichloroethene | 4.1 | | 1.0 | 0.69 | ug/L | | | 10/18/17 06:33 | 1 |
| 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.66 | ug/L | | | 10/18/17 06:33 | 1 |
| Bromodichloromethane | 1.0 | U | 1.0 | 0.64 | ug/L | | | 10/18/17 06:33 | 1 |
| cis-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.59 | ug/L | | | 10/18/17 06:33 | 1 |
| 4-Methyl-2-pentanone (MIBK) | 5.0 | U ^c | 5.0 | 3.1 | ug/L | | | 10/18/17 06:33 | 1 |
| Toluene | 1.0 | U | 1.0 | 0.46 | ug/L | | | 10/18/17 06:33 | 1 |
| trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.58 | ug/L | | | 10/18/17 06:33 | 1 |
| 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.45 | ug/L | | | 10/18/17 06:33 | 1 |
| Tetrachloroethene | 42 | | 1.0 | 0.47 | ug/L | | | 10/18/17 06:33 | 1 |
| 2-Hexanone | 5.0 | U | 5.0 | 3.3 | ug/L | | | 10/18/17 06:33 | 1 |
| Dibromochloromethane | 1.0 | U | 1.0 | 0.84 | ug/L | | | 10/18/17 06:33 | 1 |
| 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.50 | ug/L | | | 10/18/17 06:33 | 1 |
| Chlorobenzene | 1.0 | U | 1.0 | 0.50 | ug/L | | | 10/18/17 06:33 | 1 |
| 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.57 | ug/L | | | 10/18/17 06:33 | 1 |
| Ethylbenzene | 1.0 | U | 1.0 | 0.51 | ug/L | | | 10/18/17 06:33 | 1 |
| Xylenes, Total | 2.0 | U | 2.0 | 0.89 | ug/L | | | 10/18/17 06:33 | 1 |
| Styrene | 1.0 | U | 1.0 | 0.47 | ug/L | | | 10/18/17 06:33 | 1 |
| Bromoform | 1.0 | U | 1.0 | 0.98 | ug/L | | | 10/18/17 06:33 | 1 |
| 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.60 | ug/L | | | 10/18/17 06:33 | 1 |
| Acrylonitrile | 20 | U ^c | 20 | 7.8 | ug/L | | | 10/18/17 06:33 | 1 |
| 1,4-Dioxane | 200 | U | 200 | 14 | ug/L | | | 10/18/17 06:33 | 1 |
| Surrogate | %Recovery | Qualifier | Limits | | | Prepared | Analyzed | Dil Fac | |
| 1,2-Dichloroethane-d4 (Surr) | 94 | | 65 - 121 | | | | 10/18/17 06:33 | | 1 |
| Toluene-d8 (Surr) | 86 | | 73 - 120 | | | | 10/18/17 06:33 | | 1 |
| 4-Bromofluorobenzene (Surr) | 91 | | 80 - 120 | | | | 10/18/17 06:33 | | 1 |
| Dibromofluoromethane (Surr) | 93 | | 73 - 120 | | | | 10/18/17 06:33 | | 1 |

TestAmerica Pittsburgh

Client Sample Results

Client: Groundwater Sciences Corporation
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-71131-1

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

Client Sample ID: HD-CW-23-0/1-0

Date Collected: 10/06/17 09:50

Date Received: 10/10/17 09:00

Lab Sample ID: 180-71131-2

Matrix: Water

| Analyte | Result | Qualifier | RL | MDL | Unit | D | Prepared | Analyzed | Dil Fac |
|------------------------------|------------------|------------------|---------------|------|------|-----------------|-----------------|----------------|---------|
| Chloromethane | 1.0 | U | 1.0 | 0.90 | ug/L | | | 10/18/17 06:57 | 1 |
| Vinyl chloride | 1.0 | U | 1.0 | 0.88 | ug/L | | | 10/18/17 06:57 | 1 |
| Bromomethane | 1.0 | U | 1.0 | 0.89 | ug/L | | | 10/18/17 06:57 | 1 |
| Chloroethane | 1.0 | U | 1.0 | 0.90 | ug/L | | | 10/18/17 06:57 | 1 |
| 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.55 | ug/L | | | 10/18/17 06:57 | 1 |
| Acetone | 5.0 | U ^c | 5.0 | 3.4 | ug/L | | | 10/18/17 06:57 | 1 |
| Carbon disulfide | 1.0 | U | 1.0 | 0.88 | ug/L | | | 10/18/17 06:57 | 1 |
| Methylene Chloride | 1.0 | U | 1.0 | 0.36 | ug/L | | | 10/18/17 06:57 | 1 |
| trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.67 | ug/L | | | 10/18/17 06:57 | 1 |
| Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.59 | ug/L | | | 10/18/17 06:57 | 1 |
| 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.63 | ug/L | | | 10/18/17 06:57 | 1 |
| cis-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.71 | ug/L | | | 10/18/17 06:57 | 1 |
| Bromoform | 1.0 | U | 1.0 | 0.63 | ug/L | | | 10/18/17 06:57 | 1 |
| 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 | ug/L | | | 10/18/17 06:57 | 1 |
| 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.60 | ug/L | | | 10/18/17 06:57 | 1 |
| Carbon tetrachloride | 1.0 | U ^c | 1.0 | 0.88 | ug/L | | | 10/18/17 06:57 | 1 |
| Benzene | 1.0 | U | 1.0 | 0.60 | ug/L | | | 10/18/17 06:57 | 1 |
| 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.57 | ug/L | | | 10/18/17 06:57 | 1 |
| Trichloroethene | 4.4 | | 1.0 | 0.69 | ug/L | | | 10/18/17 06:57 | 1 |
| 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.66 | ug/L | | | 10/18/17 06:57 | 1 |
| Bromodichloromethane | 1.0 | U | 1.0 | 0.64 | ug/L | | | 10/18/17 06:57 | 1 |
| cis-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.59 | ug/L | | | 10/18/17 06:57 | 1 |
| 4-Methyl-2-pentanone (MIBK) | 5.0 | U ^c | 5.0 | 3.1 | ug/L | | | 10/18/17 06:57 | 1 |
| Toluene | 1.0 | U | 1.0 | 0.46 | ug/L | | | 10/18/17 06:57 | 1 |
| trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.58 | ug/L | | | 10/18/17 06:57 | 1 |
| 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.45 | ug/L | | | 10/18/17 06:57 | 1 |
| Tetrachloroethene | 43 | | 1.0 | 0.47 | ug/L | | | 10/18/17 06:57 | 1 |
| 2-Hexanone | 5.0 | U | 5.0 | 3.3 | ug/L | | | 10/18/17 06:57 | 1 |
| Dibromochloromethane | 1.0 | U | 1.0 | 0.84 | ug/L | | | 10/18/17 06:57 | 1 |
| 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.50 | ug/L | | | 10/18/17 06:57 | 1 |
| Chlorobenzene | 1.0 | U | 1.0 | 0.50 | ug/L | | | 10/18/17 06:57 | 1 |
| 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.57 | ug/L | | | 10/18/17 06:57 | 1 |
| Ethylbenzene | 1.0 | U | 1.0 | 0.51 | ug/L | | | 10/18/17 06:57 | 1 |
| Xylenes, Total | 2.0 | U | 2.0 | 0.89 | ug/L | | | 10/18/17 06:57 | 1 |
| Styrene | 1.0 | U | 1.0 | 0.47 | ug/L | | | 10/18/17 06:57 | 1 |
| Bromoform | 1.0 | U | 1.0 | 0.98 | ug/L | | | 10/18/17 06:57 | 1 |
| 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.60 | ug/L | | | 10/18/17 06:57 | 1 |
| Acrylonitrile | 20 | U ^c | 20 | 7.8 | ug/L | | | 10/18/17 06:57 | 1 |
| 1,4-Dioxane | 200 | U | 200 | 14 | ug/L | | | 10/18/17 06:57 | 1 |
| Surrogate | %Recovery | Qualifier | Limits | | | Prepared | Analyzed | Dil Fac | |
| 1,2-Dichloroethane-d4 (Surr) | 94 | | 65 - 121 | | | | 10/18/17 06:57 | | 1 |
| Toluene-d8 (Surr) | 89 | | 73 - 120 | | | | 10/18/17 06:57 | | 1 |
| 4-Bromofluorobenzene (Surr) | 91 | | 80 - 120 | | | | 10/18/17 06:57 | | 1 |
| Dibromofluoromethane (Surr) | 93 | | 73 - 120 | | | | 10/18/17 06:57 | | 1 |

TestAmerica Pittsburgh

Client Sample Results

Client: Groundwater Sciences Corporation
Project/Site: Harley Davidson

TestAmerica Job ID: 180-71131-1

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

Client Sample ID: HD-QC6-0/1-2

Date Collected: 10/05/17 12:00

Date Received: 10/10/17 09:00

Lab Sample ID: 180-71131-3

Matrix: Water

| Analyte | Result | Qualifier | RL | MDL | Unit | D | Prepared | Analyzed | Dil Fac |
|------------------------------|------------------|------------------|---------------|------|------|-----------------|-----------------|----------------|---------|
| Chloromethane | 1.0 | U | 1.0 | 0.90 | ug/L | | | 10/18/17 07:45 | 1 |
| Vinyl chloride | 1.0 | U | 1.0 | 0.88 | ug/L | | | 10/18/17 07:45 | 1 |
| Bromomethane | 1.0 | U | 1.0 | 0.89 | ug/L | | | 10/18/17 07:45 | 1 |
| Chloroethane | 1.0 | U | 1.0 | 0.90 | ug/L | | | 10/18/17 07:45 | 1 |
| 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.55 | ug/L | | | 10/18/17 07:45 | 1 |
| Acetone | 5.0 | U ^c | 5.0 | 3.4 | ug/L | | | 10/18/17 07:45 | 1 |
| Carbon disulfide | 1.0 | U | 1.0 | 0.88 | ug/L | | | 10/18/17 07:45 | 1 |
| Methylene Chloride | 1.0 | | 1.0 | 0.36 | ug/L | | | 10/18/17 07:45 | 1 |
| trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.67 | ug/L | | | 10/18/17 07:45 | 1 |
| Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.59 | ug/L | | | 10/18/17 07:45 | 1 |
| 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.63 | ug/L | | | 10/18/17 07:45 | 1 |
| cis-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.71 | ug/L | | | 10/18/17 07:45 | 1 |
| Bromochloromethane | 1.0 | U | 1.0 | 0.63 | ug/L | | | 10/18/17 07:45 | 1 |
| 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 | ug/L | | | 10/18/17 07:45 | 1 |
| Chloroform | 1.0 | U | 1.0 | 0.60 | ug/L | | | 10/18/17 07:45 | 1 |
| 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.60 | ug/L | | | 10/18/17 07:45 | 1 |
| Carbon tetrachloride | 1.0 | U ^c | 1.0 | 0.88 | ug/L | | | 10/18/17 07:45 | 1 |
| Benzene | 1.0 | U | 1.0 | 0.60 | ug/L | | | 10/18/17 07:45 | 1 |
| 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.57 | ug/L | | | 10/18/17 07:45 | 1 |
| Trichloroethene | 1.0 | U | 1.0 | 0.69 | ug/L | | | 10/18/17 07:45 | 1 |
| 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.66 | ug/L | | | 10/18/17 07:45 | 1 |
| Bromodichloromethane | 1.0 | U | 1.0 | 0.64 | ug/L | | | 10/18/17 07:45 | 1 |
| cis-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.59 | ug/L | | | 10/18/17 07:45 | 1 |
| 4-Methyl-2-pentanone (MIBK) | 5.0 | U ^c | 5.0 | 3.1 | ug/L | | | 10/18/17 07:45 | 1 |
| Toluene | 1.0 | U | 1.0 | 0.46 | ug/L | | | 10/18/17 07:45 | 1 |
| trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.58 | ug/L | | | 10/18/17 07:45 | 1 |
| 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.45 | ug/L | | | 10/18/17 07:45 | 1 |
| Tetrachloroethene | 1.0 | U | 1.0 | 0.47 | ug/L | | | 10/18/17 07:45 | 1 |
| 2-Hexanone | 5.0 | U | 5.0 | 3.3 | ug/L | | | 10/18/17 07:45 | 1 |
| Dibromochloromethane | 1.0 | U | 1.0 | 0.84 | ug/L | | | 10/18/17 07:45 | 1 |
| 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.50 | ug/L | | | 10/18/17 07:45 | 1 |
| Chlorobenzene | 1.0 | U | 1.0 | 0.50 | ug/L | | | 10/18/17 07:45 | 1 |
| 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.57 | ug/L | | | 10/18/17 07:45 | 1 |
| Ethylbenzene | 1.0 | U | 1.0 | 0.51 | ug/L | | | 10/18/17 07:45 | 1 |
| Xylenes, Total | 2.0 | U | 2.0 | 0.89 | ug/L | | | 10/18/17 07:45 | 1 |
| Styrene | 1.0 | U | 1.0 | 0.47 | ug/L | | | 10/18/17 07:45 | 1 |
| Bromoform | 1.0 | U | 1.0 | 0.98 | ug/L | | | 10/18/17 07:45 | 1 |
| 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.60 | ug/L | | | 10/18/17 07:45 | 1 |
| Acrylonitrile | 20 | U ^c | 20 | 7.8 | ug/L | | | 10/18/17 07:45 | 1 |
| 1,4-Dioxane | 200 | U | 200 | 14 | ug/L | | | 10/18/17 07:45 | 1 |
| Surrogate | %Recovery | Qualifier | Limits | | | Prepared | Analyzed | Dil Fac | |
| 1,2-Dichloroethane-d4 (Surr) | 95 | | 65 - 121 | | | | 10/18/17 07:45 | | 1 |
| Toluene-d8 (Surr) | 87 | | 73 - 120 | | | | 10/18/17 07:45 | | 1 |
| 4-Bromofluorobenzene (Surr) | 89 | | 80 - 120 | | | | 10/18/17 07:45 | | 1 |
| Dibromofluoromethane (Surr) | 95 | | 73 - 120 | | | | 10/18/17 07:45 | | 1 |

TestAmerica Pittsburgh

Default Detection Limits

Client: Groundwater Sciences Corporation
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-71131-1

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

| Analyte | RL | MDL | Units | Method |
|-----------------------------|-----|------|-------|--------|
| 1,1,1,2-Tetrachloroethane | 1.0 | 0.57 | ug/L | 8260C |
| 1,1,1-Trichloroethane | 1.0 | 0.60 | ug/L | 8260C |
| 1,1,2,2-Tetrachloroethane | 1.0 | 0.60 | ug/L | 8260C |
| 1,1,2-Trichloroethane | 1.0 | 0.45 | ug/L | 8260C |
| 1,1-Dichloroethane | 1.0 | 0.63 | ug/L | 8260C |
| 1,1-Dichloroethene | 1.0 | 0.55 | ug/L | 8260C |
| 1,2-Dibromoethane (EDB) | 1.0 | 0.50 | ug/L | 8260C |
| 1,2-Dichloroethane | 1.0 | 0.57 | ug/L | 8260C |
| 1,2-Dichloropropane | 1.0 | 0.66 | ug/L | 8260C |
| 1,4-Dioxane | 200 | 14 | ug/L | 8260C |
| 2-Butanone (MEK) | 5.0 | 2.6 | ug/L | 8260C |
| 2-Hexanone | 5.0 | 3.3 | ug/L | 8260C |
| 4-Methyl-2-pentanone (MIBK) | 5.0 | 3.1 | ug/L | 8260C |
| Acetone | 5.0 | 3.4 | ug/L | 8260C |
| Acrylonitrile | 20 | 7.8 | ug/L | 8260C |
| Benzene | 1.0 | 0.60 | ug/L | 8260C |
| Bromochloromethane | 1.0 | 0.63 | ug/L | 8260C |
| Bromodichloromethane | 1.0 | 0.64 | ug/L | 8260C |
| Bromoform | 1.0 | 0.98 | ug/L | 8260C |
| Bromomethane | 1.0 | 0.89 | ug/L | 8260C |
| Carbon disulfide | 1.0 | 0.88 | ug/L | 8260C |
| Carbon tetrachloride | 1.0 | 0.88 | ug/L | 8260C |
| Chlorobenzene | 1.0 | 0.50 | ug/L | 8260C |
| Chloroethane | 1.0 | 0.90 | ug/L | 8260C |
| Chloroform | 1.0 | 0.60 | ug/L | 8260C |
| Chloromethane | 1.0 | 0.90 | ug/L | 8260C |
| cis-1,2-Dichloroethene | 1.0 | 0.71 | ug/L | 8260C |
| cis-1,3-Dichloropropene | 1.0 | 0.59 | ug/L | 8260C |
| Dibromochloromethane | 1.0 | 0.84 | ug/L | 8260C |
| Ethylbenzene | 1.0 | 0.51 | ug/L | 8260C |
| Methyl tert-butyl ether | 1.0 | 0.59 | ug/L | 8260C |
| Methylene Chloride | 1.0 | 0.36 | ug/L | 8260C |
| Styrene | 1.0 | 0.47 | ug/L | 8260C |
| Tetrachloroethene | 1.0 | 0.47 | ug/L | 8260C |
| Toluene | 1.0 | 0.46 | ug/L | 8260C |
| trans-1,2-Dichloroethene | 1.0 | 0.67 | ug/L | 8260C |
| trans-1,3-Dichloropropene | 1.0 | 0.58 | ug/L | 8260C |
| Trichloroethene | 1.0 | 0.69 | ug/L | 8260C |
| Vinyl chloride | 1.0 | 0.88 | ug/L | 8260C |
| Xylenes, Total | 2.0 | 0.89 | ug/L | 8260C |

Surrogate Summary

Client: Groundwater Sciences Corporation
Project/Site: Harley Davidson

TestAmerica Job ID: 180-71131-1

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 (65-121) | TOL (73-120) | BFB (80-120) | DBFM (73-120) |
| 180-71131-1 | HD-SPBA-CW-23-0/1-0 | 94 | 86 | 91 | 93 |
| 180-71131-2 | HD-CW-23-0/1-0 | 94 | 89 | 91 | 93 |
| 180-71131-3 | HD-QC6-0/1-2 | 95 | 87 | 89 | 95 |
| LCS 180-226148/4 | Lab Control Sample | 92 | 93 | 94 | 92 |
| MB 180-226148/6 | Method Blank | 93 | 89 | 90 | 91 |

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: 180-71131-1

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

Lab Sample ID: MB 180-226148/6

Matrix: Water

Analysis Batch: 226148

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 | 0.90 | ug/L | 1 |
| Vinyl chloride | 1.0 | U | 1.0 | 0.88 | ug/L | 1 |
| Bromomethane | 1.0 | U | 1.0 | 0.89 | ug/L | 1 |
| Chloroethane | 1.0 | U | 1.0 | 0.90 | ug/L | 1 |
| 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.55 | ug/L | 1 |
| Acetone | 5.0 | U | 5.0 | 3.4 | ug/L | 1 |
| Carbon disulfide | 1.0 | U | 1.0 | 0.88 | ug/L | 1 |
| Methylene Chloride | 1.0 | U | 1.0 | 0.36 | ug/L | 1 |
| trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.67 | ug/L | 1 |
| Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.59 | ug/L | 1 |
| 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.63 | ug/L | 1 |
| cis-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.71 | ug/L | 1 |
| Bromochloromethane | 1.0 | U | 1.0 | 0.63 | ug/L | 1 |
| 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 | ug/L | 1 |
| Chloroform | 1.0 | U | 1.0 | 0.60 | ug/L | 1 |
| 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.60 | ug/L | 1 |
| Carbon tetrachloride | 1.0 | U | 1.0 | 0.88 | ug/L | 1 |
| Benzene | 1.0 | U | 1.0 | 0.60 | ug/L | 1 |
| 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.57 | ug/L | 1 |
| Trichloroethene | 1.0 | U | 1.0 | 0.69 | ug/L | 1 |
| 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.66 | ug/L | 1 |
| Bromodichloromethane | 1.0 | U | 1.0 | 0.64 | ug/L | 1 |
| cis-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.59 | ug/L | 1 |
| 4-Methyl-2-pentanone (MIBK) | 5.0 | U | 5.0 | 3.1 | ug/L | 1 |
| Toluene | 1.0 | U | 1.0 | 0.46 | ug/L | 1 |
| trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.58 | ug/L | 1 |
| 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.45 | ug/L | 1 |
| Tetrachloroethene | 1.0 | U | 1.0 | 0.47 | ug/L | 1 |
| 2-Hexanone | 5.0 | U | 5.0 | 3.3 | ug/L | 1 |
| Dibromochloromethane | 1.0 | U | 1.0 | 0.84 | ug/L | 1 |
| 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.50 | ug/L | 1 |
| Chlorobenzene | 1.0 | U | 1.0 | 0.50 | ug/L | 1 |
| 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.57 | ug/L | 1 |
| Ethylbenzene | 1.0 | U | 1.0 | 0.51 | ug/L | 1 |
| Xylenes, Total | 2.0 | U | 2.0 | 0.89 | ug/L | 1 |
| Styrene | 1.0 | U | 1.0 | 0.47 | ug/L | 1 |
| Bromoform | 1.0 | U | 1.0 | 0.98 | ug/L | 1 |
| 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.60 | ug/L | 1 |
| Acrylonitrile | 20 | U | 20 | 7.8 | ug/L | 1 |
| 1,4-Dioxane | 200 | U | 200 | 14 | ug/L | 1 |

| Surrogate | MB | MB | Limits | Prepared | Analyzed | Dil Fac |
|------------------------------|-----------|-----------|----------|----------|----------------|---------|
| | %Recovery | Qualifier | | | | |
| 1,2-Dichloroethane-d4 (Surr) | 93 | | 65 - 121 | | 10/18/17 02:09 | 1 |
| Toluene-d8 (Surr) | 89 | | 73 - 120 | | 10/18/17 02:09 | 1 |
| 4-Bromofluorobenzene (Surr) | 90 | | 80 - 120 | | 10/18/17 02:09 | 1 |
| Dibromofluoromethane (Surr) | 91 | | 73 - 120 | | 10/18/17 02:09 | 1 |

TestAmerica Pittsburgh

QC Sample Results

Client: Groundwater Sciences Corporation
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-71131-1

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

Lab Sample ID: LCS 180-226148/4

Matrix: Water

Analysis Batch: 226148

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

| Analyte | Spike Added | LCS Result | LCS Qualifier | Unit | D | %Rec | %Rec. Limits |
|-----------------------------|-------------|------------|---------------|------|---|------|--------------|
| Chloromethane | 10.0 | 7.89 | | ug/L | | 79 | 49 - 135 |
| Vinyl chloride | 10.0 | 8.87 | | ug/L | | 89 | 52 - 136 |
| Bromomethane | 10.0 | 7.90 | | ug/L | | 79 | 37 - 150 |
| Chloroethane | 10.0 | 9.09 | | ug/L | | 91 | 44 - 139 |
| 1,1-Dichloroethene | 10.0 | 9.66 | | ug/L | | 97 | 64 - 131 |
| Acetone | 20.0 | 22.4 | | ug/L | | 112 | 24 - 150 |
| Carbon disulfide | 10.0 | 8.89 | | ug/L | | 89 | 20 - 150 |
| Methylene Chloride | 10.0 | 8.98 | | ug/L | | 90 | 66 - 123 |
| trans-1,2-Dichloroethene | 10.0 | 9.35 | | ug/L | | 94 | 70 - 123 |
| Methyl tert-butyl ether | 10.0 | 8.28 | | ug/L | | 83 | 66 - 130 |
| 1,1-Dichloroethane | 10.0 | 9.00 | | ug/L | | 90 | 66 - 122 |
| cis-1,2-Dichloroethene | 10.0 | 9.03 | | ug/L | | 90 | 73 - 120 |
| Bromochloromethane | 10.0 | 9.10 | | ug/L | | 91 | 73 - 122 |
| 2-Butanone (MEK) | 20.0 | 18.3 | | ug/L | | 91 | 37 - 150 |
| Chloroform | 10.0 | 9.63 | | ug/L | | 96 | 72 - 123 |
| 1,1,1-Trichloroethane | 10.0 | 10.4 | | ug/L | | 104 | 66 - 129 |
| Carbon tetrachloride | 10.0 | 12.2 | | ug/L | | 122 | 58 - 145 |
| Benzene | 10.0 | 9.47 | | ug/L | | 95 | 75 - 123 |
| 1,2-Dichloroethane | 10.0 | 9.07 | | ug/L | | 91 | 63 - 130 |
| Trichloroethene | 10.0 | 9.21 | | ug/L | | 92 | 74 - 121 |
| 1,2-Dichloropropane | 10.0 | 8.45 | | ug/L | | 84 | 67 - 119 |
| Bromodichloromethane | 10.0 | 9.26 | | ug/L | | 93 | 62 - 127 |
| cis-1,3-Dichloropropene | 10.0 | 9.17 | | ug/L | | 92 | 61 - 127 |
| 4-Methyl-2-pentanone (MIBK) | 20.0 | 13.2 | | ug/L | | 66 | 41 - 135 |
| Toluene | 10.0 | 9.31 | | ug/L | | 93 | 76 - 129 |
| trans-1,3-Dichloropropene | 10.0 | 9.36 | | ug/L | | 94 | 61 - 136 |
| 1,1,2-Trichloroethane | 10.0 | 8.74 | | ug/L | | 87 | 74 - 126 |
| Tetrachloroethene | 10.0 | 9.13 | | ug/L | | 91 | 76 - 128 |
| 2-Hexanone | 20.0 | 18.9 | | ug/L | | 94 | 37 - 150 |
| Dibromochloromethane | 10.0 | 9.51 | | ug/L | | 95 | 63 - 131 |
| 1,2-Dibromoethane (EDB) | 10.0 | 8.97 | | ug/L | | 90 | 76 - 128 |
| Chlorobenzene | 10.0 | 9.44 | | ug/L | | 94 | 79 - 124 |
| 1,1,1,2-Tetrachloroethane | 10.0 | 10.1 | | ug/L | | 101 | 70 - 130 |
| Ethylbenzene | 10.0 | 9.18 | | ug/L | | 92 | 77 - 124 |
| Xylenes, Total | 20.0 | 18.2 | | ug/L | | 91 | 76 - 124 |
| Styrene | 10.0 | 9.53 | | ug/L | | 95 | 80 - 125 |
| Bromoform | 10.0 | 8.29 | | ug/L | | 83 | 54 - 136 |
| 1,1,2,2-Tetrachloroethane | 10.0 | 8.83 | | ug/L | | 88 | 72 - 128 |
| Acrylonitrile | 100 | 65.5 | | ug/L | | 65 | 60 - 130 |
| 1,4-Dioxane | 200 | 144 | J | ug/L | | 72 | 26 - 150 |

| Surrogate | LCS %Recovery | LCS Qualifier | Limits |
|------------------------------|---------------|---------------|----------|
| 1,2-Dichloroethane-d4 (Surr) | 92 | | 65 - 121 |
| Toluene-d8 (Surr) | 93 | | 73 - 120 |
| 4-Bromofluorobenzene (Surr) | 94 | | 80 - 120 |
| Dibromofluoromethane (Surr) | 92 | | 73 - 120 |

TestAmerica Pittsburgh

QC Association Summary

Client: Groundwater Sciences Corporation
Project/Site: Harley Davidson

TestAmerica Job ID: 180-71131-1

GC/MS VOA

Analysis Batch: 226148

| Lab Sample ID | Client Sample ID | Prep Type | Matrix | Method | Prep Batch |
|------------------|---------------------|-----------|--------|--------|------------|
| 180-71131-1 | HD-SPBA-CW-23-0/1-0 | Total/NA | Water | 8260C | |
| 180-71131-2 | HD-CW-23-0/1-0 | Total/NA | Water | 8260C | |
| 180-71131-3 | HD-QC6-0/1-2 | Total/NA | Water | 8260C | |
| MB 180-226148/6 | Method Blank | Total/NA | Water | 8260C | |
| LCS 180-226148/4 | Lab Control Sample | Total/NA | Water | 8260C | |

Lab Chronicle

Client: Groundwater Sciences Corporation
Project/Site: Harley Davidson

TestAmerica Job ID: 180-71131-1

Client Sample ID: HD-SPBA-CW-23-0/1-0
Date Collected: 10/05/17 09:50
Date Received: 10/10/17 09:00

Lab Sample ID: 180-71131-1
Matrix: Water

| 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 | 226148 | 10/18/17 06:33 | FBB | TAL PIT |

Client Sample ID: HD-CW-23-0/1-0
Date Collected: 10/06/17 09:50
Date Received: 10/10/17 09:00

Lab Sample ID: 180-71131-2
Matrix: Water

| 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 | 226148 | 10/18/17 06:57 | FBB | TAL PIT |

Client Sample ID: HD-QC6-0/1-2
Date Collected: 10/05/17 12:00
Date Received: 10/10/17 09:00

Lab Sample ID: 180-71131-3
Matrix: Water

| 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 | 226148 | 10/18/17 07:45 | FBB | TAL PIT |

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

FBB = Frank Bungard

Accreditation/Certification Summary

Client: Groundwater Sciences Corporation

TestAmerica Job ID: 180-71131-1

Project/Site: Harley Davidson

Laboratory: TestAmerica Pittsburgh

The accreditations/certifications listed below are applicable to this report.

| Authority | Program | EPA Region | Identification Number | Expiration Date |
|--------------|---------|------------|-----------------------|-----------------|
| Pennsylvania | NELAP | 3 | 02-00416 | 04-30-18 |

Method Summary

Client: Groundwater Sciences Corporation
Project/Site: Harley Davidson

TestAmerica Job ID: 180-71131-1

| 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: 180-71131-1

| Lab Sample ID | Client Sample ID | Matrix | Collected | Received |
|---------------|---------------------|--------|----------------|----------------|
| 180-71131-1 | HD-SPBA-CW-23-0/1-0 | Water | 10/05/17 09:50 | 10/10/17 09:00 |
| 180-71131-2 | HD-CW-23-0/1-0 | Water | 10/06/17 09:50 | 10/10/17 09:00 |
| 180-71131-3 | HD-QC6-0/1-2 | Water | 10/05/17 12:00 | 10/10/17 09:00 |

GC/MS VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

Instrument ID: CHHP6

Analysis Batch Number: 217861

Lab Sample ID: IC 180-217861/3

Client Sample ID:

Date Analyzed: 07/24/17 06:39

Lab File ID: 60724D03.D

GC Column: DB-624

ID: 0.18 (mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|------------------------|----------------|---------------------|----------|----------------|
| | | REASON | ANALYST | DATE |
| Trichlorofluoromethane | 2.50 | Poor chromatography | bungardf | 07/24/17 07:17 |

Lab Sample ID: IC 180-217861/4

Client Sample ID:

Date Analyzed: 07/24/17 07:03

Lab File ID: 60724D04.D

GC Column: DB-624

ID: 0.18 (mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---------------|----------------|---------------------|----------|----------------|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 7.80 | Poor chromatography | bungardf | 07/24/17 07:27 |

Lab Sample ID: ICIS 180-217861/5

Client Sample ID:

Date Analyzed: 07/24/17 07:27

Lab File ID: 60724D05.D

GC Column: DB-624

ID: 0.18 (mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---------------|----------------|---------------------|----------|----------------|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 7.81 | Poor chromatography | bungardf | 07/24/17 07:58 |

Lab Sample ID: IC 180-217861/6

Client Sample ID:

Date Analyzed: 07/24/17 07:52

Lab File ID: 60724D06.D

GC Column: DB-624

ID: 0.18 (mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---------------|----------------|---------------------|----------|----------------|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 7.81 | Poor chromatography | bungardf | 07/24/17 08:11 |

Lab Sample ID: IC 180-217861/7

Client Sample ID:

Date Analyzed: 07/24/17 08:16

Lab File ID: 60724D07.D

GC Column: DB-624

ID: 0.18 (mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---------------|----------------|---------------------|----------|----------------|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 7.80 | Poor chromatography | bungardf | 07/24/17 08:44 |

GC/MS VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica PittsburghJob No.: 180-71131-1

SDG No.: _____

Instrument ID: CHHP6Analysis Batch Number: 217861Lab Sample ID: IC 180-217861/9

Client Sample ID: _____

Date Analyzed: 07/24/17 09:04Lab File ID: 60724D09.DGC Column: DB-624 ID: 0.18 (mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---------------|-------------------|---------------------|----------|----------------|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 7.81 | Poor chromatography | bungardf | 07/24/17 09:26 |

Lab Sample ID: ICV 180-217861/13

Client Sample ID: _____

Date Analyzed: 07/24/17 10:40Lab File ID: 60724D13.DGC Column: DB-624 ID: 0.18 (mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---------------|-------------------|---------------------|----------|----------------|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 7.80 | Poor chromatography | bungardf | 07/25/17 01:35 |

GC/MS VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica PittsburghJob No.: 180-71131-1

SDG No.: _____

Instrument ID: CHHP6Analysis Batch Number: 226148Lab Sample ID: CCVIS 180-226148/2

Client Sample ID: _____

Date Analyzed: 10/17/17 23:55Lab File ID: 6101802D.DGC Column: DB-624ID: 0.18 (mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---------------|-------------------|---------------------|----------|----------------|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 7.79 | Poor chromatography | bungardf | 10/18/17 00:15 |

Lab Sample ID: 180-71131-1Client Sample ID: HD-SPBA-CW-23-0/1-0Date Analyzed: 10/18/17 06:33Lab File ID: 6101816D.DGC Column: DB-624ID: 0.18 (mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---------------|-------------------|---------------------|----------|----------------|
| | | REASON | ANALYST | DATE |
| Acetone | 3.19 | Poor chromatography | bungardf | 10/18/17 20:22 |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|----------------------|----------|-----------|-----------------------|----------------------|---------------------|--------------|------------------------------|---------------|
| | | | | | Reagent ID | Volume Added | | |
| VOA8260INT_00072 | 08/21/17 | 07/21/17 | Methanol, Lot 2019055 | 10 mL | VOA8260INTRES_00123 | 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_00123 | 08/31/20 | | Restek, Lot A0113246 | | (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 |
| VOA8260INT_00074 | 10/20/17 | 09/20/17 | Methanol, Lot 2469125 | 10 mL | VOA8260INTRES_00135 | 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_00135 | 01/31/22 | | Restek, Lot A0124343 | | (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_00071 | 08/21/17 | 07/21/17 | Methanol, Lot 2019055 | 100 mL | VOA8260SURRES_00118 | 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_00118 | 10/31/20 | | Restek, Lot A0114901 | | (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_00073 | 10/20/17 | 09/20/17 | Methanol, Lot 2469125 | 100 mL | VOA8260SURRES_00122 | 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_00122 | 10/31/20 | | Restek, Lot A0114901 | | (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_00253 | 08/10/17 | 07/10/17 | Methanol, Lot 2019054 | 10 mL | VOA8260GAS2ND_00200 | 0.1 mL | Bromomethane | 25 ug/mL |
| | | | | | | | Chloroethane | 25 ug/mL |
| | | | | | | | Chloromethane | 25 ug/mL |
| | | | | | | | Vinyl chloride | 25 ug/mL |
| | | | | | VOA8260VOA2ND_00252 | 1 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 |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|----------------------|----------|----------------------|-----------------------|----------------------|---------------------|--------------|---------------------------|---------------|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 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 |
| .VOA8260GAS2ND_00200 | 01/31/20 | Restek, Lot A0124116 | | | (Purchased Reagent) | | Bromomethane | 2500 ug/mL |
| .VOA8260VOA2ND_00252 | 08/10/17 | 07/10/17 | Methanol, Lot 2019055 | 10 mL | VOA8260MEGA2_00061 | 1 mL | 1,1,1,2-Tetrachloroethane | 250 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 250 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 250 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 250 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 250 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 250 ug/mL |
| | | | | | | | 1,2-Dibromoethane (EDB) | 250 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 250 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 250 ug/mL |
| | | | | | | | 1,4-Dioxane | 5000 ug/mL |
| | | | | | | | Acrylonitrile | 2500 ug/mL |
| | | | | | | | Benzene | 250 ug/mL |
| | | | | | | | Bromochloromethane | 250 ug/mL |
| | | | | | | | Bromodichloromethane | 250 ug/mL |
| | | | | | | | Bromoform | 250 ug/mL |
| | | | | | | | Carbon disulfide | 250 ug/mL |
| | | | | | | | Carbon tetrachloride | 250 ug/mL |
| | | | | | | | Chlorobenzene | 250 ug/mL |
| | | | | | | | Chloroform | 250 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 250 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 250 ug/mL |
| | | | | | | | Dibromochloromethane | 250 ug/mL |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|----------------------|----------|----------------------|-----------------------|----------------------|---------------------|--------------|---------------------------|---------------|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Ethylbenzene | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 250 ug/mL |
| | | | | | | | Methylene Chloride | 250 ug/mL |
| | | | | | | | Styrene | 250 ug/mL |
| | | | | | | | Tetrachloroethene | 250 ug/mL |
| | | | | | | | Toluene | 250 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 250 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 250 ug/mL |
| | | | | | | | Trichloroethene | 250 ug/mL |
| | | | | | | | Xylenes, Total | 500 ug/mL |
| ..VOA8260MEGA2_00061 | 12/31/18 | Restek, Lot A0123775 | | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2500 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2500 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2500 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 2500 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2500 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2500 ug/mL |
| | | | | | | | 1,2-Dibromoethane (EDB) | 2500 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2500 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2500 ug/mL |
| | | | | | | | 1,4-Dioxane | 50000 ug/mL |
| | | | | | | | Acrylonitrile | 25000 ug/mL |
| | | | | | | | Benzene | 2500 ug/mL |
| | | | | | | | Bromochloromethane | 2500 ug/mL |
| | | | | | | | Bromodichloromethane | 2500 ug/mL |
| | | | | | | | Bromoform | 2500 ug/mL |
| | | | | | | | Carbon disulfide | 2500 ug/mL |
| | | | | | | | Carbon tetrachloride | 2500 ug/mL |
| | | | | | | | Chlorobenzene | 2500 ug/mL |
| | | | | | | | Chloroform | 2500 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2500 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2500 ug/mL |
| | | | | | | | Dibromochloromethane | 2500 ug/mL |
| | | | | | | | Ethylbenzene | 2500 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2500 ug/mL |
| | | | | | | | Methylene Chloride | 2500 ug/mL |
| | | | | | | | Styrene | 2500 ug/mL |
| | | | | | | | Tetrachloroethene | 2500 ug/mL |
| | | | | | | | Toluene | 2500 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2500 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2500 ug/mL |
| | | | | | | | Trichloroethene | 2500 ug/mL |
| | | | | | | | Xylenes, Total | 5000 ug/mL |
| VOA8260VOAPRI_00263 | 07/29/17 | 07/22/17 | Methanol, Lot 2019055 | 10 mL | VOA8260GAS1ST_00203 | 0.1 mL | Bromomethane | 25 ug/mL |
| | | | | | | | Butadiene | 25 ug/mL |
| | | | | | | | Chloroethane | 25 ug/mL |
| | | | | | | | Chloromethane | 25 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 25 ug/mL |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|------------|----------|-----------|---------------|----------------------|---------------------|--------------|---------------------------------------|---------------|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | VOA8260VOAPRI_00260 | 1 mL | Trichlorofluoromethane | 25 ug/mL |
| | | | | | | | Vinyl chloride | 25 ug/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 |
| | | | | | | | 1,1,1-Trichloroethane | 25 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 25 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 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 |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|----------------------|----------|----------------------|-----------------------|----------------------|---------------------|--------------|--|---------------|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 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 |
| | | | | | | | Iodomethane | 25 ug/mL |
| | | | | | | | Isobutyl alcohol | 625 ug/mL |
| | | | | | | | Isopropylbenzene | 25 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 25 ug/mL |
| | | | | | | | Methyl acetate | 50 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_00203 | 01/31/20 | Restek, Lot A0124278 | | | (Purchased Reagent) | | Bromomethane | 2500 ug/mL |
| .VOA8260VOAPRI_00260 | 08/06/17 | 07/06/17 | Methanol, Lot 2019056 | 10 mL | VOA8260KET1ST_00100 | 0.2 mL | Butadiene | 2500 ug/mL |
| | | | | | | | Chloroethane | 2500 ug/mL |
| | | | | | | | Chloromethane | 2500 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2500 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Vinyl chloride | 2500 ug/mL |
| | | | | | | | 2-Butanone (MEK) | 250 ug/mL |
| | | | | | | | 2-Hexanone | 250 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 250 ug/mL |
| | | | | | | | Acetone | 250 ug/mL |
| | | | | | VOA8260MEGA1_00065 | 1 mL | 1,1,1,2-Tetrachloroethane | 250 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 250 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 250 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluorooethane | 250 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 250 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 250 ug/mL |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|------------|----------|-----------|---------------|----------------------|----------------|--------------|-----------------------------|---------------|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,1-Dichloroethene | 250 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 250 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 250 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 250 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 250 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 250 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 250 ug/mL |
| | | | | | | | 1,2-Dibromoethane (EDB) | 250 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 250 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 250 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 250 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 250 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 250 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 250 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 250 ug/mL |
| | | | | | | | 1,4-Dioxane | 5000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 250 ug/mL |
| | | | | | | | 2-Chlorotoluene | 250 ug/mL |
| | | | | | | | 2-Methyl-2-propanol | 2500 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 250 ug/mL |
| | | | | | | | 4-Chlorotoluene | 250 ug/mL |
| | | | | | | | 4-Isopropyltoluene | 250 ug/mL |
| | | | | | | | Acrylonitrile | 2500 ug/mL |
| | | | | | | | Benzene | 250 ug/mL |
| | | | | | | | Bromobenzene | 250 ug/mL |
| | | | | | | | Bromochloromethane | 250 ug/mL |
| | | | | | | | Bromodichloromethane | 250 ug/mL |
| | | | | | | | Bromoform | 250 ug/mL |
| | | | | | | | Carbon disulfide | 250 ug/mL |
| | | | | | | | Carbon tetrachloride | 250 ug/mL |
| | | | | | | | Chlorobenzene | 250 ug/mL |
| | | | | | | | Chloroform | 250 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 250 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 250 ug/mL |
| | | | | | | | Cyclohexane | 250 ug/mL |
| | | | | | | | Dibromochloromethane | 250 ug/mL |
| | | | | | | | Dibromomethane | 250 ug/mL |
| | | | | | | | Ethyl ether | 250 ug/mL |
| | | | | | | | Ethyl methacrylate | 250 ug/mL |
| | | | | | | | Ethylbenzene | 250 ug/mL |
| | | | | | | | Hexachlorobutadiene | 250 ug/mL |
| | | | | | | | Hexane | 250 ug/mL |
| | | | | | | | Iodomethane | 250 ug/mL |
| | | | | | | | Isobutyl alcohol | 6250 ug/mL |
| | | | | | | | Isopropylbenzene | 250 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 250 ug/mL |
| | | | | | | | Methyl acetate | 500 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 250 ug/mL |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration | | | |
|-----------------------|----------|----------------------|---------------|----------------------|----------------|---------------------|--|---------------|--|--|--|
| | | | | | Reagent ID | Volume Added | | | | | |
| ..VOA8260KET1ST_00100 | | | | | | | Methylcyclohexane | 250 ug/mL | | | |
| | | | | | | | Methylene Chloride | 250 ug/mL | | | |
| | | | | | | | n-Butylbenzene | 250 ug/mL | | | |
| | | | | | | | n-Heptane | 250 ug/mL | | | |
| | | | | | | | N-Propylbenzene | 250 ug/mL | | | |
| | | | | | | | Naphthalene | 250 ug/mL | | | |
| | | | | | | | o-Xylene | 250 ug/mL | | | |
| | | | | | | | sec-Butylbenzene | 250 ug/mL | | | |
| | | | | | | | Styrene | 250 ug/mL | | | |
| | | | | | | | tert-Butylbenzene | 250 ug/mL | | | |
| | | | | | | | Tetrachloroethene | 250 ug/mL | | | |
| | | | | | | | Tetrahydrofuran | 500 ug/mL | | | |
| | | | | | | | Toluene | 250 ug/mL | | | |
| | | | | | | | trans-1,2-Dichloroethene | 250 ug/mL | | | |
| | | | | | | | trans-1,3-Dichloropropene | 250 ug/mL | | | |
| | | | | | | | trans-1,4-Dichloro-2-butene | 250 ug/mL | | | |
| | | | | | | | Trichloroethene | 250 ug/mL | | | |
| | | Restek, Lot A0123890 | | | | (Purchased Reagent) | 2-Butanone (MEK) | 12500 ug/mL | | | |
| | | | | | | | 2-Hexanone | 12500 ug/mL | | | |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 12500 ug/mL | | | |
| | | | | | | | Acetone | 12500 ug/mL | | | |
| ..VOA8260MEGA1_00065 | | Restek, Lot A0123711 | | | | (Purchased Reagent) | 1,1,1,2-Tetrachloroethane | 2500 ug/mL | | | |
| | | | | | | | 1,1,1-Trichloroethane | 2500 ug/mL | | | |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2500 ug/mL | | | |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluor oethane | 2500 ug/mL | | | |
| | | | | | | | 1,1,2-Trichloroethane | 2500 ug/mL | | | |
| | | | | | | | 1,1-Dichloroethane | 2500 ug/mL | | | |
| | | | | | | | 1,1-Dichloroethene | 2500 ug/mL | | | |
| | | | | | | | 1,1-Dichloropropene | 2500 ug/mL | | | |
| | | | | | | | 1,2,3-Trichlorobenzene | 2500 ug/mL | | | |
| | | | | | | | 1,2,3-Trichloropropane | 2500 ug/mL | | | |
| | | | | | | | 1,2,4-Trichlorobenzene | 2500 ug/mL | | | |
| | | | | | | | 1,2,4-Trimethylbenzene | 2500 ug/mL | | | |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2500 ug/mL | | | |
| | | | | | | | 1,2-Dibromoethane (EDB) | 2500 ug/mL | | | |
| | | | | | | | 1,2-Dichlorobenzene | 2500 ug/mL | | | |
| | | | | | | | 1,2-Dichloroethane | 2500 ug/mL | | | |
| | | | | | | | 1,2-Dichloropropane | 2500 ug/mL | | | |
| | | | | | | | 1,3,5-Trimethylbenzene | 2500 ug/mL | | | |
| | | | | | | | 1,3-Dichlorobenzene | 2500 ug/mL | | | |
| | | | | | | | 1,3-Dichloropropane | 2500 ug/mL | | | |
| | | | | | | | 1,4-Dichlorobenzene | 2500 ug/mL | | | |
| | | | | | | | 1,4-Dioxane | 50000 ug/mL | | | |
| | | | | | | | 2,2-Dichloropropane | 2500 ug/mL | | | |
| | | | | | | | 2-Chlorotoluene | 2500 ug/mL | | | |
| | | | | | | | 2-Methyl-2-propanol | 25000 ug/mL | | | |
| | | | | | | | 3-Chloro-1-propene | 2500 ug/mL | | | |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---------------------|----------|-----------|-----------------------|----------------------|-----------------------------|--------------|--------------|---------------|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | 4-Chlorotoluene | 2500 ug/mL | | |
| | | | | | 4-Isopropyltoluene | 2500 ug/mL | | |
| | | | | | Acrylonitrile | 25000 ug/mL | | |
| | | | | | Benzene | 2500 ug/mL | | |
| | | | | | Bromobenzene | 2500 ug/mL | | |
| | | | | | Bromochloromethane | 2500 ug/mL | | |
| | | | | | Bromodichloromethane | 2500 ug/mL | | |
| | | | | | Bromoform | 2500 ug/mL | | |
| | | | | | Carbon disulfide | 2500 ug/mL | | |
| | | | | | Carbon tetrachloride | 2500 ug/mL | | |
| | | | | | Chlorobenzene | 2500 ug/mL | | |
| | | | | | Chloroform | 2500 ug/mL | | |
| | | | | | cis-1,2-Dichloroethene | 2500 ug/mL | | |
| | | | | | cis-1,3-Dichloropropene | 2500 ug/mL | | |
| | | | | | Cyclohexane | 2500 ug/mL | | |
| | | | | | Dibromochloromethane | 2500 ug/mL | | |
| | | | | | Dibromomethane | 2500 ug/mL | | |
| | | | | | Ethyl ether | 2500 ug/mL | | |
| | | | | | Ethyl methacrylate | 2500 ug/mL | | |
| | | | | | Ethylbenzene | 2500 ug/mL | | |
| | | | | | Hexachlorobutadiene | 2500 ug/mL | | |
| | | | | | Hexane | 2500 ug/mL | | |
| | | | | | Iodomethane | 2500 ug/mL | | |
| | | | | | Isobutyl alcohol | 62500 ug/mL | | |
| | | | | | Isopropylbenzene | 2500 ug/mL | | |
| | | | | | m-Xylene & p-Xylene | 2500 ug/mL | | |
| | | | | | Methyl acetate | 5000 ug/mL | | |
| | | | | | Methyl tert-butyl ether | 2500 ug/mL | | |
| | | | | | Methylcyclohexane | 2500 ug/mL | | |
| | | | | | Methylene Chloride | 2500 ug/mL | | |
| | | | | | n-Butylbenzene | 2500 ug/mL | | |
| | | | | | n-Heptane | 2500 ug/mL | | |
| | | | | | N-Propylbenzene | 2500 ug/mL | | |
| | | | | | Naphthalene | 2500 ug/mL | | |
| | | | | | o-Xylene | 2500 ug/mL | | |
| | | | | | sec-Butylbenzene | 2500 ug/mL | | |
| | | | | | Styrene | 2500 ug/mL | | |
| | | | | | tert-Butylbenzene | 2500 ug/mL | | |
| | | | | | Tetrachloroethene | 2500 ug/mL | | |
| | | | | | Tetrahydrofuran | 5000 ug/mL | | |
| | | | | | Toluene | 2500 ug/mL | | |
| | | | | | trans-1,2-Dichloroethene | 2500 ug/mL | | |
| | | | | | trans-1,3-Dichloropropene | 2500 ug/mL | | |
| | | | | | trans-1,4-Dichloro-2-butene | 2500 ug/mL | | |
| | | | | | Trichloroethene | 2500 ug/mL | | |
| VOA8260VOAPRI_00266 | 10/21/17 | 10/14/17 | Methanol, Lot 2469120 | 10 mL | VOA8260GAS1ST_00205 | 0.1 mL | Bromomethane | 25 ug/mL |
| | | | | | | | Chloroethane | 25 ug/mL |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration | | | |
|------------|----------|----------------------|-----------------------|----------------------|---------------------|--------------|---------------------------|---------------|--|--|--|
| | | | | | Reagent ID | Volume Added | | | | | |
| | | | | | VOA8260VOAPRI_00264 | 1 mL | Chloromethane | 25 ug/mL | | | |
| | | | | | | | Vinyl chloride | 25 ug/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 | | | |
| | | | | | | | 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 | | | |
| | 01/31/20 | Restek, Lot A0124278 | | | (Purchased Reagent) | | Tetrachloroethene | 25 ug/mL | | | |
| | | | | | | | Toluene | 25 ug/mL | | | |
| | | | | | | | trans-1,2-Dichloroethene | 25 ug/mL | | | |
| | | | | | | | trans-1,3-Dichloropropene | 25 ug/mL | | | |
| | 11/06/17 | 10/06/17 | Methanol, Lot 2469120 | 10 mL | VOA8260MEGA1_00066 | 1 mL | Trichloroethene | 25 ug/mL | | | |
| | | | | | | | Xylenes, Total | 50 ug/mL | | | |
| | | | | | | | Bromomethane | 2500 ug/mL | | | |
| | | | | | | | Chloroethane | 2500 ug/mL | | | |
| | | | | | | | Chloromethane | 2500 ug/mL | | | |
| | | | | | | | Vinyl chloride | 2500 ug/mL | | | |
| | | | | | | | 1,1,1,2-Tetrachloroethane | 250 ug/mL | | | |
| | | | | | | | 1,1,1-Trichloroethane | 250 ug/mL | | | |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 250 ug/mL | | | |
| | | | | | | | 1,1,2-Trichloroethane | 250 ug/mL | | | |
| | | | | | | | 1,1-Dichloroethane | 250 ug/mL | | | |
| | | | | | | | 1,1-Dichloroethene | 250 ug/mL | | | |
| | | | | | | | 1,2-Dibromoethane (EDB) | 250 ug/mL | | | |
| | | | | | | | 1,2-Dichloroethane | 250 ug/mL | | | |
| | | | | | | | 1,2-Dichloropropane | 250 ug/mL | | | |
| | | | | | | | 1,4-Dioxane | 5000 ug/mL | | | |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|----------------------|----------|----------------------|---------------|----------------------|---------------------|--------------|---------------------------|---------------|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Acrylonitrile | 2500 ug/mL |
| | | | | | | | Benzene | 250 ug/mL |
| | | | | | | | Bromochloromethane | 250 ug/mL |
| | | | | | | | Bromodichloromethane | 250 ug/mL |
| | | | | | | | Bromoform | 250 ug/mL |
| | | | | | | | Carbon disulfide | 250 ug/mL |
| | | | | | | | Carbon tetrachloride | 250 ug/mL |
| | | | | | | | Chlorobenzene | 250 ug/mL |
| | | | | | | | Chloroform | 250 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 250 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 250 ug/mL |
| | | | | | | | Dibromochloromethane | 250 ug/mL |
| | | | | | | | Ethylbenzene | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 250 ug/mL |
| | | | | | | | Methylene Chloride | 250 ug/mL |
| | | | | | | | Styrene | 250 ug/mL |
| | | | | | | | Tetrachloroethene | 250 ug/mL |
| | | | | | | | Toluene | 250 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 250 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 250 ug/mL |
| | | | | | | | Trichloroethene | 250 ug/mL |
| | | | | | | | Xylenes, Total | 500 ug/mL |
| ..VOA8260MEGA1_00066 | 12/31/18 | Restek, Lot A0123711 | | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2500 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2500 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2500 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 2500 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2500 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2500 ug/mL |
| | | | | | | | 1,2-Dibromoethane (EDB) | 2500 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2500 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2500 ug/mL |
| | | | | | | | 1,4-Dioxane | 50000 ug/mL |
| | | | | | | | Acrylonitrile | 25000 ug/mL |
| | | | | | | | Benzene | 2500 ug/mL |
| | | | | | | | Bromochloromethane | 2500 ug/mL |
| | | | | | | | Bromodichloromethane | 2500 ug/mL |
| | | | | | | | Bromoform | 2500 ug/mL |
| | | | | | | | Carbon disulfide | 2500 ug/mL |
| | | | | | | | Carbon tetrachloride | 2500 ug/mL |
| | | | | | | | Chlorobenzene | 2500 ug/mL |
| | | | | | | | Chloroform | 2500 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2500 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2500 ug/mL |
| | | | | | | | Dibromochloromethane | 2500 ug/mL |
| | | | | | | | Ethylbenzene | 2500 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2500 ug/mL |
| | | | | | | | Methylene Chloride | 2500 ug/mL |
| | | | | | | | Styrene | 2500 ug/mL |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|-----------------------------|----------|-----------|-----------------------|----------------------|---------------------|--------------|--|---------------|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Tetrachloroethene | 2500 ug/mL |
| | | | | | | | Toluene | 2500 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2500 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2500 ug/mL |
| | | | | | | | Trichloroethene | 2500 ug/mL |
| | | | | | | | Xylenes, Total | 5000 ug/mL |
| VOABFB25_00088 | | | | | | | 1,2-Dichloroethene, Total | |
| | | | | | | | 1,3-Dichloropropene, Total | |
| | | | | | | | Tentatively Identified Compound | |
| | | | | | | | Total BTEX | |
| | | | | | | | Xylenes, Total | |
| | | | | | VOABFB50_00091 | 5 mL | BFB | 25 ug/mL |
| .VOABFB50_00091 | 06/09/17 | 05/09/17 | Methanol, Lot 136118 | 50 mL | VOABFBRES_00052 | 1 mL | BFB | 50 ug/mL |
| ..VOABFBRES_00052 | 05/31/21 | | Restek, Lot A0119122 | | (Purchased Reagent) | | BFB | 2500 ug/mL |
| VOABFB25_00090 | | | | | | | 1,2-Dichloroethene, Total | |
| | | | | | | | 1,3-Dichloropropene, Total | |
| | | | | | | | Tentatively Identified Compound | |
| | | | | | | | Total BTEX | |
| | | | | | | | Xylenes, Total | |
| | | | | | VOABFB50_00093 | 5 mL | BFB | 25 ug/mL |
| .VOABFB50_00093 | 08/10/17 | 07/10/17 | Methanol, Lot 2019056 | 50 mL | VOABFBRES_00058 | 1 mL | BFB | 50 ug/mL |
| ..VOABFBRES_00058 | 11/30/21 | | Restek, Lot A0122647 | | (Purchased Reagent) | | BFB | 2500 ug/mL |
| VOABFB25_00094 | | | | | | | 1,2-Dichloroethene, Total | |
| | | | | | | | 1,3-Dichloropropene, Total | |
| | | | | | | | Tentatively Identified Compound | |
| | | | | | | | Total BTEX | |
| | | | | | | | Xylenes, Total | |
| | | | | | VOABFB50_00096 | 5 mL | BFB | 25 ug/mL |
| .VOABFB50_00096 | 11/09/17 | 10/09/17 | Methanol, Lot 2469125 | 50 mL | VOABFBRES_00055 | 1 mL | BFB | 50 ug/mL |
| ..VOABFBRES_00055 | 11/30/21 | | Restek, Lot A0122647 | | (Purchased Reagent) | | BFB | 2500 ug/mL |
| voaW2clev1stR_00013 | 07/31/17 | 07/24/17 | Methanol, Lot 2019056 | 10 mL | VOACEVERES_00127 | 200 uL | 2-Chloroethyl vinyl ether | 50 ug/mL |
| .VOACEVERES_00127 | 01/31/20 | | Restek, Lot A0123891 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2500 ug/mL |
| voaWAcrol1stRe_00016 | 08/17/17 | 07/17/17 | Methanol, Lot 2019056 | 100 mL | VOAACRORES_00115 | 0.125 mL | Acrolein | 25 ug/mL |
| .VOAACRORES_00115 | 09/30/17 | | Restek, Lot A0125560 | | (Purchased Reagent) | | Acrolein | 20000 ug/mL |
| voaWEEmix1stR_00009 | 08/03/17 | 07/03/17 | Methanol, Lot 127999 | 25 mL | VOARESEE1ST_00045 | 0.125 mL | 1,2-dichloro-4-(trifluoromethyl)benzene | 25 ug/mL |
| | | | | | | | 2,3,6-Trichlorotoluene | 25 ug/mL |
| | | | | | | | 2,3- & 3,4- Dichlorotoluene | 50 ug/mL |
| | | | | | | | 2,4,5-Trichlorotoluene | 25 ug/mL |
| | | | | | | | 2,4- & 2,5- & 2,6- Dichlorotoluene | 75 ug/mL |
| | | | | | | | 2,4-Dichloro-1-(trifluoromethyl)-benzene | 25 ug/mL |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|----------------------|----------|----------------------|-----------------------|----------------------|---------------------|--------------|--|---------------|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2,5-Dichlorobenzotrifluoride | 25 ug/mL |
| | | | | | | | 2-Chlorobenzotrifluoride | 25 ug/mL |
| | | | | | | | 3-Chlorobenzotrifluoride | 25 ug/mL |
| | | | | | | | 3-Chlorotoluene | 25 ug/mL |
| | | | | | | | 4-Chlorobenzotrifluoride | 25 ug/mL |
| .VOARESEE1ST_00045 | 01/31/18 | Restek, Lot A0120234 | | | (Purchased Reagent) | | 1,2-dichloro-4-(trifluoromethyl)benzene | 5000 ug/mL |
| | | | | | | | 2,3,6-Trichlorotoluene | 5000 ug/mL |
| | | | | | | | 2,3- & 3,4- Dichlorotoluene | 10000 ug/mL |
| | | | | | | | 2,4,5-Trichlorotoluene | 5000 ug/mL |
| | | | | | | | 2,4- & 2,5- & 2,6- Dichlorotoluene | 15000 ug/mL |
| | | | | | | | 2,4-Dichloro-1-(trifluoromethyl)-benzene | 5000 ug/mL |
| | | | | | | | 2,5-Dichlorobenzotrifluoride | 5000 ug/mL |
| | | | | | | | 2-Chlorobenzotrifluoride | 5000 ug/mL |
| | | | | | | | 3-Chlorobenzotrifluoride | 5000 ug/mL |
| | | | | | | | 3-Chlorotoluene | 5000 ug/mL |
| | | | | | | | 4-Chlorobenzotrifluoride | 5000 ug/mL |
| voaWKet2ndRes_00021 | 08/24/17 | 07/24/17 | Methanol, Lot 2019056 | 50 mL | VOA8260KET2ND_00098 | 100 uL | 2-Butanone (MEK) | 25 ug/mL |
| | | | | | | | 2-Hexanone | 25 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 25 ug/mL |
| | | | | | | | Acetone | 25 ug/mL |
| .VOA8260KET2ND_00098 | 03/31/19 | Restek, Lot A0123880 | | | (Purchased Reagent) | | 2-Butanone (MEK) | 12500 ug/mL |
| | | | | | | | 2-Hexanone | 12500 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 12500 ug/mL |
| | | | | | | | Acetone | 12500 ug/mL |
| voaWKetmix1st_00004 | 07/29/17 | 06/29/17 | Methanol, Lot 2019054 | 50 mL | VOA8260KET1ST_00099 | 0.1 mL | 2-Butanone (MEK) | 25 ug/mL |
| | | | | | | | 2-Hexanone | 25 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 25 ug/mL |
| | | | | | | | Acetone | 25 ug/mL |
| .VOA8260KET1ST_00099 | 01/31/20 | Restek, Lot A0123890 | | | (Purchased Reagent) | | 2-Butanone (MEK) | 12500 ug/mL |
| | | | | | | | 2-Hexanone | 12500 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 12500 ug/mL |
| | | | | | | | Acetone | 12500 ug/mL |
| voaWKetmix1st_00006 | 10/25/17 | 09/25/17 | Methanol, Lot 2469119 | 50 mL | VOA8260KET1ST_00102 | 100 uL | 2-Butanone (MEK) | 25 ug/mL |
| | | | | | | | 2-Hexanone | 25 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 25 ug/mL |
| | | | | | | | Acetone | 25 ug/mL |
| .VOA8260KET1ST_00102 | 01/31/20 | Restek, Lot A0123890 | | | (Purchased Reagent) | | 2-Butanone (MEK) | 12500 ug/mL |
| | | | | | | | 2-Hexanone | 12500 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 12500 ug/mL |
| | | | | | | | Acetone | 12500 ug/mL |
| voaWVA1stRest_00016 | 07/31/17 | 07/17/17 | Methanol, Lot 2019056 | 25 mL | VOA8260VARES_00082 | 0.125 mL | Vinyl acetate | 25 ug/mL |
| .VOA8260VARES_00082 | 07/31/17 | Restek, Lot A0124520 | | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |

Reagent

VOA8260GAS1ST_00203



CERTIFIED REFERENCE MATERIAL

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

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.: A0124278

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

8260 List 1 / Std #3 Gases (2015) 2,500 ug/ml, P&T Methanol, 1 ml/ampul

Container Size : 2 mL

Pkg Amt: > 1 mL

Expiration Date : January 31, 2020

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) CAS # 75-71-8 Purity 99% | 2,500.5 µg/mL (Lot Q167-08) | +/- 16.7232 +/- 140.4412 +/- 143.7161 | µg/mL | Gravimetric Unstressed Stressed |
| 2 | Chloromethane (methyl chloride) CAS # 74-87-3 Purity 99% | 2,498.7 µg/mL (Lot SHBG7976V) | +/- 17.4998 +/- 140.4406 +/- 143.7111 | µg/mL | Gravimetric Unstressed Stressed |
| 3 | Vinyl chloride CAS # 75-01-4 Purity 99% | 2,498.4 µg/mL (Lot 1026101231B1) | +/- 16.6753 +/- 140.3203 +/- 143.5926 | µg/mL | Gravimetric Unstressed Stressed |
| 4 | 1,3-Butadiene CAS # 106-99-0 Purity 99% | 2,496.9 µg/mL (Lot SHBF3387V) | +/- 17.0619 +/- 140.2843 +/- 143.5535 | µg/mL | Gravimetric Unstressed Stressed |
| 5 | Bromomethane (methyl bromide) CAS # 74-83-9 Purity 99% | 2,500.5 µg/mL (Lot 101604) | +/- 17.3456 +/- 140.5211 +/- 143.7944 | µg/mL | Gravimetric Unstressed Stressed |
| 6 | Chloroethane (ethyl chloride) CAS # 75-00-3 Purity 99% | 2,500.5 µg/mL (Lot 23593) | +/- 16.8189 +/- 140.4526 +/- 143.7272 | µg/mL | Gravimetric Unstressed Stressed |
| 7 | Dichlorofluoromethane (CPC-21) CAS # 75-43-4 Purity 99% | 2,500.0 µg/mL (Lot 4938100) | +/- 10.0499 +/- 139.7786 +/- 143.0675 | µg/mL | Gravimetric Unstressed Stressed |

| | | | | |
|---|---------------------------------|-----------------|--------------------|-------------|
| 8 | Trichlorofluoromethane (CFC-11) | 2,501.5 µg/mL | +/- 16.5404 µg/mL | Gravimetric |
| | CAS # 75-69-4 | (Lot SHBG7531V) | +/- 140.4793 µg/mL | Unstressed |
| | Purity 99% | | +/- 143.7562 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)

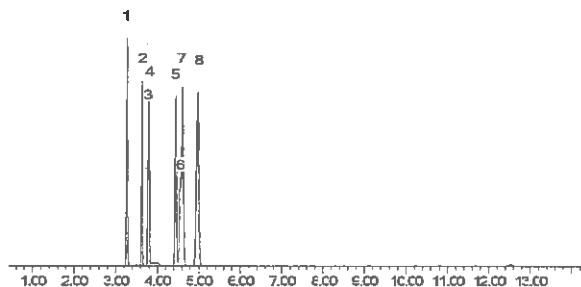
Carrier Gas:
helium-constant flow 2.0 mL/min.

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.

Joseph Jagiowski - Mix Technician

Date Mixed: 17-Jan-2017 Balance: 1125113331

Jennifer Pollino - Operations Tech-ARM QC

Date Passed: 24-Jan-2017

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

Reagent

VOA8260GAS1ST_00205



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

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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.: A0124278

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

8260 List 1 / Std #3 Gases (2015) 2,500 ug/ml, P&T Methanol, 1 ml/ampul

Container Size : 2 mL

Pkg Amt: > 1 mL

Expiration Date : January 31, 2020

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) CAS # 75-71-8 Purity 99% | 2,500.5 µg/mL (Lot Q167-08) | +/- 16.7232 +/- 140.4412 +/- 143.7161 | µg/mL | Gravimetric Unstressed Stressed |
| 2 | Chloromethane (methyl chloride) CAS # 74-87-3 Purity 99% | 2,498.7 µg/mL (Lot SHBG7976V) | +/- 17.4998 +/- 140.4406 +/- 143.7111 | µg/mL | Gravimetric Unstressed Stressed |
| 3 | Vinyl chloride CAS # 75-01-4 Purity 99% | 2,498.4 µg/mL (Lot 1026101231B1) | +/- 16.6753 +/- 140.3203 +/- 143.5926 | µg/mL | Gravimetric Unstressed Stressed |
| 4 | 1,3-Butadiene CAS # 106-99-0 Purity 99% | 2,496.9 µg/mL (Lot SHBF3387V) | +/- 17.0619 +/- 140.2843 +/- 143.5535 | µg/mL | Gravimetric Unstressed Stressed |
| 5 | Bromomethane (methyl bromide) CAS # 74-83-9 Purity 99% | 2,500.5 µg/mL (Lot 101604) | +/- 17.3456 +/- 140.5211 +/- 143.7944 | µg/mL | Gravimetric Unstressed Stressed |
| 6 | Chloroethane (ethyl chloride) CAS # 75-00-3 Purity 99% | 2,500.5 µg/mL (Lot 23593) | +/- 16.8189 +/- 140.4526 +/- 143.7272 | µg/mL | Gravimetric Unstressed Stressed |
| 7 | Dichlorofluoromethane (CPC-21) CAS # 75-43-4 Purity 99% | 2,500.0 µg/mL (Lot 4938100) | +/- 10.0499 +/- 139.7786 +/- 143.0675 | µg/mL | Gravimetric Unstressed Stressed |

| | | | | |
|--------|---------------------------------|-----------------|--------------------|-------------|
| 8 | Trichlorofluoromethane (CFC-11) | 2,501.5 µg/mL | +/- 16.5404 µg/mL | Gravimetric |
| CAS # | 75-69-4 | (Lot SHBG7531V) | +/- 140.4793 µg/mL | Unstressed |
| Purity | 99% | | +/- 143.7562 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)

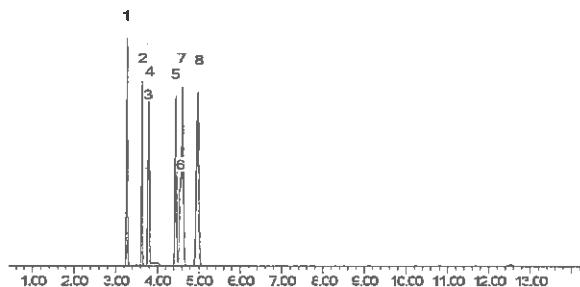
Carrier Gas:
helium-constant flow 2.0 mL/min.

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

Inj. Temp:
200°C

Det. Temp:
250°C

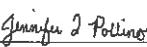
Det. Type:
MSD



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.


Joseph Jagiowski - Mix Technician

Date Mixed: 17-Jan-2017 Balance: 1125113331


Jennifer Pollino - Operations Tech-ARM QC

Date Passed: 24-Jan-2017

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

Reagent

VOA8260GAS2ND_00200



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Catalog No. : 569722.SEC

Lot No.: A0124116

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

8260 List 1 / Std #3 Gases (2015) 2,500 ug/ml, P&T Methanol, 1 ml/ampul

Container Size : 2 mL

Pkg Amt: > 1 mL

Expiration Date : January 31, 2020

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) CAS # 75-71-8.SEC Purity 99% | 2,503.4 µg/mL (Lot 23586) | +/- 19.5506 µg/mL | +/- 140.9699 µg/mL | +/- 144.2404 µg/mL |
| 2 | Chloromethane (methyl chloride) CAS # 74-87-3.SEC Purity 99% | 2,508.1 µg/mL (Lot 18343) | +/- 21.1963 µg/mL | +/- 141.4639 µg/mL | +/- 144.7353 µg/mL |
| 3 | Vinyl chloride CAS # 75-01-4.SEC Purity 99% | 2,518.6 µg/mL (Lot MKBK6872V) | +/- 19.4186 µg/mL | +/- 141.7924 µg/mL | +/- 145.0836 µg/mL |
| 4 | 1,3-Butadiene CAS # 106-99-0.SEC Purity 99% | 2,504.0 µg/mL (Lot 22331) | +/- 20.5722 µg/mL | +/- 141.1450 µg/mL | +/- 144.4130 µg/mL |
| 5 | Bromomethane (methyl bromide) CAS # 74-83-9.SEC Purity 99% | 2,498.5 µg/mL (Lot Q119-46) | +/- 19.9806 µg/mL | +/- 140.7602 µg/mL | +/- 144.0229 µg/mL |
| 6 | Chloroethane (ethyl chloride) CAS # 75-00-3.SEC Purity 99% | 2,494.9 µg/mL (Lot 00004202) | +/- 17.8868 µg/mL | +/- 140.2786 µg/mL | +/- 143.5429 µg/mL |
| 7 | Dichlorofluoromethane (CFC-21) CAS # 75-43-4.SEC Purity 99% | 2,503.4 µg/mL (Lot SHBC0858V) | +/- 20.0421 µg/mL | +/- 141.0350 µg/mL | +/- 144.3039 µg/mL |

| | | | | | | | |
|---|---------------------------------|---------------|-------|-----|----------|-------|-------------|
| 8 | Trichlorofluoromethane (CFC-11) | 2,503.2 | µg/mL | +/- | 18.7037 | µg/mL | Gravimetric |
| | CAS # 75-69-4 SEC | (Lot Q12B-59) | | +/- | 140.8450 | µg/mL | Unstressed |
| | Purity 99% | | | +/- | 144.1179 | 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)

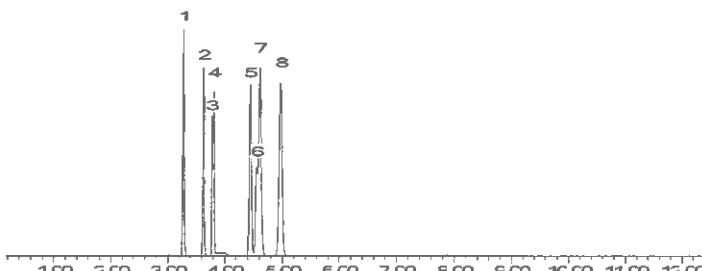
Carrier Gas:
helium-constant flow 2.0 mL/min.

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.

Brandon Reish
Brandon Reish - Mix Technician

Date Mixed: 12-Jan-2017 Balance: 1127510105

Jennifer J Pollino
Jennifer Pollino - Operations Tech-ARM QC

Date Passed: 17-Jan-2017

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

Reagent

VOA8260INTRES_00123



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| | | | |
|-------------------|---|----------|-----------------|
| Catalog No.: | <u>568718</u> | Lot No.: | <u>A0113246</u> |
| Description : | 8260 Internal Standard 2014 | | |
| | 8260 Internal Standard 2014 250-5,000 ug/ml, P&T Methanol, 5 ml/ampul | | |
| Container Size : | 5 mL | Pkg Amt: | > 5 mL |
| Expiration Date : | August 31, 2020 | Storage: | 0°C or colder |

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

| Edition Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.I., K=2) | | |
|---------------|---|--------------------------------|---|---|---------------------------------------|
| 1 | tert-Butyl-d9-alcohol CAS # 25725-11-5 Purity 99% | 5,000.4 µg/mL (Lot I201P18) | +/- 29.0712 µg/mL +/- 106.0450 µg/mL +/- 106.5155 µg/mL | +/- 29.0712 µg/mL +/- 106.0450 µg/mL +/- 106.5155 µg/mL | Gravimetric Unstressed Stressed |
| 2 | 2-Butanone-d5 CAS # 24313-50-6 Purity 99% | 1,250.2 µg/mL (Lot M276P24) | +/- 7.2688 µg/mL +/- 26.5135 µg/mL +/- 26.6311 µg/mL | +/- 7.2688 µg/mL +/- 26.5135 µg/mL +/- 26.6311 µg/mL | Gravimetric Unstressed Stressed |
| 3 | Fluorobenzene CAS # 462-06-6 Purity 99% | 250.2 µg/mL (Lot BCBK8171V) | +/- 1.4580 µg/mL +/- 5.3070 µg/mL +/- 5.3305 µg/mL | +/- 1.4580 µg/mL +/- 5.3070 µg/mL +/- 5.3305 µg/mL | Gravimetric Unstressed Stressed |
| 4 | 1,4-Dioxane-d8 CAS # 17647-74-4 Purity 98% | 5,000.6 µg/mL (Lot I-19073) | +/- 29.0727 µg/mL +/- 106.0502 µg/mL +/- 106.5208 µg/mL | +/- 29.0727 µg/mL +/- 106.0502 µg/mL +/- 106.5208 µg/mL | Gravimetric Unstressed Stressed |
| 5 | Chlorobenzene-d5 CAS # 3114-55-4 Purity 99% | 250.4 µg/mL (Lot PR-23926) | +/- 1.4592 µg/mL +/- 5.3113 µg/mL +/- 5.3348 µg/mL | +/- 1.4592 µg/mL +/- 5.3113 µg/mL +/- 5.3348 µg/mL | Gravimetric Unstressed Stressed |
| 6 | 1,4-Dichlorobenzene-d4 CAS # 3855-82-1 Purity 99% | 250.0 µg/mL (Lot PR-18488) | +/- 1.4569 µg/mL +/- 5.3028 µg/mL +/- 5.3263 µg/mL | +/- 1.4569 µg/mL +/- 5.3028 µg/mL +/- 5.3263 µg/mL | Gravimetric Unstressed Stressed |

Reagent

VOA8260INTRES_00135



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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.: 568718

Lot No.: A0124343

Description : 8260 Internal Standard 2014

8260 Internal Standard 2014 250-5,000 ug/ml, P&T Methanol, 5 ml/ampul

Container Size : 5 mL

Pkg Amt: > 5 mL

Expiration Date : January 31, 2022

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 CAS # 25725-11-5 Purity 99% | 5,050.0 µg/mL (Lot I-201) | +/- 29.3596 | µg/mL | Gravimetric |
| | | | +/- 108.1207 | µg/mL | Unstressed |
| | | | +/- 111.2640 | µg/mL | Stressed |
| 2 | 2-Butanone-d5 CAS # 24313-50-6 Purity 99% | 1,262.5 µg/mL (Lot M-276) | +/- 7.3403 | µg/mL | Gravimetric |
| | | | +/- 27.0303 | µg/mL | Unstressed |
| | | | +/- 27.8161 | µg/mL | Stressed |
| 3 | Fluorobenzene CAS # 462-06-6 Purity 99% | 251.6 µg/mL (Lot BCBK8171V) | +/- 1.4664 | µg/mL | Gravimetric |
| | | | +/- 5.3884 | µg/mL | Unstressed |
| | | | +/- 5.5450 | µg/mL | Stressed |
| 4 | 1,4-Dioxane-d8 CAS # 17647-74-4 Purity 99% | 5,048.8 µg/mL (Lot I-19942) | +/- 29.3526 | µg/mL | Gravimetric |
| | | | +/- 108.0950 | µg/mL | Unstressed |
| | | | +/- 111.2375 | µg/mL | Stressed |
| 5 | Chlorobenzene-d5 CAS # 3114-55-4 Purity 99% | 251.5 µg/mL (Lot PR-23926) | +/- 1.4654 | µg/mL | Gravimetric |
| | | | +/- 5.3849 | µg/mL | Unstressed |
| | | | +/- 5.5413 | µg/mL | Stressed |
| 6 | 1,4-Dichlorobenzene-d4 CAS # 3855-82-1 Purity 99% | 252.5 µg/mL (Lot PR-18488) | +/- 1.4714 | µg/mL | Gravimetric |
| | | | +/- 5.4070 | µg/mL | Unstressed |
| | | | +/- 5.5641 | µg/mL | Stressed |

Reagent

VOA8260KET1ST_00099



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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.: A0123890

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, 2020

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) | | | |
|---------------|---|--------------------------------|---|----------|-------|-------------|
| | | | +/- | 72.7778 | µg/mL | |
| 1 | Acetone CAS # 67-64-1 Purity 99% | (Lot SHBH0922V) | +/- | 72.7778 | µg/mL | Gravimetric |
| | | | +/- | 755.2362 | µg/mL | Unstressed |
| | | | +/- | 757.0293 | µg/mL | Stressed |
| 2 | 2-Butanone (MEK) CAS # 78-93-3 Purity 99% | (Lot SHBF2461V) | +/- | 72.8025 | µg/mL | Gravimetric |
| | | | +/- | 755.4927 | µg/mL | Unstressed |
| | | | +/- | 757.2863 | µg/mL | Stressed |
| 3 | 4-Methyl-2-pentanone (MIBK) CAS # 108-10-1 Purity 99% | (Lot SHBG3630V) | +/- | 72.7909 | µg/mL | Gravimetric |
| | | | +/- | 755.3720 | µg/mL | Unstressed |
| | | | +/- | 757.1654 | µg/mL | Stressed |
| 4 | 2-Hexanone CAS # 591-78-6 Purity 99% | (Lot MKBW0198V) | +/- | 72.7255 | µg/mL | Gravimetric |
| | | | +/- | 754.6932 | µg/mL | Unstressed |
| | | | +/- | 756.4850 | µg/mL | Stressed |
| Solvent: | P&T Methanol/Water (90:10) | | | | | |
| | CAS # 67-56-1/7732-18-5 | | | | | |
| | Purity 99% | | | | | |

Reagent

VOA8260KET1ST_00100



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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.: A0123890

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, 2020

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) | | | |
|---------------|---|--------------------------------|---|----------|-------|-------------|
| | | | +/- | 72.7778 | µg/mL | |
| 1 | Acetone CAS # 67-64-1 Purity 99% | (Lot SHBH0922V) | +/- | 72.7778 | µg/mL | Gravimetric |
| | | | +/- | 755.2362 | µg/mL | Unstressed |
| | | | +/- | 757.0293 | µg/mL | Stressed |
| 2 | 2-Butanone (MEK) CAS # 78-93-3 Purity 99% | (Lot SHBF2461V) | +/- | 72.8025 | µg/mL | Gravimetric |
| | | | +/- | 755.4927 | µg/mL | Unstressed |
| | | | +/- | 757.2863 | µg/mL | Stressed |
| 3 | 4-Methyl-2-pentanone (MIBK) CAS # 108-10-1 Purity 99% | (Lot SHBG3630V) | +/- | 72.7909 | µg/mL | Gravimetric |
| | | | +/- | 755.3720 | µg/mL | Unstressed |
| | | | +/- | 757.1654 | µg/mL | Stressed |
| 4 | 2-Hexanone CAS # 591-78-6 Purity 99% | (Lot MKBW0198V) | +/- | 72.7255 | µg/mL | Gravimetric |
| | | | +/- | 754.6932 | µg/mL | Unstressed |
| | | | +/- | 756.4850 | µg/mL | Stressed |
| Solvent: | P&T Methanol/Water (90:10) | | | | | |
| | CAS # 67-56-1/7732-18-5 | | | | | |
| | Purity 99% | | | | | |

Reagent

VOA8260KET1ST_00102



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Certificate of Analysis



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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.: A0123890

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, 2020

Storage: 0°C or colder

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

| Elation Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L.; K=2) | | |
|---------------|---|--------------------------------|---|-------|-------------|
| 1 | Acetone CAS # 67-64-1 Purity 99% | 12,517.5 µg/mL | +/- 72.7778 | µg/mL | Gravimetric |
| | | | +/- 755.2362 | µg/mL | Unstressed |
| | | | +/- 757.0293 | µg/mL | Stressed |
| 2 | 2-Butanone (MEK) CAS # 78-93-3 Purity 99% | 12,521.8 µg/mL | +/- 72.8025 | µg/mL | Gravimetric |
| | | | +/- 755.4927 | µg/mL | Unstressed |
| | | | +/- 757.2863 | µg/mL | Stressed |
| 3 | 4-Methyl-2-pentanone (MIBK) CAS # 108-10-1 Purity 99% | 12,519.8 µg/mL | +/- 72.7909 | µg/mL | Gravimetric |
| | | | +/- 755.3720 | µg/mL | Unstressed |
| | | | +/- 757.1654 | µg/mL | Stressed |
| 4 | 2-Hexanone CAS # 591-78-6 Purity 99% | 12,508.5 µg/mL | +/- 72.7255 | µg/mL | Gravimetric |
| | | | +/- 754.6932 | µg/mL | Unstressed |
| | | | +/- 756.4850 | µg/mL | Stressed |

Solvent: P&T Methanol/Water (90:10)

CAS # 67-56-1/7732-18-5

Purity 99%

Reagent

VOA8260KET2ND_00098

RESTEK CERTIFIED REFERENCE MATERIAL

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 Bellefonte, PA 16823-8812
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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>A0123880</u> |
| Description : | <u>8260 List 1/ Std #2 Ketones (2015)</u> | | |
| 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>> 1 mL</u> |
| Expiration Date : | <u>January 31, 2020</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 CAS # 67-64-1-SEC Purity 99% | 12,501.6 µg/mL | +/- 73.1996 µg/mL +/- 754.3267 µg/mL +/- 756.1173 µg/mL | µg/mL | Gravimetric Unstressed Stressed |
| 2 | 2-Butanone (MEK) CAS # 78-93-3-SEC Purity 99% | 12,503.6 µg/mL | +/- 73.2113 µg/mL +/- 754.4473 µg/mL +/- 756.2383 µg/mL | µg/mL | Gravimetric Unstressed Stressed |
| 3 | 4-Methyl-2-pentanone (MIBK) CAS # 108-10-1-SEC Purity 99% | 12,506.0 µg/mL | +/- 73.2254 µg/mL +/- 754.5921 µg/mL +/- 756.3834 µg/mL | µg/mL | Gravimetric Unstressed Stressed |
| 4 | 2-Hexanone CAS # 591-78-6-SEC Purity 99% | 12,504.0 µg/mL | +/- 73.2137 µg/mL +/- 754.4715 µg/mL +/- 756.2625 µg/mL | µg/mL | Gravimetric Unstressed Stressed |

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

Reagent

VOA8260MEGA1_00065


CERTIFIED REFERENCE MATERIAL

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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>571992</u> | Lot No.: | <u>A0123711</u> |
| Description : | 8260 List 1 / Std #1 MegaMix (2017) | | |
| | 8260 List 1 / Std #1 MegaMix (2017) 1250-62500 µg/ml, P&T Methanol, 1 ml/ampul | | |
| Container Size : | <u>2 mL</u> | Pkg Amt: | <u>> 1 mL</u> |
| Expiration Date : | <u>December 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 | Diethyl ether (ethyl ether) CAS # 60-29-7 Purity 99% | 2,501.3 µg/mL (Lot SHBG1462V) | +/- 14.5425 | µg/mL | Gravimetric |
| | | | +/- 150.9115 | µg/mL | Unstressed |
| | | | +/- 151.2698 | µg/mL | Stressed |
| 2 | 1,1,2-Trichlorotrifluoroethane (CFC-113) CAS # 76-13-1 Purity 99% | 2,505.1 µg/mL (Lot 00009482) | +/- 14.5650 | µg/mL | Gravimetric |
| | | | +/- 151.1453 | µg/mL | Unstressed |
| | | | +/- 151.5041 | µg/mL | Stressed |
| 3 | 1,1-dichloroethylene CAS # 75-35-4 Purity 99% | 2,511.5 µg/mL (Lot SHBG8609V) | +/- 14.6021 | µg/mL | Gravimetric |
| | | | +/- 151.5299 | µg/mL | Unstressed |
| | | | +/- 151.8897 | µg/mL | Stressed |
| 4 | tert-Butanol (TBA) CAS # 75-65-0 Purity 99% | 25,001.8 µg/mL (Lot SHBF0688V) | +/- 145.3547 | µg/mL | Gravimetric |
| | | | +/- 1,508.4656 | µg/mL | Unstressed |
| | | | +/- 1,512.0470 | µg/mL | Stressed |
| 5 | Methyl acetate CAS # 79-20-9 Purity 99% | 5,000.5 µg/mL (Lot SHBG4345V) | +/- 29.0733 | µg/mL | Gravimetric |
| | | | +/- 301.7023 | µg/mL | Unstressed |
| | | | +/- 302.4186 | µg/mL | Stressed |
| 6 | Iodomethane (methyl iodide) CAS # 74-88-4 Purity 99% | 2,502.9 µg/mL (Lot SHBF2149V) | +/- 14.5519 | µg/mL | Gravimetric |
| | | | +/- 151.0095 | µg/mL | Unstressed |
| | | | +/- 151.3681 | µg/mL | Stressed |
| 7 | Allyl chloride (3-chloropropene) CAS # 107-05-1 Purity 99% | 2,517.1 µg/mL (Lot SHBF8133V) | +/- 14.6348 | µg/mL | Gravimetric |
| | | | +/- 151.8693 | µg/mL | Unstressed |
| | | | +/- 152.2299 | µg/mL | Stressed |

| | | | | | | | |
|----|---|-----------------|----------|-------|--|-------------------------|---------------------------------------|
| 8 | Methylene chloride (dichloromethane) CAS # 75-09-2 Purity 99% | (Lot SHBH2578V) | 2,502.1 | µg/mL | +/- 14.5476 +/- 150.9643 +/- 151.3227 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 9 | Carbon disulfide CAS # 75-15-0 Purity 99% | (Lot S20A856) | 2,501.4 | µg/mL | +/- 14.5432 +/- 150.9190 +/- 151.2773 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 10 | Acrylonitrile CAS # 107-13-1 Purity 99% | (Lot T07B2030) | 25,001.3 | µg/mL | +/- 145.3518 +/- 1,508.4355 +/- 1,512.0167 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 11 | Methyl-tert-butyl ether (MTBE) CAS # 1634-04-4 Purity 99% | (Lot SHBG2655V) | 2,505.3 | µg/mL | +/- 14.5657 +/- 151.1528 +/- 151.5117 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 12 | cis-1,2-Dichloroethene CAS # 156-59-2 Purity 98% | (Lot MKBV2831V) | 2,500.5 | µg/mL | +/- 14.5379 +/- 150.8644 +/- 151.2226 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 13 | n-Hexane (C6) CAS # 110-54-3 Purity 99% | (Lot SHBG2674V) | 2,503.8 | µg/mL | +/- 14.5570 +/- 151.0623 +/- 151.4210 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 14 | 1,1-Dichloroethane CAS # 75-34-3 Purity 99% | (Lot 00008621) | 2,500.4 | µg/mL | +/- 14.5374 +/- 150.8587 +/- 151.2169 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 15 | 2,2-Dichloropropane CAS # 594-20-7 Purity 98% | (Lot BCBR0622V) | 2,501.0 | µg/mL | +/- 14.5408 +/- 150.8940 +/- 151.2522 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 16 | trans-1,2-Dichloroethene CAS # 156-60-5 Purity 99% | (Lot 09431AEV) | 2,503.8 | µg/mL | +/- 14.5570 +/- 151.0623 +/- 151.4210 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 17 | Isobutanol (2-Methyl-1-propanol) CAS # 78-83-1 Purity 99% | (Lot SHBG8201V) | 62,512.5 | µg/mL | +/- 363.4341 +/- 3,771.6543 +/- 3,780.6088 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 18 | chloroform CAS # 67-66-3 Purity 99% | (Lot MKBV2089V) | 2,501.9 | µg/mL | +/- 14.5461 +/- 150.9492 +/- 151.3076 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 19 | Bromochloromethane CAS # 74-97-5 Purity 99% | (Lot 00004559) | 2,503.3 | µg/mL | +/- 14.5541 +/- 151.0322 +/- 151.3907 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 20 | Tetrahydrofuran CAS # 109-99-9 Purity 99% | (Lot SHBG2910V) | 5,001.3 | µg/mL | +/- 29.0777 +/- 301.7476 +/- 302.4640 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 21 | 1,1,1-trichloroethane CAS # 71-55-6 Purity 99% | (Lot B15W12061) | 2,500.3 | µg/mL | +/- 14.5367 +/- 150.8512 +/- 151.2093 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 22 | Cyclohexane CAS # 110-82-7 Purity 99% | (Lot MKBX4768V) | 2,502.0 | µg/mL | +/- 14.5468 +/- 150.9567 +/- 151.3151 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 23 | 1,1-Dichloropropene CAS # 563-58-6 Purity 99% | (Lot 160727JLM) | 2,500.5 | µg/mL | +/- 14.5381 +/- 150.8662 +/- 151.2244 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |

| | | | | | | | |
|----|---|-----------------|----------|-------|--|-------------------------|---------------------------------------|
| 24 | carbon tetrachloride CAS # 56-23-5 Purity 99% | (Lot SHBG1763V) | 2,503.3 | µg/mL | +/- 14.5541 +/- 151.0322 +/- 151.3907 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 25 | n-Heptane (C7) CAS # 142-82-5 Purity 99% | (Lot SHBG6171V) | 2,505.5 | µg/mL | +/- 14.5672 +/- 151.1679 +/- 151.5268 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 26 | 1,2-Dichloroethane CAS # 107-06-2 Purity 99% | (Lot SHBF9313V) | 2,504.8 | µg/mL | +/- 14.5628 +/- 151.1227 +/- 151.4815 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 27 | Benzene CAS # 71-43-2 Purity 99% | (Lot SHBH2056V) | 2,506.9 | µg/mL | +/- 14.5752 +/- 151.2509 +/- 151.6100 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 28 | Trichloroethylene CAS # 79-01-6 Purity 99% | (Lot SHBH1955V) | 2,502.4 | µg/mL | +/- 14.5490 +/- 150.9794 +/- 151.3378 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 29 | Methylcyclohexane CAS # 108-87-2 Purity 98% | (Lot SHBG0634V) | 2,500.3 | µg/mL | +/- 14.5372 +/- 150.8570 +/- 151.2152 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 30 | 1,2-Dichloropropane CAS # 78-87-5 Purity 99% | (Lot 01113D0V) | 2,503.0 | µg/mL | +/- 14.5527 +/- 151.0171 +/- 151.3756 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 31 | 1,4-Dioxane CAS # 123-91-1 Purity 99% | (Lot SHBH2584V) | 50,011.4 | µg/mL | +/- 290.7552 +/- 3,017.4064 +/- 3,024.5702 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 32 | Dibromomethane CAS # 74-95-3 Purity 98% | (Lot 10183283) | 2,501.9 | µg/mL | +/- 14.5465 +/- 150.9531 +/- 151.3115 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 33 | cis-1,3-Dichloropropene CAS # 10061-01-5 Purity 99% | (Lot 22622) | 2,501.0 | µg/mL | +/- 14.5410 +/- 150.8964 +/- 151.2547 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 34 | Toluene CAS # 108-88-3 Purity 99% | (Lot SHBH1932V) | 2,504.3 | µg/mL | +/- 14.5599 +/- 151.0925 +/- 151.4512 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 35 | Ethyl methacrylate CAS # 97-63-2 Purity 99% | (Lot SHBD9190V) | 2,506.9 | µg/mL | +/- 14.5752 +/- 151.2509 +/- 151.6100 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 36 | trans-1,3-Dichloropropene CAS # 10061-02-6 Purity 99% | (Lot C584177) | 2,503.6 | µg/mL | +/- 14.5563 +/- 151.0548 +/- 151.4134 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 37 | 1,1,2-Trichloroethane CAS # 79-00-5 Purity 99% | (Lot FGB01) | 2,501.0 | µg/mL | +/- 14.5410 +/- 150.8964 +/- 151.2547 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 38 | 1,3-Dichloropropane CAS # 142-28-9 Purity 99% | (Lot BCBG2162V) | 2,503.5 | µg/mL | +/- 14.5556 +/- 151.0472 +/- 151.4059 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 39 | Tetrachloroethylene CAS # 127-18-4 Purity 99% | (Lot SHBD9374V) | 2,500.9 | µg/mL | +/- 14.5403 +/- 150.8889 +/- 151.2471 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |

| | | | | | | | | |
|----|---|-----------------|---------|-------|-----|----------|-------|-------------|
| 40 | dibromochloromethane CAS # 124-48-1 Purity 98% | (Lot MKBW3597V) | 2,500.2 | µg/mL | +/- | 14.5365 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8497 | µg/mL | Unstressed |
| | | | | | +/- | 151.2078 | µg/mL | Stressed |
| 41 | 1,2-Dibromoethane (EDB) CAS # 106-93-4 Purity 99% | (Lot BCBH3877V) | 2,501.3 | µg/mL | +/- | 14.5425 | µg/mL | Gravimetric |
| | | | | | +/- | 150.9115 | µg/mL | Unstressed |
| | | | | | +/- | 151.2698 | µg/mL | Stressed |
| 42 | Chlorobenzene CAS # 108-90-7 Purity 99% | (Lot SHBF0505V) | 2,500.1 | µg/mL | +/- | 14.5359 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8436 | µg/mL | Unstressed |
| | | | | | +/- | 151.2017 | µg/mL | Stressed |
| 43 | m-Xylene CAS # 108-38-3 Purity 99% | (Lot SHBG4347V) | 1,250.3 | µg/mL | +/- | 7.2691 | µg/mL | Gravimetric |
| | | | | | +/- | 75.4331 | µg/mL | Unstressed |
| | | | | | +/- | 75.6122 | µg/mL | Stressed |
| 44 | p-Xylene CAS # 106-42-3 Purity 99% | (Lot SHBG3928V) | 1,251.3 | µg/mL | +/- | 7.2749 | µg/mL | Gravimetric |
| | | | | | +/- | 75.4935 | µg/mL | Unstressed |
| | | | | | +/- | 75.6727 | µg/mL | Stressed |
| 45 | Ethylbenzene CAS # 100-41-4 Purity 99% | (Lot SHBG5920V) | 2,503.3 | µg/mL | +/- | 14.5541 | µg/mL | Gravimetric |
| | | | | | +/- | 151.0322 | µg/mL | Unstressed |
| | | | | | +/- | 151.3907 | µg/mL | Stressed |
| 46 | 1,1,1,2-Tetrachloroethane CAS # 630-20-6 Purity 99% | (Lot MKBS3769V) | 2,500.3 | µg/mL | +/- | 14.5367 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8512 | µg/mL | Unstressed |
| | | | | | +/- | 151.2093 | µg/mL | Stressed |
| 47 | o-Xylene CAS # 95-47-6 Purity 99% | (Lot SHBH3432V) | 2,504.9 | µg/mL | +/- | 14.5636 | µg/mL | Gravimetric |
| | | | | | +/- | 151.1302 | µg/mL | Unstressed |
| | | | | | +/- | 151.4890 | µg/mL | Stressed |
| 48 | Styrene CAS # 100-42-5 Purity 99% | (Lot MKBS7097V) | 2,506.3 | µg/mL | +/- | 14.5716 | µg/mL | Gravimetric |
| | | | | | +/- | 151.2132 | µg/mL | Unstressed |
| | | | | | +/- | 151.5722 | µg/mL | Stressed |
| 49 | Isopropylbenzene (cumene) CAS # 98-82-8 Purity 99% | (Lot 10185056) | 2,501.6 | µg/mL | +/- | 14.5447 | µg/mL | Gravimetric |
| | | | | | +/- | 150.9341 | µg/mL | Unstressed |
| | | | | | +/- | 151.2925 | µg/mL | Stressed |
| 50 | bromoform CAS # 75-25-2 Purity 99% | (Lot SHBD8459V) | 2,502.9 | µg/mL | +/- | 14.5519 | µg/mL | Gravimetric |
| | | | | | +/- | 151.0095 | µg/mL | Unstressed |
| | | | | | +/- | 151.3681 | µg/mL | Stressed |
| 51 | bromodichloromethane CAS # 75-27-4 Purity 97% | (Lot MKBW5506V) | 2,506.8 | µg/mL | +/- | 14.5750 | µg/mL | Gravimetric |
| | | | | | +/- | 151.2490 | µg/mL | Unstressed |
| | | | | | +/- | 151.6081 | µg/mL | Stressed |
| 52 | 1,1,2,2-Tetrachloroethane CAS # 79-34-5 Purity 99% | (Lot CFA4D) | 2,501.3 | µg/mL | +/- | 14.5425 | µg/mL | Gravimetric |
| | | | | | +/- | 150.9115 | µg/mL | Unstressed |
| | | | | | +/- | 151.2698 | µg/mL | Stressed |
| 53 | 1,2,3-Trichloropropane CAS # 96-18-4 Purity 99% | (Lot BCBH8722V) | 2,508.5 | µg/mL | +/- | 14.5846 | µg/mL | Gravimetric |
| | | | | | +/- | 151.3489 | µg/mL | Unstressed |
| | | | | | +/- | 151.7082 | µg/mL | Stressed |
| 54 | trans-1,4-dichloro-2-butene CAS # 110-57-6 Purity 95% | (Lot MKBP6041V) | 2,500.8 | µg/mL | +/- | 14.5396 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8817 | µg/mL | Unstressed |
| | | | | | +/- | 151.2399 | µg/mL | Stressed |
| 55 | n-Propylbenzene CAS # 103-65-1 Purity 99% | (Lot MKBJ0332V) | 2,501.9 | µg/mL | +/- | 14.5461 | µg/mL | Gravimetric |
| | | | | | +/- | 150.9492 | µg/mL | Unstressed |
| | | | | | +/- | 151.3076 | µg/mL | Stressed |

| | | | | | | | |
|----|--|-----------------|---------|-------|---|-------------------------|---------------------------------------|
| 56 | Bromobenzene CAS # 108-86-1 Purity 99% | (Lot MKBD4032V) | 2,507.0 | µg/mL | +/- 14.5759 +/- 151.2584 +/- 151.6175 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 57 | 1,3,5-Trimethylbenzene CAS # 108-67-8 Purity 99% | (Lot BCBQ2165V) | 2,501.1 | µg/mL | +/- 14.5418 +/- 150.9040 +/- 151.2622 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 58 | 2-Chlorotoluene CAS # 95-49-8 Purity 99% | (Lot MKBW5554V) | 2,500.6 | µg/mL | +/- 14.5388 +/- 150.8738 +/- 151.2320 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 59 | 4-Chlorotoluene CAS # 106-43-4 Purity 99% | (Lot MKBL7753V) | 2,501.3 | µg/mL | +/- 14.5425 +/- 150.9115 +/- 151.2698 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 60 | tert-Butylbenzene CAS # 98-06-6 Purity 99% | (Lot S52237V) | 2,507.0 | µg/mL | +/- 14.5759 +/- 151.2584 +/- 151.6175 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 61 | 1,2,4-Trimethylbenzene CAS # 95-63-6 Purity 98% | (Lot MKBJ6229V) | 2,500.8 | µg/mL | +/- 14.5401 +/- 150.8866 +/- 151.2448 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 62 | sec-Butylbenzene CAS # 135-98-8 Purity 99% | (Lot MKBR9260V) | 2,505.4 | µg/mL | +/- 14.5665 +/- 151.1604 +/- 151.5193 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 63 | p-Isopropyltoluene (p-Cymene) CAS # 99-87-6 Purity 99% | (Lot MKBS2604V) | 2,503.8 | µg/mL | +/- 14.5570 +/- 151.0623 +/- 151.4210 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 64 | 1,3-Dichlorobenzene CAS # 541-73-1 Purity 99% | (Lot BCBM5751V) | 2,503.9 | µg/mL | +/- 14.5577 +/- 151.0699 +/- 151.4285 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 65 | 1,4-Dichlorobenzene CAS # 106-46-7 Purity 99% | (Lot MKBS1350V) | 2,509.9 | µg/mL | +/- 14.5926 +/- 151.4319 +/- 151.7914 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 66 | n-Butylbenzene CAS # 104-51-8 Purity 99% | (Lot 09418JJV) | 2,503.3 | µg/mL | +/- 14.5541 +/- 151.0322 +/- 151.3907 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 67 | 1,2-Dichlorobenzene CAS # 95-50-1 Purity 99% | (Lot SHBD7331V) | 2,503.8 | µg/mL | +/- 14.5570 +/- 151.0623 +/- 151.4210 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 68 | 1,2-Dibromo-3-chloropropane CAS # 96-12-8 Purity 99% | (Lot FBL01) | 2,505.0 | µg/mL | +/- 14.5643 +/- 151.1378 +/- 151.4966 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 69 | 1,2,4-Trichlorobenzene CAS # 120-82-1 Purity 99% | (Lot SHBC5541V) | 2,505.3 | µg/mL | +/- 14.5657 +/- 151.1528 +/- 151.5117 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 70 | Hexachlorobutadiene CAS # 87-68-3 Purity 98% | (Lot J31X013) | 2,506.5 | µg/mL | +/- 14.5728 +/- 151.2266 +/- 151.5856 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 71 | Naphthalene CAS # 91-20-3 Purity 99% | (Lot MKBW2603V) | 2,500.9 | µg/mL | +/- 14.5403 +/- 150.8889 +/- 151.2471 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |

| | | | | | |
|----|---|----------------|---------------|---|---------------------------------------|
| 72 | 1,2,3-Trichlorobenzene CAS # 87-61-6 Purity 99% | (Lot 12912PFV) | 2,511.1 µg/mL | +/- 14.5999 µg/mL +/- 151.5073 µg/mL +/- 151.8670 µg/mL | Gravimetric Unstressed Stressed |
|----|---|----------------|---------------|---|---------------------------------------|

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

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

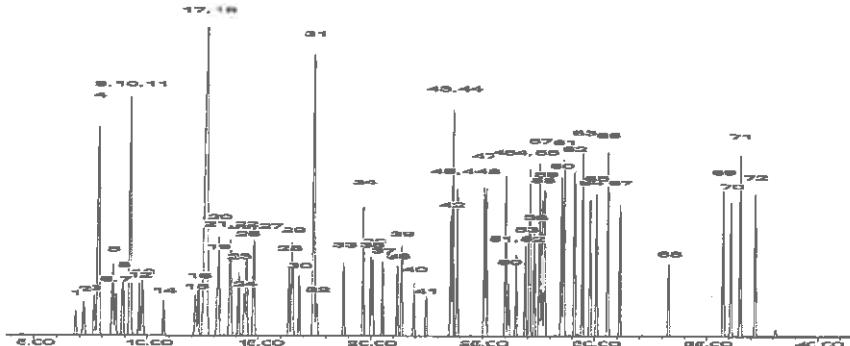
Carrier Gas:
helium-constant pressure 30 psi

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



Reagent

VOA8260MEGA1_00066



CERTIFIED REFERENCE MATERIAL

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

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.: 571992

Lot No.: A0123711

Description : 8260 List 1 / Std #1 MegaMix (2017)

8260 List 1 / Std #1 MegaMix (2017) 1250-62500 µg/ml, P&T Methanol, 1 ml/ampul

Container Size : 2 mL

Pkg Amt: > 1 mL

Expiration Date : December 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 | Diethyl ether (ethyl ether) CAS # 60-29-7 Purity 99% | 2,501.3 µg/mL (Lot SHBG1462V) | +/- 14.5425 | µg/mL | Gravimetric |
| | | | +/- 150.9115 | µg/mL | Unstressed |
| | | | +/- 151.2698 | µg/mL | Stressed |
| 2 | 1,1,2-Trichlorotrifluoroethane (CFC-113) CAS # 76-13-1 Purity 99% | 2,505.1 µg/mL (Lot 00009482) | +/- 14.5650 | µg/mL | Gravimetric |
| | | | +/- 151.1453 | µg/mL | Unstressed |
| | | | +/- 151.5041 | µg/mL | Stressed |
| 3 | 1,1-dichloroethylene CAS # 75-35-4 Purity 99% | 2,511.5 µg/mL (Lot SHBG8609V) | +/- 14.6021 | µg/mL | Gravimetric |
| | | | +/- 151.5299 | µg/mL | Unstressed |
| | | | +/- 151.8897 | µg/mL | Stressed |
| 4 | tert-Butanol (TBA) CAS # 75-65-0 Purity 99% | 25,001.8 µg/mL (Lot SHBF0688V) | +/- 145.3547 | µg/mL | Gravimetric |
| | | | +/- 1,508.4656 | µg/mL | Unstressed |
| | | | +/- 1,512.0470 | µg/mL | Stressed |
| 5 | Methyl acetate CAS # 79-20-9 Purity 99% | 5,000.5 µg/mL (Lot SHBG4345V) | +/- 29.0733 | µg/mL | Gravimetric |
| | | | +/- 301.7023 | µg/mL | Unstressed |
| | | | +/- 302.4186 | µg/mL | Stressed |
| 6 | Iodomethane (methyl iodide) CAS # 74-88-4 Purity 99% | 2,502.9 µg/mL (Lot SHBF2149V) | +/- 14.5519 | µg/mL | Gravimetric |
| | | | +/- 151.0095 | µg/mL | Unstressed |
| | | | +/- 151.3681 | µg/mL | Stressed |
| 7 | Allyl chloride (3-chloropropene) CAS # 107-05-1 Purity 99% | 2,517.1 µg/mL (Lot SHBF8133V) | +/- 14.6348 | µg/mL | Gravimetric |
| | | | +/- 151.8693 | µg/mL | Unstressed |
| | | | +/- 152.2299 | µg/mL | Stressed |

| | | | | | | | |
|----|---|-----------------|----------|-------|--|-------------------------|---------------------------------------|
| 8 | Methylene chloride (dichloromethane) CAS # 75-09-2 Purity 99% | (Lot SHBH2578V) | 2,502.1 | µg/mL | +/- 14.5476 +/- 150.9643 +/- 151.3227 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 9 | Carbon disulfide CAS # 75-15-0 Purity 99% | (Lot S20A856) | 2,501.4 | µg/mL | +/- 14.5432 +/- 150.9190 +/- 151.2773 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 10 | Acrylonitrile CAS # 107-13-1 Purity 99% | (Lot T07B2030) | 25,001.3 | µg/mL | +/- 145.3518 +/- 1,508.4355 +/- 1,512.0167 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 11 | Methyl-tert-butyl ether (MTBE) CAS # 1634-04-4 Purity 99% | (Lot SHBG2655V) | 2,505.3 | µg/mL | +/- 14.5657 +/- 151.1528 +/- 151.5117 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 12 | cis-1,2-Dichloroethene CAS # 156-59-2 Purity 98% | (Lot MKBV2831V) | 2,500.5 | µg/mL | +/- 14.5379 +/- 150.8644 +/- 151.2226 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 13 | n-Hexane (C6) CAS # 110-54-3 Purity 99% | (Lot SHBG2674V) | 2,503.8 | µg/mL | +/- 14.5570 +/- 151.0623 +/- 151.4210 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 14 | 1,1-Dichloroethane CAS # 75-34-3 Purity 99% | (Lot 00008621) | 2,500.4 | µg/mL | +/- 14.5374 +/- 150.8587 +/- 151.2169 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 15 | 2,2-Dichloropropane CAS # 594-20-7 Purity 98% | (Lot BCBR0622V) | 2,501.0 | µg/mL | +/- 14.5408 +/- 150.8940 +/- 151.2522 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 16 | trans-1,2-Dichloroethene CAS # 156-60-5 Purity 99% | (Lot 09431AEV) | 2,503.8 | µg/mL | +/- 14.5570 +/- 151.0623 +/- 151.4210 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 17 | Isobutanol (2-Methyl-1-propanol) CAS # 78-83-1 Purity 99% | (Lot SHBG8201V) | 62,512.5 | µg/mL | +/- 363.4341 +/- 3,771.6543 +/- 3,780.6088 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 18 | chloroform CAS # 67-66-3 Purity 99% | (Lot MKBV2089V) | 2,501.9 | µg/mL | +/- 14.5461 +/- 150.9492 +/- 151.3076 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 19 | Bromochloromethane CAS # 74-97-5 Purity 99% | (Lot 00004559) | 2,503.3 | µg/mL | +/- 14.5541 +/- 151.0322 +/- 151.3907 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 20 | Tetrahydrofuran CAS # 109-99-9 Purity 99% | (Lot SHBG2910V) | 5,001.3 | µg/mL | +/- 29.0777 +/- 301.7476 +/- 302.4640 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 21 | 1,1,1-trichloroethane CAS # 71-55-6 Purity 99% | (Lot B15W12061) | 2,500.3 | µg/mL | +/- 14.5367 +/- 150.8512 +/- 151.2093 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 22 | Cyclohexane CAS # 110-82-7 Purity 99% | (Lot MKBX4768V) | 2,502.0 | µg/mL | +/- 14.5468 +/- 150.9567 +/- 151.3151 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 23 | 1,1-Dichloropropene CAS # 563-58-6 Purity 99% | (Lot 160727JLM) | 2,500.5 | µg/mL | +/- 14.5381 +/- 150.8662 +/- 151.2244 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |

| | | | | | | | |
|----|---|-----------------|----------|-------|--|-------------------------|---------------------------------------|
| 24 | carbon tetrachloride CAS # 56-23-5 Purity 99% | (Lot SHBG1763V) | 2,503.3 | µg/mL | +/- 14.5541 +/- 151.0322 +/- 151.3907 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 25 | n-Heptane (C7) CAS # 142-82-5 Purity 99% | (Lot SHBG6171V) | 2,505.5 | µg/mL | +/- 14.5672 +/- 151.1679 +/- 151.5268 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 26 | 1,2-Dichloroethane CAS # 107-06-2 Purity 99% | (Lot SHBF9313V) | 2,504.8 | µg/mL | +/- 14.5628 +/- 151.1227 +/- 151.4815 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 27 | Benzene CAS # 71-43-2 Purity 99% | (Lot SHBH2056V) | 2,506.9 | µg/mL | +/- 14.5752 +/- 151.2509 +/- 151.6100 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 28 | Trichloroethylene CAS # 79-01-6 Purity 99% | (Lot SHBH1955V) | 2,502.4 | µg/mL | +/- 14.5490 +/- 150.9794 +/- 151.3378 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 29 | Methylcyclohexane CAS # 108-87-2 Purity 98% | (Lot SHBG0634V) | 2,500.3 | µg/mL | +/- 14.5372 +/- 150.8570 +/- 151.2152 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 30 | 1,2-Dichloropropane CAS # 78-87-5 Purity 99% | (Lot 01113D0V) | 2,503.0 | µg/mL | +/- 14.5527 +/- 151.0171 +/- 151.3756 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 31 | 1,4-Dioxane CAS # 123-91-1 Purity 99% | (Lot SHBH2584V) | 50,011.4 | µg/mL | +/- 290.7552 +/- 3,017.4064 +/- 3,024.5702 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 32 | Dibromomethane CAS # 74-95-3 Purity 98% | (Lot 10183283) | 2,501.9 | µg/mL | +/- 14.5465 +/- 150.9531 +/- 151.3115 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 33 | cis-1,3-Dichloropropene CAS # 10061-01-5 Purity 99% | (Lot 22622) | 2,501.0 | µg/mL | +/- 14.5410 +/- 150.8964 +/- 151.2547 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 34 | Toluene CAS # 108-88-3 Purity 99% | (Lot SHBH1932V) | 2,504.3 | µg/mL | +/- 14.5599 +/- 151.0925 +/- 151.4512 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 35 | Ethyl methacrylate CAS # 97-63-2 Purity 99% | (Lot SHBD9190V) | 2,506.9 | µg/mL | +/- 14.5752 +/- 151.2509 +/- 151.6100 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 36 | trans-1,3-Dichloropropene CAS # 10061-02-6 Purity 99% | (Lot C584177) | 2,503.6 | µg/mL | +/- 14.5563 +/- 151.0548 +/- 151.4134 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 37 | 1,1,2-Trichloroethane CAS # 79-00-5 Purity 99% | (Lot FGB01) | 2,501.0 | µg/mL | +/- 14.5410 +/- 150.8964 +/- 151.2547 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 38 | 1,3-Dichloropropane CAS # 142-28-9 Purity 99% | (Lot BCBG2162V) | 2,503.5 | µg/mL | +/- 14.5556 +/- 151.0472 +/- 151.4059 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 39 | Tetrachloroethylene CAS # 127-18-4 Purity 99% | (Lot SHBD9374V) | 2,500.9 | µg/mL | +/- 14.5403 +/- 150.8889 +/- 151.2471 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |

| | | | | | | | | |
|----|---|-----------------|---------|-------|-----|----------|-------|-------------|
| 40 | dibromochloromethane CAS # 124-48-1 Purity 98% | (Lot MKBW3597V) | 2,500.2 | µg/mL | +/- | 14.5365 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8497 | µg/mL | Unstressed |
| | | | | | +/- | 151.2078 | µg/mL | Stressed |
| 41 | 1,2-Dibromoethane (EDB) CAS # 106-93-4 Purity 99% | (Lot BCBH3877V) | 2,501.3 | µg/mL | +/- | 14.5425 | µg/mL | Gravimetric |
| | | | | | +/- | 150.9115 | µg/mL | Unstressed |
| | | | | | +/- | 151.2698 | µg/mL | Stressed |
| 42 | Chlorobenzene CAS # 108-90-7 Purity 99% | (Lot SHBF0505V) | 2,500.1 | µg/mL | +/- | 14.5359 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8436 | µg/mL | Unstressed |
| | | | | | +/- | 151.2017 | µg/mL | Stressed |
| 43 | m-Xylene CAS # 108-38-3 Purity 99% | (Lot SHBG4347V) | 1,250.3 | µg/mL | +/- | 7.2691 | µg/mL | Gravimetric |
| | | | | | +/- | 75.4331 | µg/mL | Unstressed |
| | | | | | +/- | 75.6122 | µg/mL | Stressed |
| 44 | p-Xylene CAS # 106-42-3 Purity 99% | (Lot SHBG3928V) | 1,251.3 | µg/mL | +/- | 7.2749 | µg/mL | Gravimetric |
| | | | | | +/- | 75.4935 | µg/mL | Unstressed |
| | | | | | +/- | 75.6727 | µg/mL | Stressed |
| 45 | Ethylbenzene CAS # 100-41-4 Purity 99% | (Lot SHBG5920V) | 2,503.3 | µg/mL | +/- | 14.5541 | µg/mL | Gravimetric |
| | | | | | +/- | 151.0322 | µg/mL | Unstressed |
| | | | | | +/- | 151.3907 | µg/mL | Stressed |
| 46 | 1,1,1,2-Tetrachloroethane CAS # 630-20-6 Purity 99% | (Lot MKBS3769V) | 2,500.3 | µg/mL | +/- | 14.5367 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8512 | µg/mL | Unstressed |
| | | | | | +/- | 151.2093 | µg/mL | Stressed |
| 47 | o-Xylene CAS # 95-47-6 Purity 99% | (Lot SHBH3432V) | 2,504.9 | µg/mL | +/- | 14.5636 | µg/mL | Gravimetric |
| | | | | | +/- | 151.1302 | µg/mL | Unstressed |
| | | | | | +/- | 151.4890 | µg/mL | Stressed |
| 48 | Styrene CAS # 100-42-5 Purity 99% | (Lot MKBS7097V) | 2,506.3 | µg/mL | +/- | 14.5716 | µg/mL | Gravimetric |
| | | | | | +/- | 151.2132 | µg/mL | Unstressed |
| | | | | | +/- | 151.5722 | µg/mL | Stressed |
| 49 | Isopropylbenzene (cumene) CAS # 98-82-8 Purity 99% | (Lot 10185056) | 2,501.6 | µg/mL | +/- | 14.5447 | µg/mL | Gravimetric |
| | | | | | +/- | 150.9341 | µg/mL | Unstressed |
| | | | | | +/- | 151.2925 | µg/mL | Stressed |
| 50 | bromoform CAS # 75-25-2 Purity 99% | (Lot SHBD8459V) | 2,502.9 | µg/mL | +/- | 14.5519 | µg/mL | Gravimetric |
| | | | | | +/- | 151.0095 | µg/mL | Unstressed |
| | | | | | +/- | 151.3681 | µg/mL | Stressed |
| 51 | bromodichloromethane CAS # 75-27-4 Purity 97% | (Lot MKBW5506V) | 2,506.8 | µg/mL | +/- | 14.5750 | µg/mL | Gravimetric |
| | | | | | +/- | 151.2490 | µg/mL | Unstressed |
| | | | | | +/- | 151.6081 | µg/mL | Stressed |
| 52 | 1,1,2,2-Tetrachloroethane CAS # 79-34-5 Purity 99% | (Lot CFA4D) | 2,501.3 | µg/mL | +/- | 14.5425 | µg/mL | Gravimetric |
| | | | | | +/- | 150.9115 | µg/mL | Unstressed |
| | | | | | +/- | 151.2698 | µg/mL | Stressed |
| 53 | 1,2,3-Trichloropropane CAS # 96-18-4 Purity 99% | (Lot BCBH8722V) | 2,508.5 | µg/mL | +/- | 14.5846 | µg/mL | Gravimetric |
| | | | | | +/- | 151.3489 | µg/mL | Unstressed |
| | | | | | +/- | 151.7082 | µg/mL | Stressed |
| 54 | trans-1,4-dichloro-2-butene CAS # 110-57-6 Purity 95% | (Lot MKBP6041V) | 2,500.8 | µg/mL | +/- | 14.5396 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8817 | µg/mL | Unstressed |
| | | | | | +/- | 151.2399 | µg/mL | Stressed |
| 55 | n-Propylbenzene CAS # 103-65-1 Purity 99% | (Lot MKBJ0332V) | 2,501.9 | µg/mL | +/- | 14.5461 | µg/mL | Gravimetric |
| | | | | | +/- | 150.9492 | µg/mL | Unstressed |
| | | | | | +/- | 151.3076 | µg/mL | Stressed |

| | | | | | | | |
|----|--|-----------------|---------|-------|---|-------------------------|---------------------------------------|
| 56 | Bromobenzene CAS # 108-86-1 Purity 99% | (Lot MKBD4032V) | 2,507.0 | µg/mL | +/- 14.5759 +/- 151.2584 +/- 151.6175 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 57 | 1,3,5-Trimethylbenzene CAS # 108-67-8 Purity 99% | (Lot BCBQ2165V) | 2,501.1 | µg/mL | +/- 14.5418 +/- 150.9040 +/- 151.2622 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 58 | 2-Chlorotoluene CAS # 95-49-8 Purity 99% | (Lot MKBW5554V) | 2,500.6 | µg/mL | +/- 14.5388 +/- 150.8738 +/- 151.2320 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 59 | 4-Chlorotoluene CAS # 106-43-4 Purity 99% | (Lot MKBL7753V) | 2,501.3 | µg/mL | +/- 14.5425 +/- 150.9115 +/- 151.2698 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 60 | tert-Butylbenzene CAS # 98-06-6 Purity 99% | (Lot S52237V) | 2,507.0 | µg/mL | +/- 14.5759 +/- 151.2584 +/- 151.6175 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 61 | 1,2,4-Trimethylbenzene CAS # 95-63-6 Purity 98% | (Lot MKBJ6229V) | 2,500.8 | µg/mL | +/- 14.5401 +/- 150.8866 +/- 151.2448 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 62 | sec-Butylbenzene CAS # 135-98-8 Purity 99% | (Lot MKBR9260V) | 2,505.4 | µg/mL | +/- 14.5665 +/- 151.1604 +/- 151.5193 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 63 | p-Isopropyltoluene (p-Cymene) CAS # 99-87-6 Purity 99% | (Lot MKBS2604V) | 2,503.8 | µg/mL | +/- 14.5570 +/- 151.0623 +/- 151.4210 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 64 | 1,3-Dichlorobenzene CAS # 541-73-1 Purity 99% | (Lot BCBM5751V) | 2,503.9 | µg/mL | +/- 14.5577 +/- 151.0699 +/- 151.4285 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 65 | 1,4-Dichlorobenzene CAS # 106-46-7 Purity 99% | (Lot MKBS1350V) | 2,509.9 | µg/mL | +/- 14.5926 +/- 151.4319 +/- 151.7914 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 66 | n-Butylbenzene CAS # 104-51-8 Purity 99% | (Lot 09418JJV) | 2,503.3 | µg/mL | +/- 14.5541 +/- 151.0322 +/- 151.3907 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 67 | 1,2-Dichlorobenzene CAS # 95-50-1 Purity 99% | (Lot SHBD7331V) | 2,503.8 | µg/mL | +/- 14.5570 +/- 151.0623 +/- 151.4210 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 68 | 1,2-Dibromo-3-chloropropane CAS # 96-12-8 Purity 99% | (Lot FBL01) | 2,505.0 | µg/mL | +/- 14.5643 +/- 151.1378 +/- 151.4966 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 69 | 1,2,4-Trichlorobenzene CAS # 120-82-1 Purity 99% | (Lot SHBC5541V) | 2,505.3 | µg/mL | +/- 14.5657 +/- 151.1528 +/- 151.5117 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 70 | Hexachlorobutadiene CAS # 87-68-3 Purity 98% | (Lot J31X013) | 2,506.5 | µg/mL | +/- 14.5728 +/- 151.2266 +/- 151.5856 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 71 | Naphthalene CAS # 91-20-3 Purity 99% | (Lot MKBW2603V) | 2,500.9 | µg/mL | +/- 14.5403 +/- 150.8889 +/- 151.2471 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |

72 1,2,3-Trichlorobenzene 2,511.1 µg/mL +/- 14.5999 µg/mL Gravimetric
 CAS # 87-61-6 (Lot 12912PFV) +/- 151.5073 µg/mL Unstressed
 Purity 99% +/- 151.8670 µg/mL Stressed

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

Column 3

60m x 0.25mm 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.)

Ini. Temp:

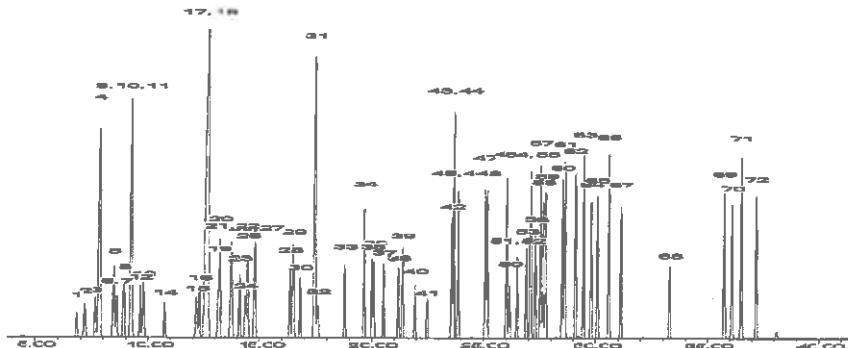
300°c

Det. Temp:

DCL
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.

E. Joseph Tally - Mix Technician

Date Mixed: 22-Dec-2016 Balance: B251644995

Jennifer Pollino - Operations Tech-ARM QC

Date Passed: 04-Jan-2017

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

Reagent

VOA8260MEGA2_00061



CERTIFIED REFERENCE MATERIAL

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

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. : | <u>571992.sec</u> | Lot No.: <u>A0123775</u> | | | |
| Description : | 8260 List 1 / Std #1 MegaMix (2017) | | | | |
| | 8260 List 1 / Std #1 MegaMix (2017) 1250-62500 µg/ml, P&T Methanol, 1 ml/ampul | | | | |
| Container Size : | <u>2 mL</u> | Pkg Amt: | <u>> 1 mL</u> | | |
| Expiration Date : | <u>December 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 | Diethyl ether (ethyl ether) CAS # 60-29-7.SEC Purity 98% | 2,501.2 µg/mL (Lot F23X068) | +/- 14.5422 | µg/mL | Gravimetric |
| | | | +/- 150.9088 | µg/mL | Unstressed |
| | | | +/- 151.2671 | µg/mL | Stressed |
| 2 | 1,1,2-Trichlorotrifluoroethane (CFC-113) CAS # 76-13-1.SEC Purity 99% | 2,501.1 µg/mL (Lot 18342) | +/- 14.5418 | µg/mL | Gravimetric |
| | | | +/- 150.9040 | µg/mL | Unstressed |
| | | | +/- 151.2622 | µg/mL | Stressed |
| 3 | 1,1-Dichloroethene CAS # 75-35-4.SEC Purity 99% | 2,500.5 µg/mL (Lot 2767000) | +/- 14.5381 | µg/mL | Gravimetric |
| | | | +/- 150.8662 | µg/mL | Unstressed |
| | | | +/- 151.2244 | µg/mL | Stressed |
| 4 | tert-Butanol (TBA) CAS # 75-65-0.SEC Purity 98% | 25,003.1 µg/mL (Lot XYXDO) | +/- 145.3626 | µg/mL | Gravimetric |
| | | | +/- 1,508.5475 | µg/mL | Unstressed |
| | | | +/- 1,512.1291 | µg/mL | Stressed |
| 5 | Methyl acetate CAS # 79-20-9.SEC Purity 99% | 5,000.4 µg/mL (Lot YDGVD) | +/- 29.0726 | µg/mL | Gravimetric |
| | | | +/- 301.6948 | µg/mL | Unstressed |
| | | | +/- 302.4111 | µg/mL | Stressed |
| 6 | Iodomethane (methyl iodide) CAS # 74-88-4.SEC Purity 99% | 2,500.4 µg/mL (Lot Y25A027) | +/- 14.5374 | µg/mL | Gravimetric |
| | | | +/- 150.8587 | µg/mL | Unstressed |
| | | | +/- 151.2169 | µg/mL | Stressed |
| 7 | Allyl chloride (3-chloropropene) CAS # 107-05-1.SEC Purity 98% | 2,500.1 µg/mL (Lot VEBOC) | +/- 14.5358 | µg/mL | Gravimetric |
| | | | +/- 150.8423 | µg/mL | Unstressed |
| | | | +/- 151.2004 | µg/mL | Stressed |

| | | | | | | |
|----|---|------------------|----------------|--|-------------------------|---------------------------------------|
| 8 | Methylene chloride (dichloromethane) CAS # 75-09-2-SEC Purity 99% | (Lot FGM02) | 2,500.8 µg/mL | +/- 14.5396 +/- 150.8813 +/- 151.2395 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 9 | Carbon disulfide CAS # 75-15-0-SEC Purity 99% | (Lot MKBL1376V) | 2,500.9 µg/mL | +/- 14.5403 +/- 150.8889 +/- 151.2471 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 10 | Acrylonitrile CAS # 107-13-1-SEC Purity 99% | (Lot UERIL) | 25,000.9 µg/mL | +/- 145.3496 +/- 1,508.4128 +/- 1,511.9941 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 11 | Methyl-tert-butyl ether (MTBE) CAS # 1634-04-4-SEC Purity 99% | (Lot ZAQTA-MS) | 2,500.0 µg/mL | +/- 14.5352 +/- 150.8361 +/- 151.1942 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 12 | cis-1,2-Dichloroethene CAS # 156-59-2-SEC Purity 98% | (Lot HGC01-BLKT) | 2,500.7 µg/mL | +/- 14.5394 +/- 150.8792 +/- 151.2374 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 13 | n-Hexane (C6) CAS # 110-54-3-SEC Purity 99% | (Lot 10188491) | 2,501.5 µg/mL | +/- 14.5439 +/- 150.9266 +/- 151.2849 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 14 | 1,1-Dichloroethane CAS # 75-34-3-SEC Purity 99% | (Lot 5379000) | 2,500.3 µg/mL | +/- 14.5367 +/- 150.8512 +/- 151.2093 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 15 | 2,2-Dichloropropane CAS # 594-20-7-SEC Purity 98% | (Lot I7E8E) | 2,500.1 µg/mL | +/- 14.5358 +/- 150.8423 +/- 151.2004 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 16 | trans-1,2-Dichloroethene CAS # 156-60-5-SEC Purity 97% | (Lot TS5UB) | 2,500.2 µg/mL | +/- 14.5362 +/- 150.8466 +/- 151.2048 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 17 | Isobutanol (2-Methyl-1-propanol) CAS # 78-83-1-SEC Purity 99% | (Lot 83NHH) | 62,506.9 µg/mL | +/- 363.4014 +/- 3,771.3149 +/- 3,780.2687 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 18 | Chloroform CAS # 67-66-3-SEC Purity 99% | (Lot 1297547) | 2,500.1 µg/mL | +/- 14.5359 +/- 150.8436 +/- 151.2017 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 19 | Bromochloromethane CAS # 74-97-5-SEC Purity 99% | (Lot 5670200) | 2,501.1 µg/mL | +/- 14.5418 +/- 150.9040 +/- 151.2622 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 20 | Tetrahydrofuran CAS # 109-99-9-SEC Purity 99% | (Lot K3V7J-SJ) | 5,002.3 µg/mL | +/- 29.0835 +/- 301.8079 +/- 302.5245 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 21 | 1,1,1-Trichloroethane CAS # 71-55-6-SEC Purity 98% | (Lot CS160712) | 2,500.7 µg/mL | +/- 14.5394 +/- 150.8792 +/- 151.2374 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 22 | Cyclohexane CAS # 110-82-7-SEC Purity 99% | (Lot YADRA) | 2,501.0 µg/mL | +/- 14.5410 +/- 150.8964 +/- 151.2547 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 23 | 1,1-Dichloropropene CAS # 563-58-6-SEC Purity 96% | (Lot 5221100) | 2,501.3 µg/mL | +/- 14.5427 +/- 150.9133 +/- 151.2716 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |

| | | | | | | |
|----|---|------------------|----------------|--|-------------------------|---------------------------------------|
| 24 | Carbon tetrachloride CAS # 56-23-5-SEC Purity 99% | (Lot 11466) | 2,500.5 µg/mL | +/- 14.5381 +/- 150.8662 +/- 151.2244 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 25 | n-Heptane (C7) CAS # 142-82-5-SEC Purity 99% | (Lot OGM01) | 2,500.5 µg/mL | +/- 14.5381 +/- 150.8662 +/- 151.2244 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 26 | 1,2-Dichloroethane CAS # 107-06-2-SEC Purity 99% | (Lot FO6PK) | 2,500.1 µg/mL | +/- 14.5359 +/- 150.8436 +/- 151.2017 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 27 | Benzene CAS # 71-43-2-SEC Purity 99% | (Lot B28Y008) | 2,501.5 µg/mL | +/- 14.5439 +/- 150.9266 +/- 151.2849 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 28 | Trichloroethylene CAS # 79-01-6-SEC Purity 99% | (Lot H04X050) | 2,501.0 µg/mL | +/- 14.5410 +/- 150.8964 +/- 151.2547 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 29 | Methylcyclohexane CAS # 108-87-2-SEC Purity 99% | (Lot 24MSD-CD) | 2,500.9 µg/mL | +/- 14.5403 +/- 150.8889 +/- 151.2471 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 30 | 1,2-Dichloropropane CAS # 78-87-5-SEC Purity 99% | (Lot OGG01) | 2,501.1 µg/mL | +/- 14.5418 +/- 150.9040 +/- 151.2622 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 31 | 1,4-Dioxane CAS # 123-91-1-SEC Purity 99% | (Lot MUFZH) | 50,007.1 µg/mL | +/- 290.7305 +/- 3,017.1500 +/- 3,024.3132 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 32 | Dibromomethane CAS # 74-95-3-SEC Purity 99% | (Lot FGI01-OICH) | 2,501.6 µg/mL | +/- 14.5447 +/- 150.9341 +/- 151.2925 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 33 | cis-1,3-Dichloropropene CAS # 10061-01-5-SEC Purity 99% | (Lot 487OA) | 2,500.1 µg/mL | +/- 14.5359 +/- 150.8436 +/- 151.2017 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 34 | Toluene CAS # 108-88-3-SEC Purity 99% | (Lot YND2B-BD) | 2,500.0 µg/mL | +/- 14.5352 +/- 150.8361 +/- 151.1942 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 35 | Ethyl methacrylate CAS # 97-63-2-SEC Purity 99% | (Lot MLWYK-LS) | 2,500.5 µg/mL | +/- 14.5381 +/- 150.8662 +/- 151.2244 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 36 | trans-1,3-Dichloropropene CAS # 10061-02-6-SEC Purity 99% | (Lot ZDMSL) | 2,500.5 µg/mL | +/- 14.5381 +/- 150.8662 +/- 151.2244 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 37 | 1,1,2-Trichloroethane CAS # 79-00-5-SEC Purity 98% | (Lot 5034600) | 2,500.8 µg/mL | +/- 14.5401 +/- 150.8866 +/- 151.2448 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 38 | 1,3-Dichloropropane CAS # 142-28-9-SEC Purity 99% | (Lot AGN01-EFPC) | 2,500.5 µg/mL | +/- 14.5381 +/- 150.8662 +/- 151.2244 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 39 | Tetrachloroethylene CAS # 127-18-4-SEC Purity 99% | (Lot F09W014) | 2,501.3 µg/mL | +/- 14.5425 +/- 150.9115 +/- 151.2698 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |

| | | | | | | | | |
|----|---|------------------|---------|-------|-----|----------|-------|-------------|
| 40 | Dibromochloromethane CAS # 124-48-1-SEC Purity 97% | (Lot 10181507) | 2,500.4 | µg/mL | +/- | 14.5376 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8613 | µg/mL | Unstressed |
| | | | | | +/- | 151.2194 | µg/mL | Stressed |
| 41 | 1,2-Dibromoethane (EDB) CAS # 106-93-4-SEC Purity 99% | (Lot 3505900) | 2,500.5 | µg/mL | +/- | 14.5381 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8662 | µg/mL | Unstressed |
| | | | | | +/- | 151.2244 | µg/mL | Stressed |
| 42 | Chlorobenzene CAS # 108-90-7-SEC Purity 99% | (Lot 1161936) | 2,501.0 | µg/mL | +/- | 14.5410 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8964 | µg/mL | Unstressed |
| | | | | | +/- | 151.2547 | µg/mL | Stressed |
| 43 | m-Xylene CAS # 108-38-3-SEC Purity 99% | (Lot OUKMG-GB) | 1,250.9 | µg/mL | +/- | 7.2727 | µg/mL | Gravimetric |
| | | | | | +/- | 75.4708 | µg/mL | Unstressed |
| | | | | | +/- | 75.6500 | µg/mL | Stressed |
| 44 | p-Xylene CAS # 106-42-3-SEC Purity 99% | (Lot GM01) | 1,250.5 | µg/mL | +/- | 7.2705 | µg/mL | Gravimetric |
| | | | | | +/- | 75.4482 | µg/mL | Unstressed |
| | | | | | +/- | 75.6273 | µg/mL | Stressed |
| 45 | Ethylbenzene CAS # 100-41-4-SEC Purity 99% | (Lot PI4SE) | 2,500.9 | µg/mL | +/- | 14.5403 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8889 | µg/mL | Unstressed |
| | | | | | +/- | 151.2471 | µg/mL | Stressed |
| 46 | 1,1,1,2-Tetrachloroethane CAS # 630-20-6-SEC Purity 99% | (Lot GC01) | 2,501.1 | µg/mL | +/- | 14.5418 | µg/mL | Gravimetric |
| | | | | | +/- | 150.9040 | µg/mL | Unstressed |
| | | | | | +/- | 151.2622 | µg/mL | Stressed |
| 47 | o-Xylene CAS # 95-47-6-SEC Purity 99% | (Lot FGL01-KTPK) | 2,500.9 | µg/mL | +/- | 14.5403 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8889 | µg/mL | Unstressed |
| | | | | | +/- | 151.2471 | µg/mL | Stressed |
| 48 | Styrene CAS # 100-42-5-SEC Purity 99% | (Lot OFIOL-IA) | 2,500.4 | µg/mL | +/- | 14.5374 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8587 | µg/mL | Unstressed |
| | | | | | +/- | 151.2169 | µg/mL | Stressed |
| 49 | Isopropylbenzene (cumene) CAS # 98-82-8-SEC Purity 99% | (Lot 2PHXG-IH) | 2,500.5 | µg/mL | +/- | 14.5381 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8662 | µg/mL | Unstressed |
| | | | | | +/- | 151.2244 | µg/mL | Stressed |
| 50 | Bromoform CAS # 75-25-2-SEC Purity 99% | (Lot 5139000) | 2,502.3 | µg/mL | +/- | 14.5483 | µg/mL | Gravimetric |
| | | | | | +/- | 150.9718 | µg/mL | Unstressed |
| | | | | | +/- | 151.3303 | µg/mL | Stressed |
| 51 | Bromodichloromethane CAS # 75-27-4-SEC Purity 98% | (Lot 13780) | 2,500.1 | µg/mL | +/- | 14.5358 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8423 | µg/mL | Unstressed |
| | | | | | +/- | 151.2004 | µg/mL | Stressed |
| 52 | 1,1,2,2-Tetrachloroethane CAS # 79-34-5-SEC Purity 99% | (Lot CFA4D-AQ) | 2,501.3 | µg/mL | +/- | 14.5425 | µg/mL | Gravimetric |
| | | | | | +/- | 150.9115 | µg/mL | Unstressed |
| | | | | | +/- | 151.2698 | µg/mL | Stressed |
| 53 | 1,2,3-Trichloropropane CAS # 96-18-4-SEC Purity 98% | (Lot OGI01) | 2,500.1 | µg/mL | +/- | 14.5358 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8423 | µg/mL | Unstressed |
| | | | | | +/- | 151.2004 | µg/mL | Stressed |
| 54 | trans-1,4-Dichloro-2-butene CAS # 110-57-6-SEC Purity 98% | (Lot 100700-3) | 2,501.0 | µg/mL | +/- | 14.5408 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8940 | µg/mL | Unstressed |
| | | | | | +/- | 151.2522 | µg/mL | Stressed |
| 55 | n-Propylbenzene CAS # 103-65-1-SEC Purity 99% | (Lot T2HFC-II) | 2,500.0 | µg/mL | +/- | 14.5352 | µg/mL | Gravimetric |
| | | | | | +/- | 150.8361 | µg/mL | Unstressed |
| | | | | | +/- | 151.1942 | µg/mL | Stressed |

| | | | | | | | |
|----|--|-----------------|---------|-------|---|-------------------------|---------------------------------------|
| 56 | Bromobenzene CAS # 108-86-1-SEC Purity 99% | (Lot 2FUHG-EM) | 2,500.1 | µg/mL | +/- 14.5359 +/- 150.8436 +/- 151.2017 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 57 | 1,3,5-Trimethylbenzene CAS # 108-67-8-SEC Purity 99% | (Lot TOOOF) | 2,500.3 | µg/mL | +/- 14.5367 +/- 150.8512 +/- 151.2093 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 58 | 2-Chlorotoluene CAS # 95-49-8-SEC Purity 99% | (Lot SW8QG-AO) | 2,500.9 | µg/mL | +/- 14.5403 +/- 150.8889 +/- 151.2471 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 59 | 4-Chlorotoluene CAS # 106-43-4-SEC Purity 99% | (Lot P4XHJ-AO) | 2,500.5 | µg/mL | +/- 14.5381 +/- 150.8662 +/- 151.2244 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 60 | tert-Butylbenzene CAS # 98-06-6-SEC Purity 99% | (Lot OGN01-CAI) | 2,500.1 | µg/mL | +/- 14.5359 +/- 150.8436 +/- 151.2017 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 61 | 1,2,4-Trimethylbenzene CAS # 95-63-6-SEC Purity 99% | (Lot SC7LO-QA) | 2,500.4 | µg/mL | +/- 14.5374 +/- 150.8587 +/- 151.2169 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 62 | sec-Butylbenzene CAS # 135-98-8-SEC Purity 99% | (Lot OGN01-IMA) | 2,501.4 | µg/mL | +/- 14.5432 +/- 150.9190 +/- 151.2773 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 63 | 4-Isopropyltoluene (p-cymene) CAS # 99-87-6-SEC Purity 99% | (Lot 5221800) | 2,501.3 | µg/mL | +/- 14.5425 +/- 150.9115 +/- 151.2698 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 64 | 1,3-Dichlorobenzene CAS # 541-73-1-SEC Purity 99% | (Lot FMDFD) | 2,500.9 | µg/mL | +/- 14.5403 +/- 150.8889 +/- 151.2471 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 65 | 1,4-Dichlorobenzene CAS # 106-46-7-SEC Purity 99% | (Lot 4Y5DC) | 2,500.8 | µg/mL | +/- 14.5396 +/- 150.8813 +/- 151.2395 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 66 | n-Butylbenzene CAS # 104-51-8-SEC Purity 99% | (Lot OGN01-PNP) | 2,500.8 | µg/mL | +/- 14.5396 +/- 150.8813 +/- 151.2395 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 67 | 1,2-Dichlorobenzene CAS # 95-50-1-SEC Purity 99% | (Lot R6QDM) | 2,501.0 | µg/mL | +/- 14.5410 +/- 150.8964 +/- 151.2547 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 68 | 1,2-Dibromo-3-chloropropane CAS # 96-12-8-SEC Purity 98% | (Lot LC00408V) | 2,501.5 | µg/mL | +/- 14.5436 +/- 150.9236 +/- 151.2819 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 69 | 1,2,4-Trichlorobenzene CAS # 120-82-1-SEC Purity 99% | (Lot 3LYYC) | 2,502.5 | µg/mL | +/- 14.5498 +/- 150.9869 +/- 151.3454 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 70 | Hexachlorobutadiene CAS # 87-68-3-SEC Purity 97% | (Lot 5526800) | 2,501.4 | µg/mL | +/- 14.5433 +/- 150.9198 +/- 151.2781 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 71 | Naphthalene CAS # 91-20-3-SEC Purity 99% | (Lot SKZ5N) | 2,501.8 | µg/mL | +/- 14.5454 +/- 150.9417 +/- 151.3000 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |

| | | | | | | | |
|----|------------------------|----------------|---------|-------|--------------|-------|-------------|
| 72 | 1,2,3-Trichlorobenzene | | 2,500.7 | µg/mL | +/- 14.5394 | µg/mL | Gravimetric |
| | CAS # 87-61-6.SEC | (Lot A0043055) | | | +/- 150.8792 | µg/mL | Unstressed |
| | Purity 98% | | | | +/- 151.2374 | 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)

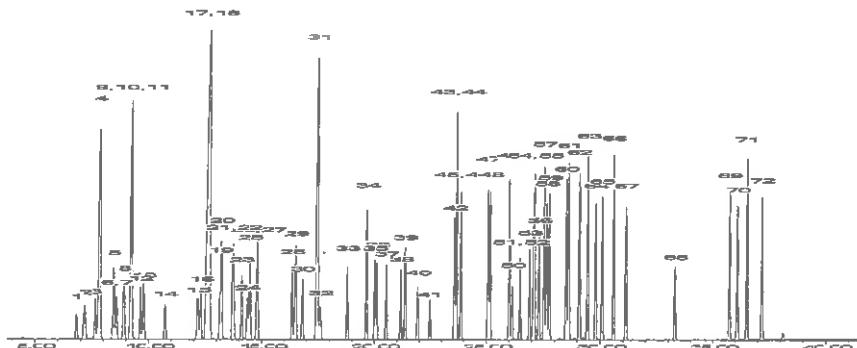
Carrier Gas:
 helium-constant pressure 30 psi

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



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 Maye

Date Mixed: 28-Dec-2016 Balance: 1127510105

Jennifer J. Pollino
 Jennifer Pollino - Operations Tech-ARM QC

Date Passed: 04-Jan-2017

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

Reagent

VOA8260SURRES_00118



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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>567650</u> | Lot No.: | <u>A0114901</u> |
| Description : | <u>8260 Surrogate Standard</u> | | |
| 8260 Surrogate Standard 2,500 ug/ml, P&T Methanol, 5 ml/ampul | | | |
| Container Size : | <u>5 mL</u> | Pkg Amt: | <u>> 5 mL</u> |
| Expiration Date : | <u>October 31, 2020</u> | | |
| | Storage: <u>0°C or colder</u> | | |

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

| Elation Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L.; K=2) | | |
|---------------|-------------------------------|--------------------------------|---|----------|-------|
| 1 | Dibromofluoromethane | 2,509.4 µg/mL | +/- | 14.5899 | µg/mL |
| | CAS # 1868-53-7 | | +/- | 140.6996 | µg/mL |
| | Purity 99% | | +/- | 143.9918 | µg/mL |
| 2 | 1,2-Dichloroethane-d4 | 2,509.0 µg/mL | +/- | 14.5875 | µg/mL |
| | CAS # 17060-07-0 | | +/- | 140.6769 | µg/mL |
| | Purity 98% | | +/- | 143.9686 | µg/mL |
| 3 | Toluene-d8 | 2,507.0 µg/mL | +/- | 14.5759 | µg/mL |
| | CAS # 2037-26-5 | | +/- | 140.5650 | µg/mL |
| | Purity 99% | | +/- | 143.8540 | µg/mL |
| 4 | 1-Bromo-4-fluorobenzene (BFB) | 2,503.6 µg/mL | +/- | 14.5561 | µg/mL |
| | CAS # 460-00-4 | | +/- | 140.3744 | µg/mL |
| | Purity 99% | | +/- | 143.6590 | µg/mL |
| Solvent: | P&T Methanol | | | | |
| | CAS # 67-56-1 | | | | |
| | Purity 99% | | | | |

Reagent

VOA8260SURRES_00122



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Catalog No.: 567650 **Lot No.:** A0114901
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 : October 31, 2020 **Storage:** 0°C or colder

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

| Elation Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L.; K=2) | | |
|---|-------------------------------|--------------------------------|---|----------|-------|
| 1 | Dibromofluoromethane | 2,509.4 µg/mL | +/- | 14.5899 | µg/mL |
| | CAS # 1868-53-7 | | +/- | 140.6996 | µg/mL |
| | Purity 99% | | +/- | 143.9918 | µg/mL |
| 2 | 1,2-Dichloroethane-d4 | 2,509.0 µg/mL | +/- | 14.5875 | µg/mL |
| | CAS # 17060-07-0 | | +/- | 140.6769 | µg/mL |
| | Purity 98% | | +/- | 143.9686 | µg/mL |
| 3 | Toluene-d8 | 2,507.0 µg/mL | +/- | 14.5759 | µg/mL |
| | CAS # 2037-26-5 | | +/- | 140.5650 | µg/mL |
| | Purity 99% | | +/- | 143.8540 | µg/mL |
| 4 | 1-Bromo-4-fluorobenzene (BFB) | 2,503.6 µg/mL | +/- | 14.5561 | µg/mL |
| | CAS # 460-00-4 | | +/- | 140.3744 | µg/mL |
| | Purity 99% | | +/- | 143.6590 | µg/mL |
| Solvent: P&T Methanol CAS # 67-56-1 Purity 99% | | | | | |

Reagent

VOA8260VARES_00082



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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.:** A0124520

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, 2017 **Storage:** 0°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 | Vinyl acetate CAS # 108-05-4 Purity 99% | 5,027.0 µg/mL | +/- 29.5013 µg/mL | +/- 303.3277 µg/mL | +/- 304.0477 µg/mL |

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

VOAACRORES_00115



CERTIFIED REFERENCE MATERIAL

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Catalog No. : 568720

Lot No.: A0125560

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 : September 30, 2017

Storage: 0°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 | 19,779.0 µg/mL | +/- 115.8104 | µg/mL | Gravimetric |
| | CAS # 107-02-8 | (Lot 170123JLM) | +/- 634.1769 | µg/mL | Unstressed |
| | Purity 99% | | +/- 737.1613 | µg/mL | Stressed |

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

Reagent

VOABFBRES_00052

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Catalog No.: 30067

Lot No.: A0119122

Description : 4-Bromofluorobenzene Standard

4-Bromofluorobenzene Standard 2,500 μ g/mL, P&T Methanol,
 1mL/ampul

Container Size : 2 mL

Pkg Amt: > 1 mL

Expiration Date : May 31, 2021

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 | 1-Bromo-4-fluorobenzene (BFB) CAS # 460-00-4 Purity 99% | 2,501.0 μ g/mL | +/- 14.6773 μ g/mL | +/- 140.2428 μ g/mL | +/- 143.5236 μ g/mL |

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

Reagent

VOABFBRES_00055



CERTIFIED REFERENCE MATERIAL

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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. : 30067

Lot No.: A0122647

Description : 4-Bromofluorobenzene Standard

4-Bromofluorobenzene Standard 2,500 μ g/mL, P&T Methanol,
1mL/ampul

Container Size : 2 mL

Pkg Amt: > 1 mL

Expiration Date : November 30, 2021

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 | 1-Bromo-4-fluorobenzene (BFB) CAS # 460-00-4 Purity 99% | 2,524.0 μ g/mL | +/- 14.8122 μ g/mL | +/- 141.5325 μ g/mL | Gravimetric Unstressed Stressed |

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

Reagent

VOABFBRES_00058



CERTIFIED REFERENCE MATERIAL

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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. : 30067

Lot No.: A0122647

Description : 4-Bromofluorobenzene Standard

4-Bromofluorobenzene Standard 2,500 μ g/mL, P&T Methanol,
1mL/ampul

Container Size : 2 mL

Pkg Amt: > 1 mL

Expiration Date : November 30, 2021

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 | 1-Bromo-4-fluorobenzene (BFB) CAS # 460-00-4 Purity 99% | 2,524.0 μ g/mL | +/- 14.8122 μ g/mL | +/- 141.5325 μ g/mL | Gravimetric Unstressed Stressed |

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

Reagent

VOACEVERES_00127



CERTIFIED REFERENCE MATERIAL

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

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. : 569723

Lot No.: A0123891

Description : 8260 List 1 / Std #4 2-CEVE (2015)

8260 List 1 / Std #4 2-CEVE (2015) 2,500 ug/ml, P&T Methanol, 1 ml/ampul

Container Size : 2 mL

Pkg Amt: > 1 mL

Expiration Date : January 31, 2020

Storage: 0°C or colder

1/24/2027

28

29

30

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 | 2-Chloroethyl vinyl ether CAS # 110-75-8 Purity 98% | 2,503.5 µg/mL (Lot MKBS6526V) | +/- 14.5556 µg/mL | +/- 53.6004 µg/mL | +/- 55.1587 µg/mL |

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

Tech Tips:

Degradation of tetrachloroethylene to pentachloroethane may occur if solutions containing 2-chloroethyl vinyl ether are combined with solutions that contain tetrachloroethylene.

Reagent

VOARESEE1ST_00045



CERTIFIED REFERENCE MATERIAL

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Catalog No.: 568363-FL **Lot No.:** A0120234 239675

Description : Custom EE Standard

Custom EE Standard 5,000 μ g/mL, P&T Methanol, 1mL/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 | 3-Chlorobenzotrifluoride CAS # 98-15-7 Purity 99% | 5,025.0 μ g/mL | +/- 29.4895 μ g/mL | +/- 281.7753 μ g/mL | +/- 288.3671 μ g/mL |
| | (Lot 21324DO) | | | | |
| 2 | 4-Chlorobenzotrifluoride CAS # 98-56-6 Purity 99% | 5,031.0 μ g/mL | +/- 29.5247 μ g/mL | +/- 282.1117 μ g/mL | +/- 288.7115 μ g/mL |
| | (Lot 08507BO) | | | | |
| 3 | 2-Chlorobenzotrifluoride CAS # 88-16-4 Purity 99% | 5,011.0 μ g/mL | +/- 29.4074 μ g/mL | +/- 280.9902 μ g/mL | +/- 287.5637 μ g/mL |
| | (Lot I0316DQ) | | | | |
| 4 | 3-Chlorotoluene CAS # 108-41-8 Purity 99% | 5,046.0 μ g/mL | +/- 29.6128 μ g/mL | +/- 282.9528 μ g/mL | +/- 289.5723 μ g/mL |
| | (Lot 13528LX) | | | | |
| 5 | 2,4-Dichlorobenzotrifluoride CAS # 320-60-5 Purity 99% | 5,018.0 μ g/mL | +/- 29.4484 μ g/mL | +/- 281.3828 μ g/mL | +/- 287.9654 μ g/mL |
| | (Lot MKBL3552V) | | | | |
| 6 | 3,4-Dichlorobenzotrifluoride CAS # 328-84-7 Purity 99% | 5,031.0 μ g/mL | +/- 29.5247 μ g/mL | +/- 282.1117 μ g/mL | +/- 288.7115 μ g/mL |
| | (Lot 11105EJV) | | | | |
| 7 | 2,5-Dichlorobenzotrifluoride CAS # 320-50-3 Purity 99% | 5,047.0 μ g/mL | +/- 29.6186 μ g/mL | +/- 283.0089 μ g/mL | +/- 289.6296 μ g/mL |
| | (Lot 04415DSV) | | | | |

Method 8260C Low Level

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

FORM II
GC/MS VOA SURROGATE RECOVERY

Lab Name: TestAmerica Pittsburgh Job No.: 180-71131-1
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-SPBA-CW-23-0/1-0 | 180-71131-1 | 93 | 94 | 86 | 91 |
| HD-CW-23-0/1-0 | 180-71131-2 | 93 | 94 | 89 | 91 |
| HD-QC6-0/1-2 | 180-71131-3 | 95 | 95 | 87 | 89 |
| | MB 180-226148/6 | 91 | 93 | 89 | 90 |
| | LCS 180-226148/4 | 92 | 92 | 93 | 94 |

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

QC LIMITS

| |
|--------|
| 73-120 |
| 65-121 |
| 73-120 |
| 80-120 |

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.: 180-71131-1
SDG No.: _____
Matrix: Water Level: Low Lab File ID: 6101804D.D
Lab ID: LCS 180-226148/4 Client ID: _____

| COMPOUND | SPIKE ADDED (ug/L) | LCS CONCENTRATION (ug/L) | LCS % REC | QC LIMITS REC | # |
|-----------------------------|--------------------------|--------------------------------|-----------------|---------------------|---|
| Chloromethane | 10.0 | 7.89 | 79 | 49-135 | |
| Vinyl chloride | 10.0 | 8.87 | 89 | 52-136 | |
| Bromomethane | 10.0 | 7.90 | 79 | 37-150 | |
| Chloroethane | 10.0 | 9.09 | 91 | 44-139 | |
| 1,1-Dichloroethene | 10.0 | 9.66 | 97 | 64-131 | |
| Acetone | 20.0 | 22.4 | 112 | 24-150 | |
| Carbon disulfide | 10.0 | 8.89 | 89 | 20-150 | |
| Methylene Chloride | 10.0 | 8.98 | 90 | 66-123 | |
| trans-1,2-Dichloroethene | 10.0 | 9.35 | 94 | 70-123 | |
| Methyl tert-butyl ether | 10.0 | 8.28 | 83 | 66-130 | |
| 1,1-Dichloroethane | 10.0 | 9.00 | 90 | 66-122 | |
| cis-1,2-Dichloroethene | 10.0 | 9.03 | 90 | 73-120 | |
| Bromochloromethane | 10.0 | 9.10 | 91 | 73-122 | |
| 2-Butanone (MEK) | 20.0 | 18.3 | 91 | 37-150 | |
| Chloroform | 10.0 | 9.63 | 96 | 72-123 | |
| 1,1,1-Trichloroethane | 10.0 | 10.4 | 104 | 66-129 | |
| Carbon tetrachloride | 10.0 | 12.2 | 122 | 58-145 | |
| Benzene | 10.0 | 9.47 | 95 | 75-123 | |
| 1,2-Dichloroethane | 10.0 | 9.07 | 91 | 63-130 | |
| Trichloroethene | 10.0 | 9.21 | 92 | 74-121 | |
| 1,2-Dichloropropane | 10.0 | 8.45 | 84 | 67-119 | |
| Bromodichloromethane | 10.0 | 9.26 | 93 | 62-127 | |
| cis-1,3-Dichloropropene | 10.0 | 9.17 | 92 | 61-127 | |
| 4-Methyl-2-pentanone (MIBK) | 20.0 | 13.2 | 66 | 41-135 | |
| Toluene | 10.0 | 9.31 | 93 | 76-129 | |
| trans-1,3-Dichloropropene | 10.0 | 9.36 | 94 | 61-136 | |
| 1,1,2-Trichloroethane | 10.0 | 8.74 | 87 | 74-126 | |
| Tetrachloroethene | 10.0 | 9.13 | 91 | 76-128 | |
| 2-Hexanone | 20.0 | 18.9 | 94 | 37-150 | |
| Dibromochloromethane | 10.0 | 9.51 | 95 | 63-131 | |
| 1,2-Dibromoethane (EDB) | 10.0 | 8.97 | 90 | 76-128 | |
| Chlorobenzene | 10.0 | 9.44 | 94 | 79-124 | |
| 1,1,1,2-Tetrachloroethane | 10.0 | 10.1 | 101 | 70-130 | |
| Ethylbenzene | 10.0 | 9.18 | 92 | 77-124 | |
| Xylenes, Total | 20.0 | 18.2 | 91 | 76-124 | |
| Styrene | 10.0 | 9.53 | 95 | 80-125 | |
| Bromoform | 10.0 | 8.29 | 83 | 54-136 | |
| 1,1,2,2-Tetrachloroethane | 10.0 | 8.83 | 88 | 72-128 | |
| Acrylonitrile | 100 | 65.5 | 65 | 60-130 | |
| 1,4-Dioxane | 200 | 144 J | 72 | 26-150 | |

Column to be used to flag recovery and RPD values

FORM III 8260C

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Pittsburgh Job No.: 180-71131-1
SDG No.: _____
Lab File ID: 6101806D.D Lab Sample ID: MB 180-226148/6
Matrix: Water Heated Purge: (Y/N) N
Instrument ID: CHHP6 Date Analyzed: 10/18/2017 02:09
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 180-226148/4 | 6101804D.D | 10/18/2017 01:04 |
| HD-SPBA-CW-23-0/1-0 | 180-71131-1 | 6101816D.D | 10/18/2017 06:33 |
| HD-CW-23-0/1-0 | 180-71131-2 | 6101817D.D | 10/18/2017 06:57 |
| HD-QC6-0/1-2 | 180-71131-3 | 6101819D.D | 10/18/2017 07:45 |

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

Lab Name: TestAmerica Pittsburgh Job No.: 180-71131-1
SDG No.: _____
Lab File ID: 60602003.D BFB Injection Date: 06/02/2017
Instrument ID: CHHP6 BFB Injection Time: 06:02
Analysis Batch No.: 213005

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 50 | 15.0 - 40.0 % of mass 95 | 18.8 |
| 75 | 30.0 - 60.0 % of mass 95 | 48.4 |
| 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 | 0.0 (0.0) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 65.4 |
| 175 | 5.0 - 9.0 % of mass 174 | 5.1 (7.9) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 65.5 (100.3) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 5.0 (7.7) 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 |
|------------------|-------------------|-------------|---------------|---------------|
| | ICV 180-213005/23 | 60602023.D | 06/02/2017 | 15:14 |

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

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

Lab File ID: 60724D01.D BFB Injection Date: 07/24/2017

Instrument ID: CHHP6 BFB Injection Time: 04:58

Analysis Batch No.: 217861

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 50 | 15.0 - 40.0 % of mass 95 | 17.3 |
| 75 | 30.0 - 60.0 % of mass 95 | 52.6 |
| 95 | Base Peak, 100% relative abundance | 100.0 |
| 96 | 5.0 - 9.0 % of mass 95 | 7.7 |
| 173 | Less than 2.0 % of mass 174 | 0.3 (0.4) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 68.6 |
| 175 | 5.0 - 9.0 % of mass 174 | 4.4 (6.4) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 68.1 (99.3) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.9 (7.1) 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 180-217861/3 | 60724D03.D | 07/24/2017 | 06:39 |
| | IC 180-217861/4 | 60724D04.D | 07/24/2017 | 07:03 |
| | ICIS 180-217861/5 | 60724D05.D | 07/24/2017 | 07:27 |
| | IC 180-217861/6 | 60724D06.D | 07/24/2017 | 07:52 |
| | IC 180-217861/7 | 60724D07.D | 07/24/2017 | 08:16 |
| | IC 180-217861/8 | 60724D08.D | 07/24/2017 | 08:40 |
| | IC 180-217861/9 | 60724D09.D | 07/24/2017 | 09:04 |
| | IC 180-217861/10 | 60724D10.D | 07/24/2017 | 09:28 |
| | ICV 180-217861/13 | 60724D13.D | 07/24/2017 | 10:40 |

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

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

Lab File ID: 6101801D.D BFB Injection Date: 10/17/2017

Instrument ID: CHHP6 BFB Injection Time: 21:58

Analysis Batch No.: 226148

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 50 | 15.0 - 40.0 % of mass 95 | 15.4 |
| 75 | 30.0 - 60.0 % of mass 95 | 52.8 |
| 95 | Base Peak, 100% relative abundance | 100.0 |
| 96 | 5.0 - 9.0 % of mass 95 | 7.6 |
| 173 | Less than 2.0 % of mass 174 | 0.0 (0.0) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 62.8 |
| 175 | 5.0 - 9.0 % of mass 174 | 4.8 (7.7) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 60.6 (96.6) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 3.4 (5.5) 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 180-226148/2 | 6101802D.D | 10/17/2017 | 23:55 |
| | ICS 180-226148/4 | 6101804D.D | 10/18/2017 | 01:04 |
| | MB 180-226148/6 | 6101806D.D | 10/18/2017 | 02:09 |
| HD-SPBA-CW-23-0/1-0 | 180-71131-1 | 6101816D.D | 10/18/2017 | 06:33 |
| HD-CW-23-0/1-0 | 180-71131-2 | 6101817D.D | 10/18/2017 | 06:57 |
| HD-QC6-0/1-2 | 180-71131-3 | 6101819D.D | 10/18/2017 | 07:45 |

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

Lab Name: TestAmerica Pittsburgh Job No.: 180-71131-1
SDG No.: _____
Sample No.: ICIS 180-217861/5 Date Analyzed: 07/24/2017 07:27
Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm)
Lab File ID (Standard): 60724D05.D Heated Purge: (Y/N) N
Calibration ID: 35029

| | TBAd9 | | FB | | CBNZd5 | | |
|-------------------------------|------------------|--------|---------|--------|--------|--------|-------|
| | AREA # | RT # | AREA # | RT # | AREA # | RT # | |
| INITIAL CALIBRATION MID-POINT | 281180 | 3.98 | 854988 | 7.05 | 187443 | 10.17 | |
| UPPER LIMIT | 562360 | 4.48 | 1709976 | 7.55 | 374886 | 10.67 | |
| LOWER LIMIT | 140590 | 3.48 | 427494 | 6.55 | 93722 | 9.67 | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 180-217861/13 | | 317387 | 3.96 | 870770 | 7.05 | 217522 | 10.17 |

TBAd9 = TBA-d9 (IS)

FB = Fluorobenzene (IS)

CBNZd5 = 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.: 180-71131-1
SDG No.: _____
Sample No.: ICIS 180-217861/5 Date Analyzed: 07/24/2017 07:27
Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm)
Lab File ID (Standard): 60724D05.D Heated Purge: (Y/N) N
Calibration ID: 35029

| | DCBd4 | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
|-------------------------------|------------------|--------|--------|------|--------|------|--------|------|
| | AREA # | RT # | | | | | | |
| INITIAL CALIBRATION MID-POINT | 265638 | 12.51 | | | | | | |
| UPPER LIMIT | 531276 | 13.01 | | | | | | |
| LOWER LIMIT | 132819 | 12.01 | | | | | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | | |
| ICV 180-217861/13 | | 360072 | 12.51 | | | | | |

DCBd4 = 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.: 180-71131-1
SDG No.: _____
Sample No.: CCVIS 180-226148/2 Date Analyzed: 10/17/2017 23:55
Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm)
Lab File ID (Standard): 6101802D.D Heated Purge: (Y/N) N
Calibration ID: 35029

| | TBAd9 | | FB | | CBNZd5 | |
|------------------|---------------------|--------|---------|---------|--------|--------|
| | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | 360770 | 3.95 | 1145374 | 7.04 | 290716 | 10.17 |
| UPPER LIMIT | 721540 | 4.45 | 2290748 | 7.54 | 581432 | 10.67 |
| LOWER LIMIT | 180385 | 3.45 | 572687 | 6.54 | 145358 | 9.67 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | |
| LCS 180-226148/4 | | 373208 | 3.96 | 1246060 | 7.05 | 301504 |
| MB 180-226148/6 | | 312973 | 3.96 | 1062992 | 7.05 | 286612 |
| 180-71131-1 | HD-SPBA-CW-23-0/1-0 | 317182 | 3.95 | 1010234 | 7.05 | 279053 |
| 180-71131-2 | HD-CW-23-0/1-0 | 286138 | 3.95 | 994351 | 7.05 | 272894 |
| 180-71131-3 | HD-QC6-0/1-2 | 274765 | 3.96 | 988524 | 7.05 | 275892 |

TBAd9 = TBA-d9 (IS)

FB = Fluorobenzene (IS)

CBNZd5 = 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.: 180-71131-1
SDG No.: _____
Sample No.: CCVIS 180-226148/2 Date Analyzed: 10/17/2017 23:55
Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm)
Lab File ID (Standard): 6101802D.D Heated Purge: (Y/N) N
Calibration ID: 35029

| | DCBd4 | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
|------------------|---------------------|--------|--------|------|--------|------|--------|------|
| | AREA # | RT # | | | | | | |
| 12/24 HOUR STD | 448308 | 12.51 | | | | | | |
| UPPER LIMIT | 896616 | 13.01 | | | | | | |
| LOWER LIMIT | 224154 | 12.01 | | | | | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | | |
| LCS 180-226148/4 | | 469268 | 12.51 | | | | | |
| MB 180-226148/6 | | 425493 | 12.51 | | | | | |
| 180-71131-1 | HD-SPBA-CW-23-0/1-0 | 406630 | 12.51 | | | | | |
| 180-71131-2 | HD-CW-23-0/1-0 | 392535 | 12.51 | | | | | |
| 180-71131-3 | HD-QC6-0/1-2 | 388863 | 12.51 | | | | | |

DCBd4 = 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.: 180-71131-1

SDG No.: _____

Client Sample ID: HD-SPBA-CW-23-0/1-0

Lab Sample ID: 180-71131-1

Matrix: Water

Lab File ID: 6101816D.D

Analysis Method: 8260C

Date Collected: 10/05/2017 09:50

Sample wt/vol: 5 (mL)

Date Analyzed: 10/18/2017 06:33

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.: 226148

Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|------------|-----------------------------|--------|------|-----|------|
| 74-87-3 | Chloromethane | 1.0 | U | 1.0 | 0.90 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 1.0 | 0.88 |
| 74-83-9 | Bromomethane | 1.0 | U | 1.0 | 0.89 |
| 75-00-3 | Chloroethane | 1.0 | U | 1.0 | 0.90 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.55 |
| 67-64-1 | Acetone | 5.0 | U ^c | 5.0 | 3.4 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 1.0 | 0.88 |
| 75-09-2 | Methylene Chloride | 1.0 | U | 1.0 | 0.36 |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.67 |
| 1634-04-4 | Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.59 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.63 |
| 156-59-2 | cis-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.71 |
| 74-97-5 | Bromochloromethane | 1.0 | U | 1.0 | 0.63 |
| 78-93-3 | 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 |
| 67-66-3 | Chloroform | 1.0 | U | 1.0 | 0.60 |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.60 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U ^c | 1.0 | 0.88 |
| 71-43-2 | Benzene | 1.0 | U | 1.0 | 0.60 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.57 |
| 79-01-6 | Trichloroethene | 4.1 | | 1.0 | 0.69 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.66 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 1.0 | 0.64 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.59 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 5.0 | U ^c | 5.0 | 3.1 |
| 108-88-3 | Toluene | 1.0 | U | 1.0 | 0.46 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.58 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.45 |
| 127-18-4 | Tetrachloroethene | 42 | | 1.0 | 0.47 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 5.0 | 3.3 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 1.0 | 0.84 |
| 106-93-4 | 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.50 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 1.0 | 0.50 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.57 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 1.0 | 0.51 |
| 1330-20-7 | Xylenes, Total | 2.0 | U | 2.0 | 0.89 |
| 100-42-5 | Styrene | 1.0 | U | 1.0 | 0.47 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.: _____

Client Sample ID: HD-SPBA-CW-23-0/1-0

Lab Sample ID: 180-71131-1

Matrix: Water

Lab File ID: 6101816D.D

Analysis Method: 8260C

Date Collected: 10/05/2017 09:50

Sample wt/vol: 5 (mL)

Date Analyzed: 10/18/2017 06:33

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.: 226148

Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|----------|---------------------------|--------|------|-----|------|
| 75-25-2 | Bromoform | 1.0 | U | 1.0 | 0.98 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.60 |
| 107-13-1 | Acrylonitrile | 20 | U ^c | 20 | 7.8 |
| 123-91-1 | 1,4-Dioxane | 200 | U | 200 | 14 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|------------|------------------------------|------|---|--------|
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 94 | | 65-121 |
| 2037-26-5 | Toluene-d8 (Surr) | 86 | | 73-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 91 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 93 | | 73-120 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101816.D.D
 Lims ID: 180-71131-B-1
 Client ID: HD-SPBA-CW-23-0/1-0
 Sample Type: Client
 Inject. Date: 18-Oct-2017 06:33:30 ALS Bottle#: 25 Worklist Smp#: 16
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018914-016
 Misc. Info.: 180-71131-B-1
 Operator ID: 034635 Instrument ID: CHHP6
 Method: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 18-Oct-2017 20:30:42 Calib Date: 28-Sep-2017 15:13:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICAL File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170928-18631.b\\60928P06.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK031

First Level Reviewer: bungardf Date: 18-Oct-2017 20:23:32

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 3.954 | 3.949 | 0.005 | 91 | 317182 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.045 | 7.045 | 0.000 | 98 | 1010234 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.165 | 10.166 | -0.001 | 88 | 279053 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.508 | 12.508 | 0.000 | 96 | 406630 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr) | 113 | 6.315 | 6.314 | 0.001 | 92 | 244987 | 46.7 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur) | 65 | 6.692 | 6.685 | 0.007 | 70 | 353516 | 47.1 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.712 | 8.711 | 0.001 | 93 | 974331 | 43.1 | |
| \$ 8 4-Bromofluorobenzene (Surr) | 95 | 11.352 | 11.351 | 0.001 | 82 | 436416 | 45.3 | |
| 12 Chloromethane | 50 | | 1.624 | | | | ND | |
| 13 Vinyl chloride | 62 | | 1.740 | | | | ND | |
| 15 Bromomethane | 94 | | 2.068 | | | | ND | |
| 16 Chloroethane | 64 | | 2.202 | | | | ND | |
| 22 1,1-Dichloroethene | 96 | | 3.090 | | | | ND | |
| 24 Acetone | 43 | 3.188 | 3.181 | 0.007 | 78 | 9414 | 4.34 | M |
| 26 Carbon disulfide | 76 | | 3.364 | | | | ND | |
| 31 Methylene Chloride | 84 | | 3.838 | | | | ND | |
| 33 Acrylonitrile | 53 | | 4.222 | | | | ND | |
| 34 trans-1,2-Dichloroethene | 96 | | 4.270 | | | | ND | |
| 35 Methyl tert-butyl ether | 73 | | 4.276 | | | | ND | |
| 37 1,1-Dichloroethane | 63 | | 4.921 | | | | ND | |
| 43 cis-1,2-Dichloroethene | 96 | | 5.688 | | | | ND | |
| 44 2-Butanone (MEK) | 43 | | 5.694 | | | | ND | |
| 48 Chlorobromomethane | 128 | | 5.974 | | | | ND | |
| 50 Chloroform | 83 | 6.126 | 6.126 | 0.000 | 88 | 21828 | 2.09 | |
| 51 1,1,1-Trichloroethane | 97 | | 6.290 | | | | ND | |
| 53 Carbon tetrachloride | 117 | | 6.454 | | | | ND | |
| 56 Benzene | 78 | | 6.691 | | | | ND | |
| 57 1,2-Dichloroethane | 62 | | 6.777 | | | | ND | |
| 61 Trichloroethene | 130 | 7.440 | 7.440 | 0.000 | 97 | 117362 | 20.7 | |
| 64 1,2-Dichloropropane | 63 | | 7.713 | | | | ND | |
| 65 1,4-Dioxane | 88 | | 7.793 | | | | ND | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | OnCol Amt ng | Flags |
|--------------------------------|-----|--------------|------------------|------------------|----|----------|-----------------|-------|
| 68 Dichlorobromomethane | 83 | | 7.999 | | | | ND | |
| 71 cis-1,3-Dichloropropene | 75 | | 8.450 | | | | ND | |
| 72 4-Methyl-2-pentanone (MIBK) | 43 | | 8.608 | | | | ND | |
| 73 Toluene | 91 | 8.785 | 8.778 | 0.007 | 98 | 10505 | 0.3820 | |
| 74 trans-1,3-Dichloropropene | 75 | | 9.034 | | | | ND | |
| 76 1,1,2-Trichloroethane | 97 | | 9.222 | | | | ND | |
| 77 Tetrachloroethene | 164 | 9.296 | 9.295 | 0.001 | 90 | 1033054 | 209.8 | |
| 79 2-Hexanone | 43 | | 9.447 | | | | ND | |
| 81 Chlorodibromomethane | 129 | | 9.593 | | | | ND | |
| 82 Ethylene Dibromide | 107 | | 9.709 | | | | ND | |
| 84 Chlorobenzene | 112 | | 10.196 | | | | ND | |
| 86 1,1,1,2-Tetrachloroethane | 131 | | 10.293 | | | | ND | |
| 87 Ethylbenzene | 106 | | 10.299 | | | | ND | |
| 88 m-Xylene & p-Xylene | 106 | | 10.433 | | | | ND | |
| 89 o-Xylene | 106 | | 10.816 | | | | ND | |
| 90 Styrene | 104 | | 10.834 | | | | ND | |
| 91 Bromoform | 173 | | 11.011 | | | | ND | |
| 96 1,1,2,2-Tetrachloroethane | 83 | | 11.497 | | | | ND | |
| S 131 Xylenes, Total | 106 | | 1.000 | | | | ND | |

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

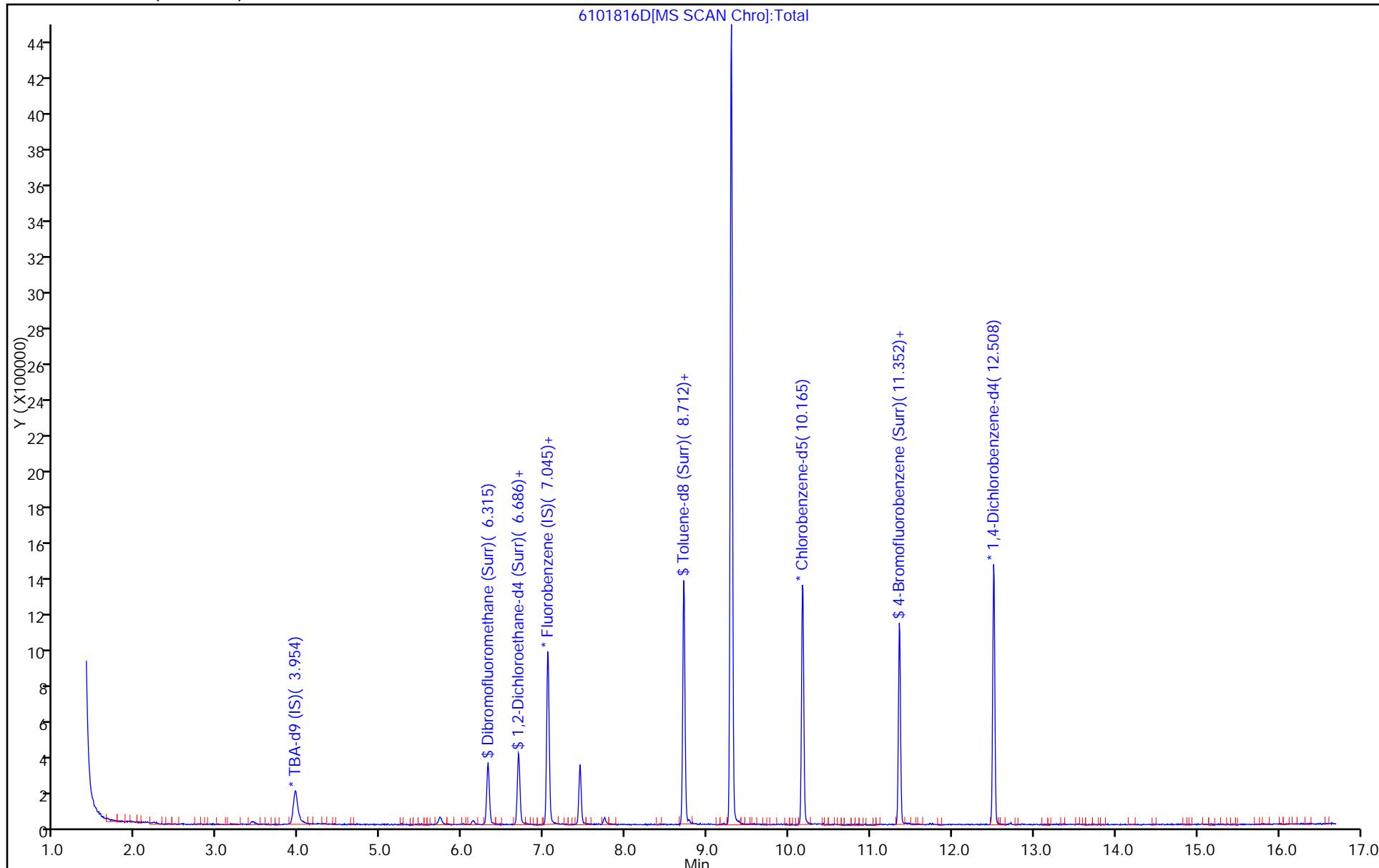
| | | | |
|-------------------|--------------------|-----------|-------------|
| VOA8260INT_00074 | Amount Added: 2.00 | Units: uL | Run Reagent |
| VOA8260SURR_00073 | Amount Added: 2.00 | Units: uL | Run Reagent |

Report Date: 18-Oct-2017 20:31:27

Chrom Revision: 2.2 16-Aug-2017 16:24:46

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101816.D.D
Injection Date: 18-Oct-2017 06:33:30 Instrument ID: CHHP6 Operator ID: 034635
Lims ID: 180-71131-B-1 Lab Sample ID: 180-71131-1 Worklist Smp#: 16
Client ID: HD-SPBA-CW-23-0/1-0
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 25
Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Recovery Report

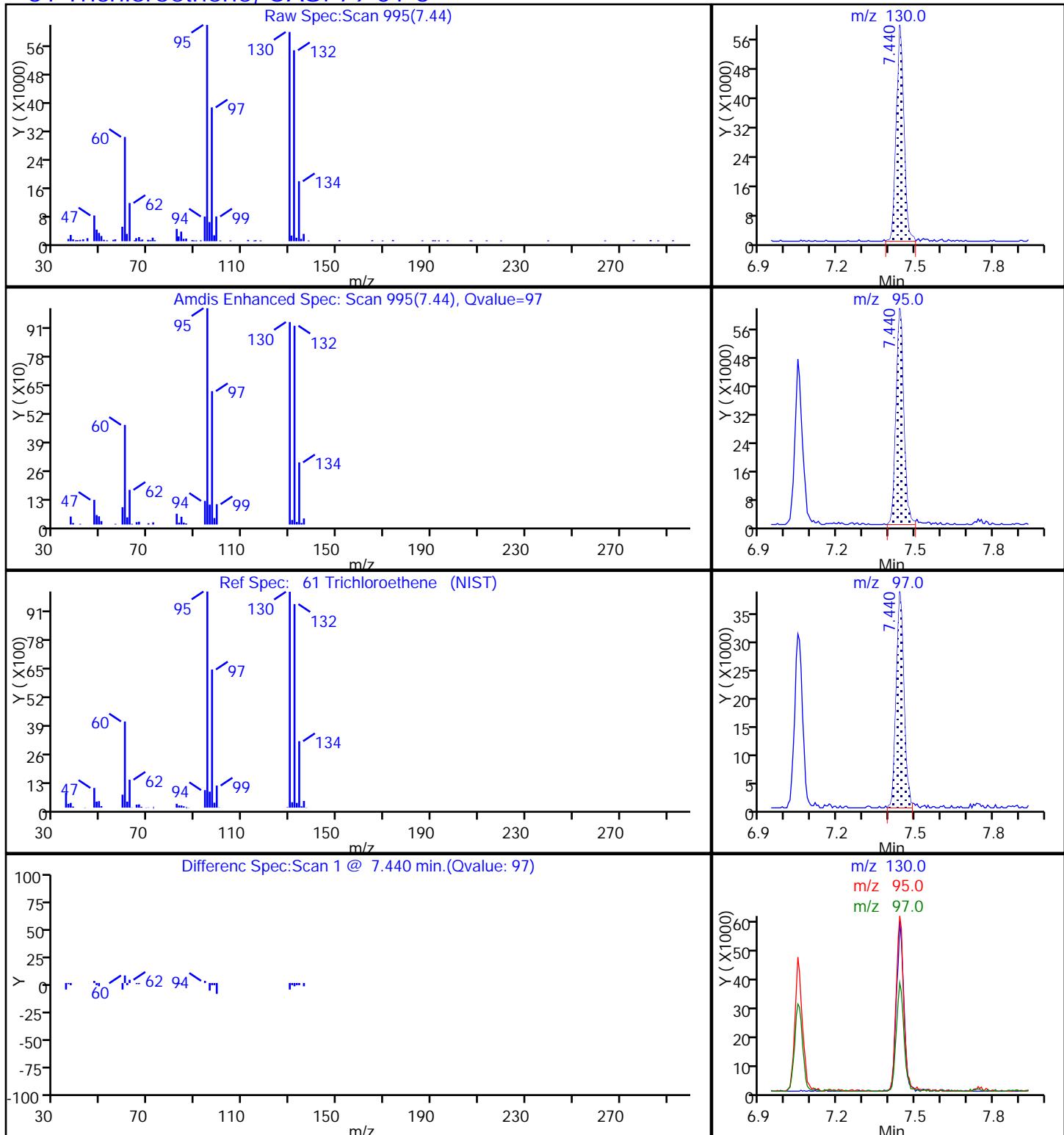
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 Lims ID: 180-71131-B-1
 Client ID: HD-SPBA-CW-23-0/1-0
 Sample Type: Client
 Inject. Date: 18-Oct-2017 06:33:30 ALS Bottle#: 25 Worklist Smp#: 16
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018914-016
 Misc. Info.: 180-71131-B-1
 Operator ID: 034635 Instrument ID: CHHP6
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 18-Oct-2017 20:30:42 Calib Date: 28-Sep-2017 15:13:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICAL File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170928-18631.b\60928P06.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK031

First Level Reviewer: bungardf Date: 18-Oct-2017 20:23:32

| Compound | Amount Added | Amount Recovered | % Rec. |
|-----------------------------------|--------------|------------------|--------|
| \$ 5 Dibromofluoromethane (Surr) | 50.0 | 46.7 | 93.33 |
| \$ 6 1,2-Dichloroethane-d4 (Surr) | 50.0 | 47.1 | 94.23 |
| \$ 7 Toluene-d8 (Surr) | 50.0 | 43.1 | 86.27 |
| \$ 8 4-Bromofluorobenzene (Surr) | 50.0 | 45.3 | 90.59 |

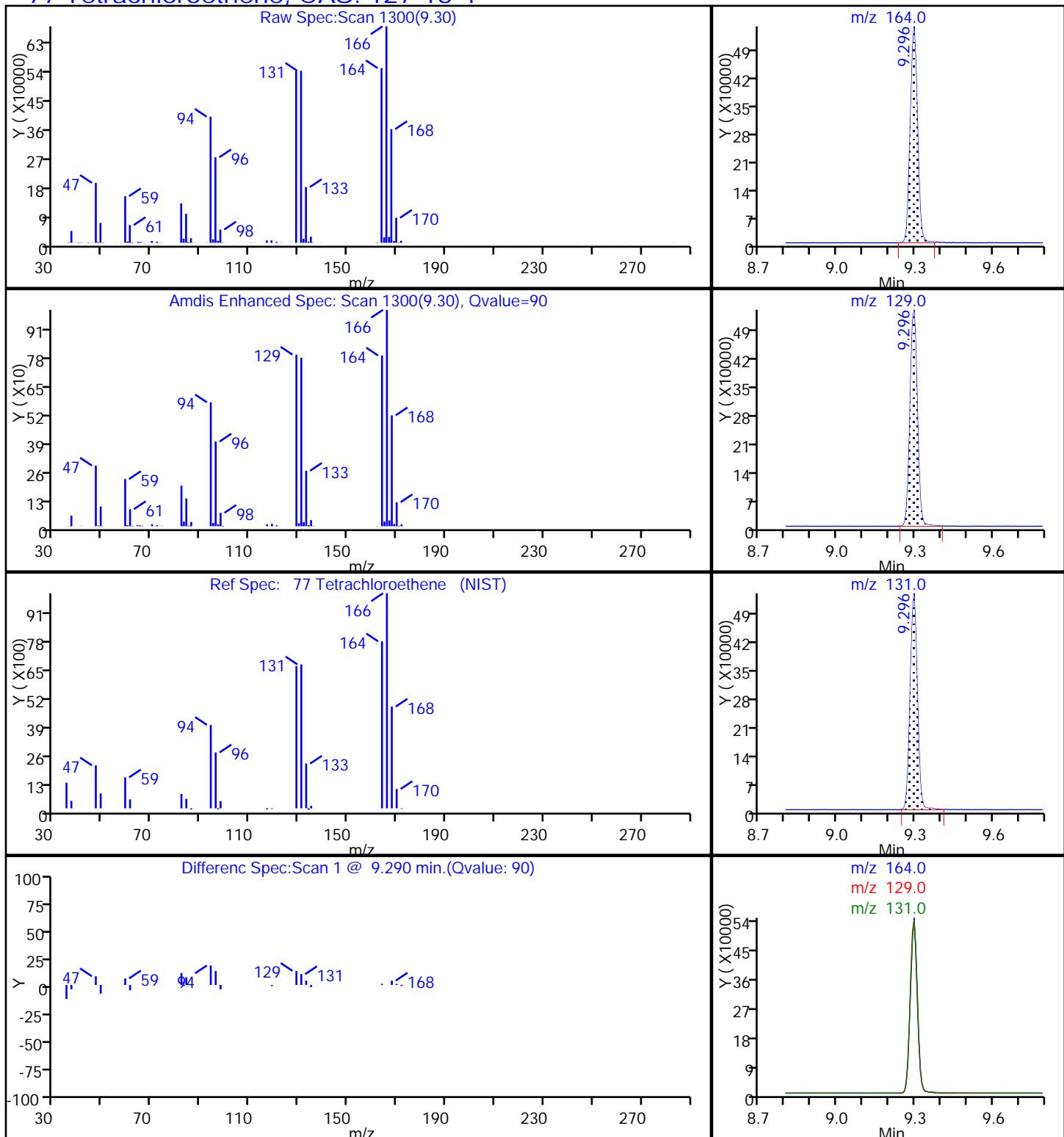
TestAmerica Pittsburgh
 Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101816.D.D
 Injection Date: 18-Oct-2017 06:33:30 Instrument ID: CHHP6
 Lims ID: 180-71131-B-1 Lab Sample ID: 180-71131-1
 Client ID: HD-SPBA-CW-23-0/1-0
 Operator ID: 034635 ALS Bottle#: 25 Worklist Smp#: 16
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
 Column: DB-624 (0.18 mm) Detector: MS SCAN

61 Trichloroethene, CAS: 79-01-6



TestAmerica Pittsburgh
 Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101816.D.D
 Injection Date: 18-Oct-2017 06:33:30 Instrument ID: CHHP6
 Lims ID: 180-71131-B-1 Lab Sample ID: 180-71131-1
 Client ID: HD-SPBA-CW-23-0/1-0
 Operator ID: 034635 ALS Bottle#: 25 Worklist Smp#: 16
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
 Column: DB-624 (0.18 mm) Detector: MS SCAN

77 Tetrachloroethene, CAS: 127-18-4



TestAmerica Pittsburgh

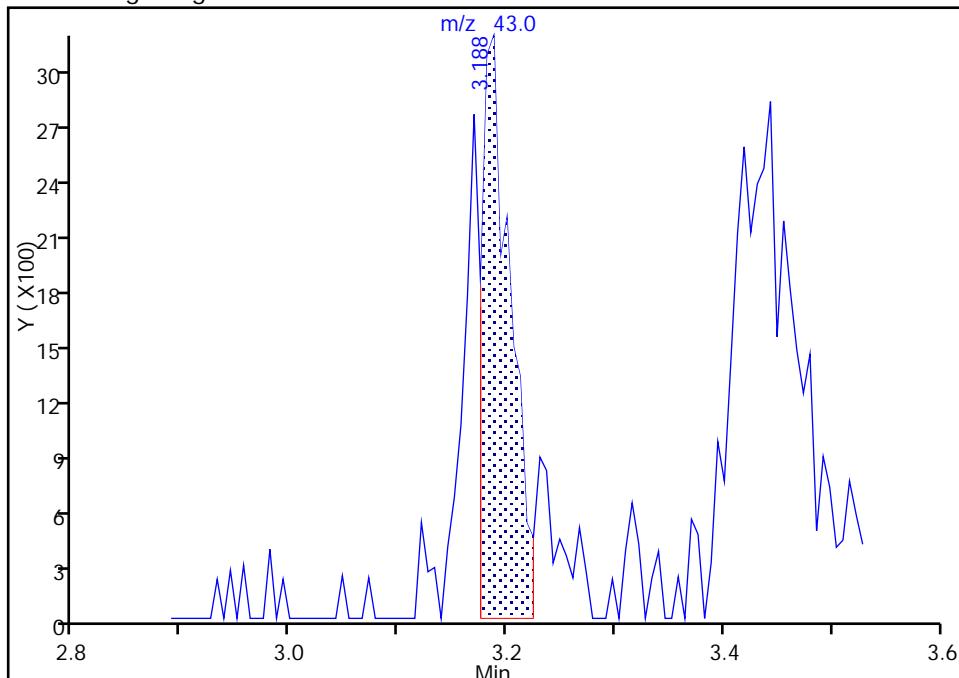
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 Injection Date: 18-Oct-2017 06:33:30 Instrument ID: CHHP6
 Lims ID: 180-71131-B-1 Lab Sample ID: 180-71131-1
 Client ID: HD-SPBA-CW-23-0/1-0
 Operator ID: 034635 ALS Bottle#: 25 Worklist Smp#: 16
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
 Column: DB-624 (0.18 mm) Detector: MS SCAN

24 Acetone, CAS: 67-64-1

Signal: 1

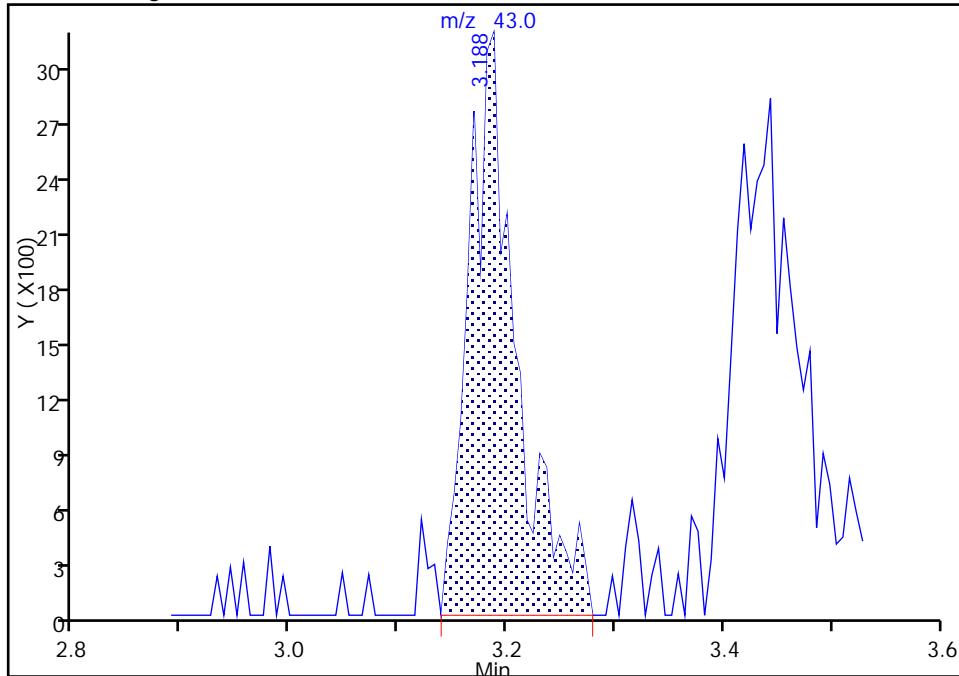
RT: 3.19
 Area: 5710
 Amount: 2.634638
 Amount Units: ng

Processing Integration Results



RT: 3.19
 Area: 9414
 Amount: 4.343692
 Amount Units: ng

Manual Integration Results



Reviewer: bungardf, 18-Oct-2017 20:22:53

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.: _____

Client Sample ID: HD-CW-23-01/1-0

Lab Sample ID: 180-71131-2

Matrix: Water

Lab File ID: 6101817D.D

Analysis Method: 8260C

Date Collected: 10/06/2017 09:50

Sample wt/vol: 5 (mL)

Date Analyzed: 10/18/2017 06:57

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.: 226148

Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|------------|-----------------------------|--------|------|-----|------|
| 74-87-3 | Chloromethane | 1.0 | U | 1.0 | 0.90 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 1.0 | 0.88 |
| 74-83-9 | Bromomethane | 1.0 | U | 1.0 | 0.89 |
| 75-00-3 | Chloroethane | 1.0 | U | 1.0 | 0.90 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.55 |
| 67-64-1 | Acetone | 5.0 | U ^c | 5.0 | 3.4 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 1.0 | 0.88 |
| 75-09-2 | Methylene Chloride | 1.0 | U | 1.0 | 0.36 |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.67 |
| 1634-04-4 | Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.59 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.63 |
| 156-59-2 | cis-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.71 |
| 74-97-5 | Bromochloromethane | 1.0 | U | 1.0 | 0.63 |
| 78-93-3 | 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 |
| 67-66-3 | Chloroform | 1.0 | U | 1.0 | 0.60 |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.60 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U ^c | 1.0 | 0.88 |
| 71-43-2 | Benzene | 1.0 | U | 1.0 | 0.60 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.57 |
| 79-01-6 | Trichloroethene | 4.4 | | 1.0 | 0.69 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.66 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 1.0 | 0.64 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.59 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 5.0 | U ^c | 5.0 | 3.1 |
| 108-88-3 | Toluene | 1.0 | U | 1.0 | 0.46 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.58 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.45 |
| 127-18-4 | Tetrachloroethene | 43 | | 1.0 | 0.47 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 5.0 | 3.3 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 1.0 | 0.84 |
| 106-93-4 | 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.50 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 1.0 | 0.50 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.57 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 1.0 | 0.51 |
| 1330-20-7 | Xylenes, Total | 2.0 | U | 2.0 | 0.89 |
| 100-42-5 | Styrene | 1.0 | U | 1.0 | 0.47 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-71131-1
SDG No.: _____
Client Sample ID: HD-CW-23-0/1-0 Lab Sample ID: 180-71131-2
Matrix: Water Lab File ID: 6101817D.D
Analysis Method: 8260C Date Collected: 10/06/2017 09:50
Sample wt/vol: 5 (mL) Date Analyzed: 10/18/2017 06:57
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.: 226148 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|----------|---------------------------|--------|------|-----|------|
| 75-25-2 | Bromoform | 1.0 | U | 1.0 | 0.98 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.60 |
| 107-13-1 | Acrylonitrile | 20 | U ^c | 20 | 7.8 |
| 123-91-1 | 1,4-Dioxane | 200 | U | 200 | 14 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|------------|------------------------------|------|---|--------|
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 94 | | 65-121 |
| 2037-26-5 | Toluene-d8 (Surr) | 89 | | 73-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 91 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 93 | | 73-120 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101817D.D
 Lims ID: 180-71131-A-2
 Client ID: HD-CW-23-0/1-0
 Sample Type: Client
 Inject. Date: 18-Oct-2017 06:57:30 ALS Bottle#: 26 Worklist Smp#: 17
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018914-017
 Misc. Info.: 180-71131-A-2
 Operator ID: 034635 Instrument ID: CHHP6
 Method: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 18-Oct-2017 20:30:42 Calib Date: 28-Sep-2017 15:13:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICAL File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170928-18631.b\\60928P06.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK031

First Level Reviewer: bungardf Date: 18-Oct-2017 20:24:26

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 3.951 | 3.949 | 0.002 | 91 | 286138 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.048 | 7.045 | 0.003 | 98 | 994351 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.169 | 10.166 | 0.003 | 88 | 272894 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.511 | 12.508 | 0.003 | 97 | 392535 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr) | 113 | 6.318 | 6.314 | 0.004 | 91 | 241266 | 46.7 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur) | 65 | 6.689 | 6.685 | 0.004 | 70 | 347514 | 47.1 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.715 | 8.711 | 0.004 | 93 | 982090 | 44.6 | |
| \$ 8 4-Bromofluorobenzene (Surr) | 95 | 11.355 | 11.351 | 0.004 | 82 | 430032 | 45.6 | |
| 12 Chloromethane | 50 | | 1.624 | | | | ND | |
| 13 Vinyl chloride | 62 | | 1.740 | | | | ND | |
| 15 Bromomethane | 94 | | 2.068 | | | | ND | |
| 16 Chloroethane | 64 | | 2.202 | | | | ND | |
| 22 1,1-Dichloroethene | 96 | | 3.090 | | | | ND | |
| 24 Acetone | 43 | 3.179 | 3.181 | -0.002 | 90 | 9919 | 4.65 | |
| 26 Carbon disulfide | 76 | | 3.364 | | | | ND | |
| 31 Methylene Chloride | 84 | | 3.838 | | | | ND | |
| 33 Acrylonitrile | 53 | | 4.222 | | | | ND | |
| 34 trans-1,2-Dichloroethene | 96 | | 4.270 | | | | ND | |
| 35 Methyl tert-butyl ether | 73 | | 4.276 | | | | ND | |
| 37 1,1-Dichloroethane | 63 | | 4.921 | | | | ND | |
| 43 cis-1,2-Dichloroethene | 96 | 5.691 | 5.688 | 0.003 | 28 | 3038 | 0.4435 | |
| 44 2-Butanone (MEK) | 43 | | 5.694 | | | | ND | |
| 48 Chlorobromomethane | 128 | | 5.974 | | | | ND | |
| 50 Chloroform | 83 | 6.135 | 6.126 | 0.009 | 92 | 20285 | 1.97 | |
| 51 1,1,1-Trichloroethane | 97 | | 6.290 | | | | ND | |
| 53 Carbon tetrachloride | 117 | | 6.454 | | | | ND | |
| 56 Benzene | 78 | | 6.691 | | | | ND | |
| 57 1,2-Dichloroethane | 62 | | 6.777 | | | | ND | |
| 61 Trichloroethene | 130 | 7.437 | 7.440 | -0.003 | 98 | 122690 | 22.0 | |
| 64 1,2-Dichloropropane | 63 | | 7.713 | | | | ND | |
| 65 1,4-Dioxane | 88 | | 7.793 | | | | ND | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | OnCol Amt ng | Flags |
|--------------------------------|-----|--------------|------------------|------------------|----|----------|-----------------|-------|
| 68 Dichlorobromomethane | 83 | | 7.999 | | | | ND | |
| 71 cis-1,3-Dichloropropene | 75 | | 8.450 | | | | ND | |
| 72 4-Methyl-2-pentanone (MIBK) | 43 | | 8.608 | | | | ND | |
| 73 Toluene | 91 | 8.782 | 8.778 | 0.004 | 92 | 8805 | 0.3274 | |
| 74 trans-1,3-Dichloropropene | 75 | | 9.034 | | | | ND | |
| 76 1,1,2-Trichloroethane | 97 | | 9.222 | | | | ND | |
| 77 Tetrachloroethene | 164 | 9.293 | 9.295 | -0.002 | 90 | 1025103 | 212.9 | |
| 79 2-Hexanone | 43 | | 9.447 | | | | ND | |
| 81 Chlorodibromomethane | 129 | | 9.593 | | | | ND | |
| 82 Ethylene Dibromide | 107 | | 9.709 | | | | ND | |
| 84 Chlorobenzene | 112 | | 10.196 | | | | ND | |
| 86 1,1,1,2-Tetrachloroethane | 131 | | 10.293 | | | | ND | |
| 87 Ethylbenzene | 106 | | 10.299 | | | | ND | |
| 88 m-Xylene & p-Xylene | 106 | | 10.433 | | | | ND | |
| 89 o-Xylene | 106 | | 10.816 | | | | ND | |
| 90 Styrene | 104 | | 10.834 | | | | ND | |
| 91 Bromoform | 173 | | 11.011 | | | | ND | |
| 96 1,1,2,2-Tetrachloroethane | 83 | | 11.497 | | | | ND | |
| S 131 Xylenes, Total | 106 | | 1.000 | | | | ND | |

Reagents:

VOA8260INT_00074
 VOA8260SURR_00073

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

Report Date: 18-Oct-2017 20:31:29

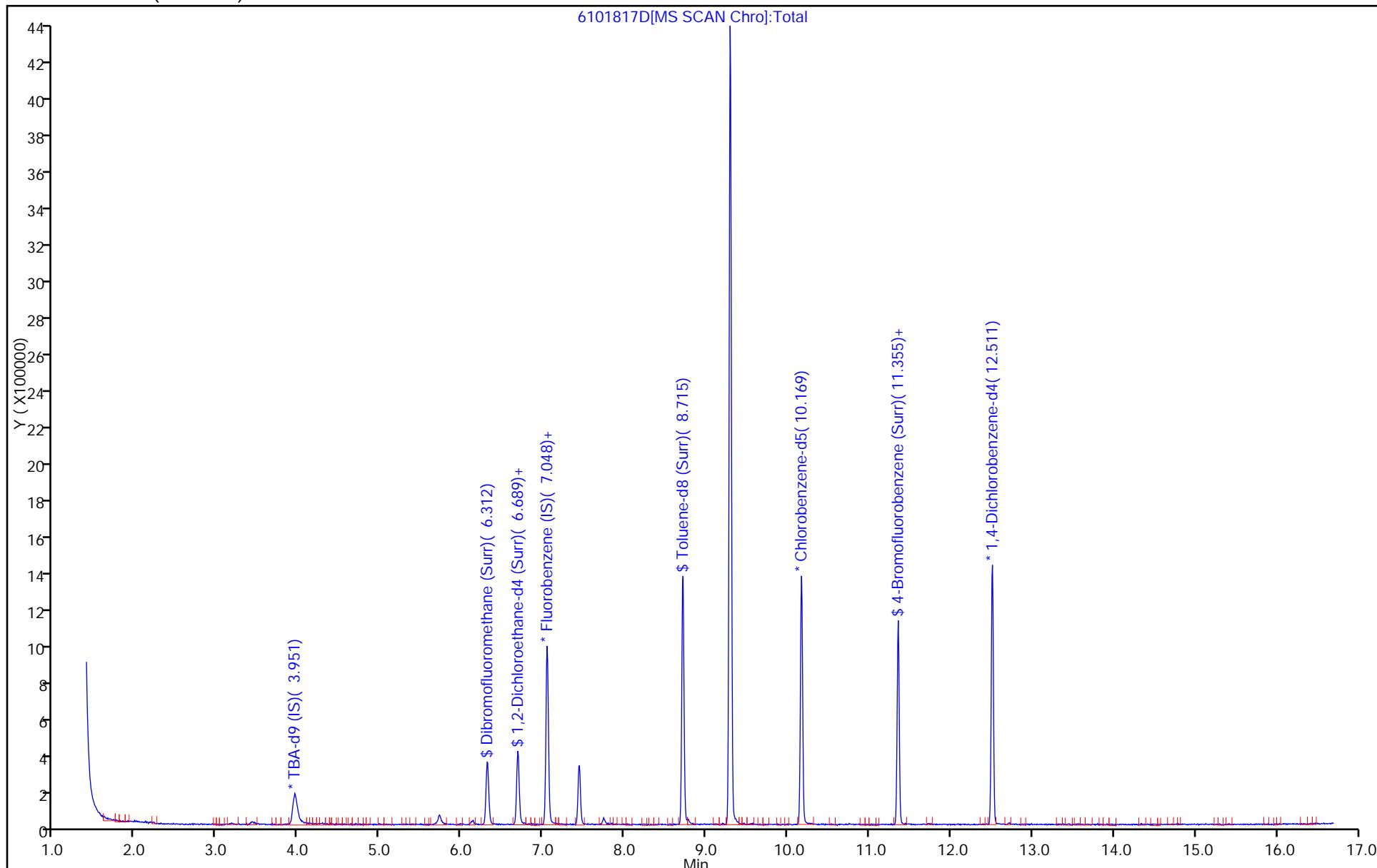
Chrom Revision: 2.2 16-Aug-2017 16:24:46

TestAmerica Pittsburgh

| | | | |
|-----------------|---|----------------|----------------|
| Data File: | \ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\6101817D.D | | |
| Injection Date: | 18-Oct-2017 06:57:30 | Instrument ID: | CHHP6 |
| Lims ID: | 180-71131-A-2 | Lab Sample ID: | 180-71131-2 |
| Client ID: | HD-CW-23-01-0 | Dil. Factor: | 1.0000 |
| Purge Vol: | 5.000 mL | Limit Group: | VOA 8260C ICAL |
| Method: | MSVOA_LL_CHHP6 | | |
| Column: | DB-624 (0.18 mm) | | |

Operator ID: 034635
Worklist Smp#: 17

ALS Bottle#: 26



TestAmerica Pittsburgh
Recovery Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\6101817D.D
 Lims ID: 180-71131-A-2
 Client ID: HD-CW-23-0/1-0
 Sample Type: Client
 Inject. Date: 18-Oct-2017 06:57:30 ALS Bottle#: 26 Worklist Smp#: 17
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018914-017
 Misc. Info.: 180-71131-A-2
 Operator ID: 034635 Instrument ID: CHHP6
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 18-Oct-2017 20:30:42 Calib Date: 28-Sep-2017 15:13:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170928-18631.b\60928P06.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK031

First Level Reviewer: bungardf Date: 18-Oct-2017 20:24:26

| Compound | Amount Added | Amount Recovered | % Rec. |
|-----------------------------------|--------------|------------------|--------|
| \$ 5 Dibromofluoromethane (Surr) | 50.0 | 46.7 | 93.38 |
| \$ 6 1,2-Dichloroethane-d4 (Surr) | 50.0 | 47.1 | 94.11 |
| \$ 7 Toluene-d8 (Surr) | 50.0 | 44.6 | 89.11 |
| \$ 8 4-Bromofluorobenzene (Surr) | 50.0 | 45.6 | 91.28 |

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101817D.D

Injection Date: 18-Oct-2017 06:57:30

Instrument ID: CHHP6

Lims ID: 180-71131-A-2

Lab Sample ID: 180-71131-2

Client ID: HD-CW-23-0/1-0

Operator ID: 034635

ALS Bottle#: 26 Worklist Smp#: 17

Purge Vol: 5.000 mL

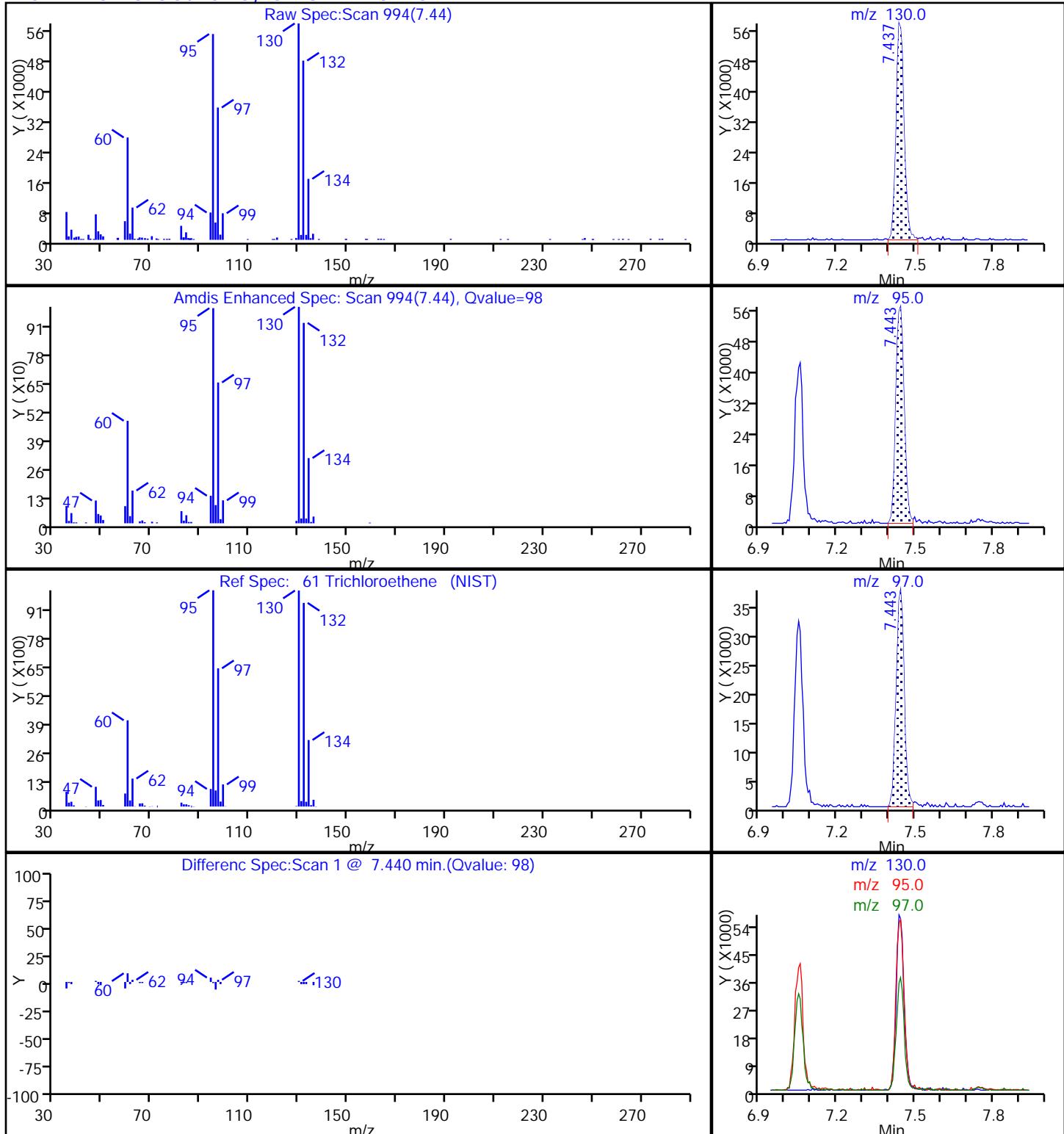
Dil. Factor: 1.0000

Method: MSVOA_LL_CHHP6

Limit Group: VOA 8260C ICAL

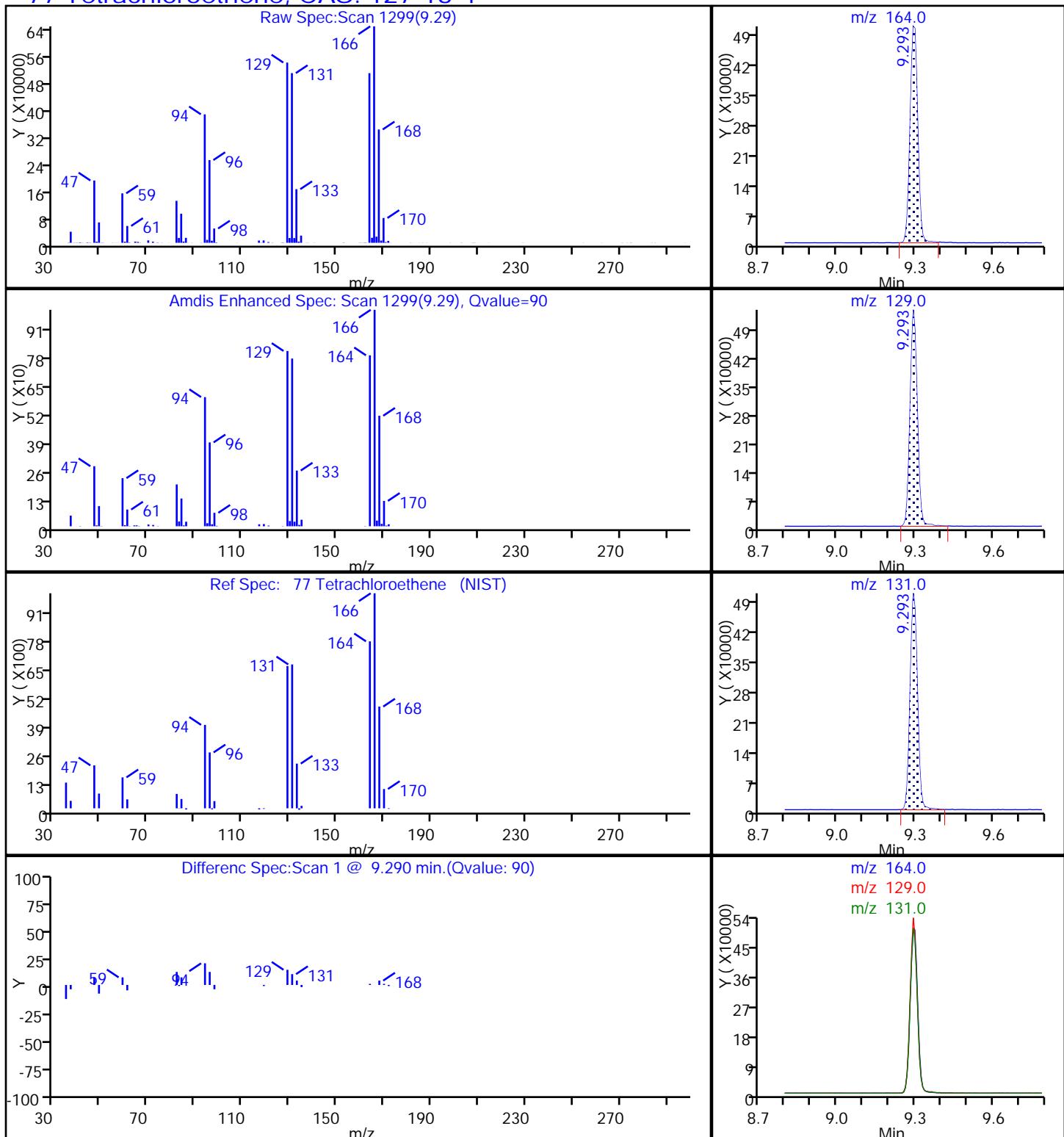
Column: DB-624 (0.18 mm)

Detector: MS SCAN

61 Trichloroethene, CAS: 79-01-6

TestAmerica Pittsburgh
 Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101817D.D
 Injection Date: 18-Oct-2017 06:57:30 Instrument ID: CHHP6
 Lims ID: 180-71131-A-2 Lab Sample ID: 180-71131-2
 Client ID: HD-CW-23-0/1-0
 Operator ID: 034635 ALS Bottle#: 26 Worklist Smp#: 17
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
 Column: DB-624 (0.18 mm) Detector: MS SCAN

77 Tetrachloroethene, CAS: 127-18-4



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.: _____

Client Sample ID: HD-QC6-0/1-2

Lab Sample ID: 180-71131-3

Matrix: Water

Lab File ID: 6101819D.D

Analysis Method: 8260C

Date Collected: 10/05/2017 12:00

Sample wt/vol: 5 (mL)

Date Analyzed: 10/18/2017 07:45

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.: 226148

Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|------------|-----------------------------|--------|------|-----|------|
| 74-87-3 | Chloromethane | 1.0 | U | 1.0 | 0.90 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 1.0 | 0.88 |
| 74-83-9 | Bromomethane | 1.0 | U | 1.0 | 0.89 |
| 75-00-3 | Chloroethane | 1.0 | U | 1.0 | 0.90 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.55 |
| 67-64-1 | Acetone | 5.0 | U ^c | 5.0 | 3.4 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 1.0 | 0.88 |
| 75-09-2 | Methylene Chloride | 1.0 | | 1.0 | 0.36 |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.67 |
| 1634-04-4 | Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.59 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.63 |
| 156-59-2 | cis-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.71 |
| 74-97-5 | Bromochloromethane | 1.0 | U | 1.0 | 0.63 |
| 78-93-3 | 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 |
| 67-66-3 | Chloroform | 1.0 | U | 1.0 | 0.60 |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.60 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U ^c | 1.0 | 0.88 |
| 71-43-2 | Benzene | 1.0 | U | 1.0 | 0.60 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.57 |
| 79-01-6 | Trichloroethene | 1.0 | U | 1.0 | 0.69 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.66 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 1.0 | 0.64 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.59 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 5.0 | U ^c | 5.0 | 3.1 |
| 108-88-3 | Toluene | 1.0 | U | 1.0 | 0.46 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.58 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.45 |
| 127-18-4 | Tetrachloroethene | 1.0 | U | 1.0 | 0.47 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 5.0 | 3.3 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 1.0 | 0.84 |
| 106-93-4 | 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.50 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 1.0 | 0.50 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.57 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 1.0 | 0.51 |
| 1330-20-7 | Xylenes, Total | 2.0 | U | 2.0 | 0.89 |
| 100-42-5 | Styrene | 1.0 | U | 1.0 | 0.47 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-71131-1
SDG No.:
Client Sample ID: HD-QC6-0/1-2 Lab Sample ID: 180-71131-3
Matrix: Water Lab File ID: 6101819D.D
Analysis Method: 8260C Date Collected: 10/05/2017 12:00
Sample wt/vol: 5 (mL) Date Analyzed: 10/18/2017 07:45
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.: 226148 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|----------|---------------------------|--------|------|-----|------|
| 75-25-2 | Bromoform | 1.0 | U | 1.0 | 0.98 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.60 |
| 107-13-1 | Acrylonitrile | 20 | U ^c | 20 | 7.8 |
| 123-91-1 | 1,4-Dioxane | 200 | U | 200 | 14 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|------------|------------------------------|------|---|--------|
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 95 | | 65-121 |
| 2037-26-5 | Toluene-d8 (Surr) | 87 | | 73-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 89 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 95 | | 73-120 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101819D.D
 Lims ID: 180-71131-B-3
 Client ID: HD-QC6-0/1-2
 Sample Type: Client
 Inject. Date: 18-Oct-2017 07:45:30 ALS Bottle#: 28 Worklist Smp#: 19
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018914-019
 Misc. Info.: 180-71131-B-3
 Operator ID: 034635 Instrument ID: CHHP6
 Method: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 18-Oct-2017 20:30:42 Calib Date: 28-Sep-2017 15:13:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICAL File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170928-18631.b\\60928P06.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK031

First Level Reviewer: bungardf Date: 18-Oct-2017 20:25:23

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 3.955 | 3.949 | 0.006 | 89 | 274765 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.045 | 7.045 | 0.000 | 98 | 988524 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.166 | 10.166 | 0.000 | 88 | 275892 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.508 | 12.508 | 0.000 | 97 | 388863 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr) | 113 | 6.315 | 6.314 | 0.001 | 92 | 244625 | 47.6 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur) | 65 | 6.687 | 6.685 | 0.002 | 70 | 347543 | 47.3 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.712 | 8.711 | 0.001 | 93 | 966982 | 43.3 | |
| \$ 8 4-Bromofluorobenzene (Surr) | 95 | 11.353 | 11.351 | 0.002 | 81 | 425709 | 44.7 | |
| 12 Chloromethane | 50 | 1.624 | | | | | ND | |
| 13 Vinyl chloride | 62 | 1.740 | | | | | ND | |
| 15 Bromomethane | 94 | 2.068 | | | | | ND | |
| 16 Chloroethane | 64 | 2.202 | | | | | ND | |
| 22 1,1-Dichloroethene | 96 | 3.090 | | | | | ND | |
| 24 Acetone | 43 | 3.176 | 3.181 | -0.005 | 92 | 18530 | 8.74 | |
| 26 Carbon disulfide | 76 | 3.364 | | | | | ND | |
| 31 Methylene Chloride | 84 | 3.839 | 3.838 | 0.001 | 81 | 34183 | 5.02 | |
| 33 Acrylonitrile | 53 | 4.222 | | | | | ND | |
| 34 trans-1,2-Dichloroethene | 96 | 4.270 | | | | | ND | |
| 35 Methyl tert-butyl ether | 73 | 4.276 | | | | | ND | |
| 37 1,1-Dichloroethane | 63 | 4.921 | | | | | ND | |
| 43 cis-1,2-Dichloroethene | 96 | 5.688 | | | | | ND | |
| 44 2-Butanone (MEK) | 43 | 5.694 | | | | | ND | |
| 48 Chlorobromomethane | 128 | 5.974 | | | | | ND | |
| 50 Chloroform | 83 | 6.126 | | | | | ND | |
| 51 1,1,1-Trichloroethane | 97 | 6.290 | | | | | ND | |
| 53 Carbon tetrachloride | 117 | 6.454 | | | | | ND | |
| 56 Benzene | 78 | 6.691 | | | | | ND | |
| 57 1,2-Dichloroethane | 62 | 6.777 | | | | | ND | |
| 61 Trichloroethene | 130 | 7.440 | | | | | ND | |
| 64 1,2-Dichloropropane | 63 | 7.713 | | | | | ND | |
| 65 1,4-Dioxane | 88 | 7.793 | | | | | ND | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | OnCol Amt ng | Flags |
|--------------------------------|-----|--------------|------------------|------------------|---|----------|-----------------|-------|
| 68 Dichlorobromomethane | 83 | | 7.999 | | | | ND | |
| 71 cis-1,3-Dichloropropene | 75 | | 8.450 | | | | ND | |
| 72 4-Methyl-2-pentanone (MIBK) | 43 | | 8.608 | | | | ND | |
| 73 Toluene | 91 | | 8.778 | | | | ND | |
| 74 trans-1,3-Dichloropropene | 75 | | 9.034 | | | | ND | |
| 76 1,1,2-Trichloroethane | 97 | | 9.222 | | | | ND | |
| 77 Tetrachloroethene | 164 | | 9.295 | | | | ND | |
| 79 2-Hexanone | 43 | | 9.447 | | | | ND | |
| 81 Chlorodibromomethane | 129 | | 9.593 | | | | ND | |
| 82 Ethylene Dibromide | 107 | | 9.709 | | | | ND | |
| 84 Chlorobenzene | 112 | | 10.196 | | | | ND | |
| 86 1,1,1,2-Tetrachloroethane | 131 | | 10.293 | | | | ND | |
| 87 Ethylbenzene | 106 | | 10.299 | | | | ND | |
| 88 m-Xylene & p-Xylene | 106 | | 10.433 | | | | ND | |
| 89 o-Xylene | 106 | | 10.816 | | | | ND | |
| 90 Styrene | 104 | | 10.834 | | | | ND | |
| 91 Bromoform | 173 | | 11.011 | | | | ND | |
| 96 1,1,2,2-Tetrachloroethane | 83 | | 11.497 | | | | ND | |
| S 131 Xylenes, Total | 106 | | 1.000 | | | | ND | |

Reagents:

VOA8260INT_00074
 VOA8260SURR_00073

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

Report Date: 18-Oct-2017 20:31:31

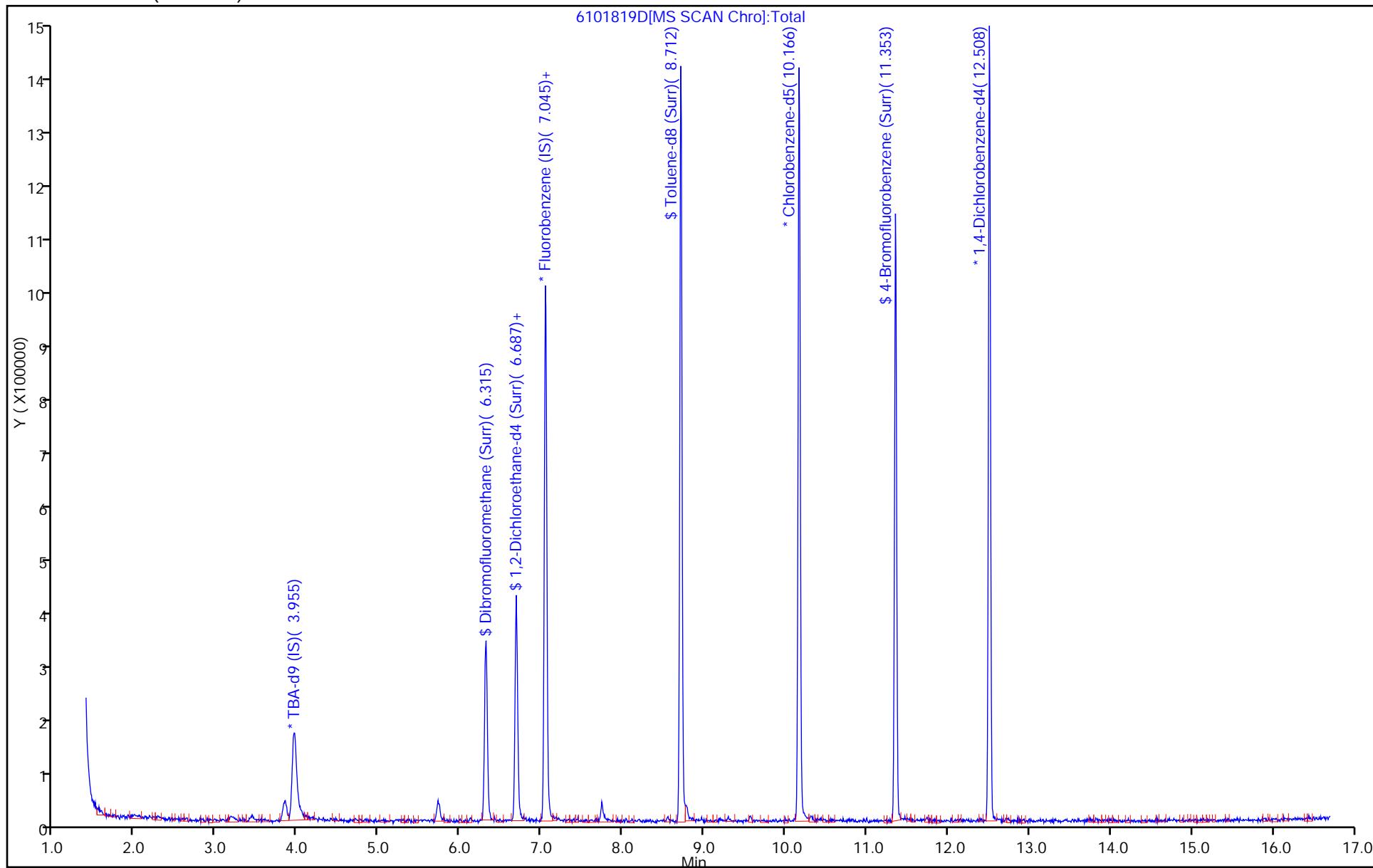
Chrom Revision: 2.2 16-Aug-2017 16:24:46

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101819D.D
Injection Date: 18-Oct-2017 07:45:30 Instrument ID: CHHP6
Lims ID: 180-71131-B-3 Lab Sample ID: 180-71131-3
Client ID: HD-QC6-01-2
Purge Vol: 5.000 mL Dil. Factor: 1.0000
Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm)

Operator ID: 034635
Worklist Smp#: 19

ALS Bottle#: 28



TestAmerica Pittsburgh
Recovery Report

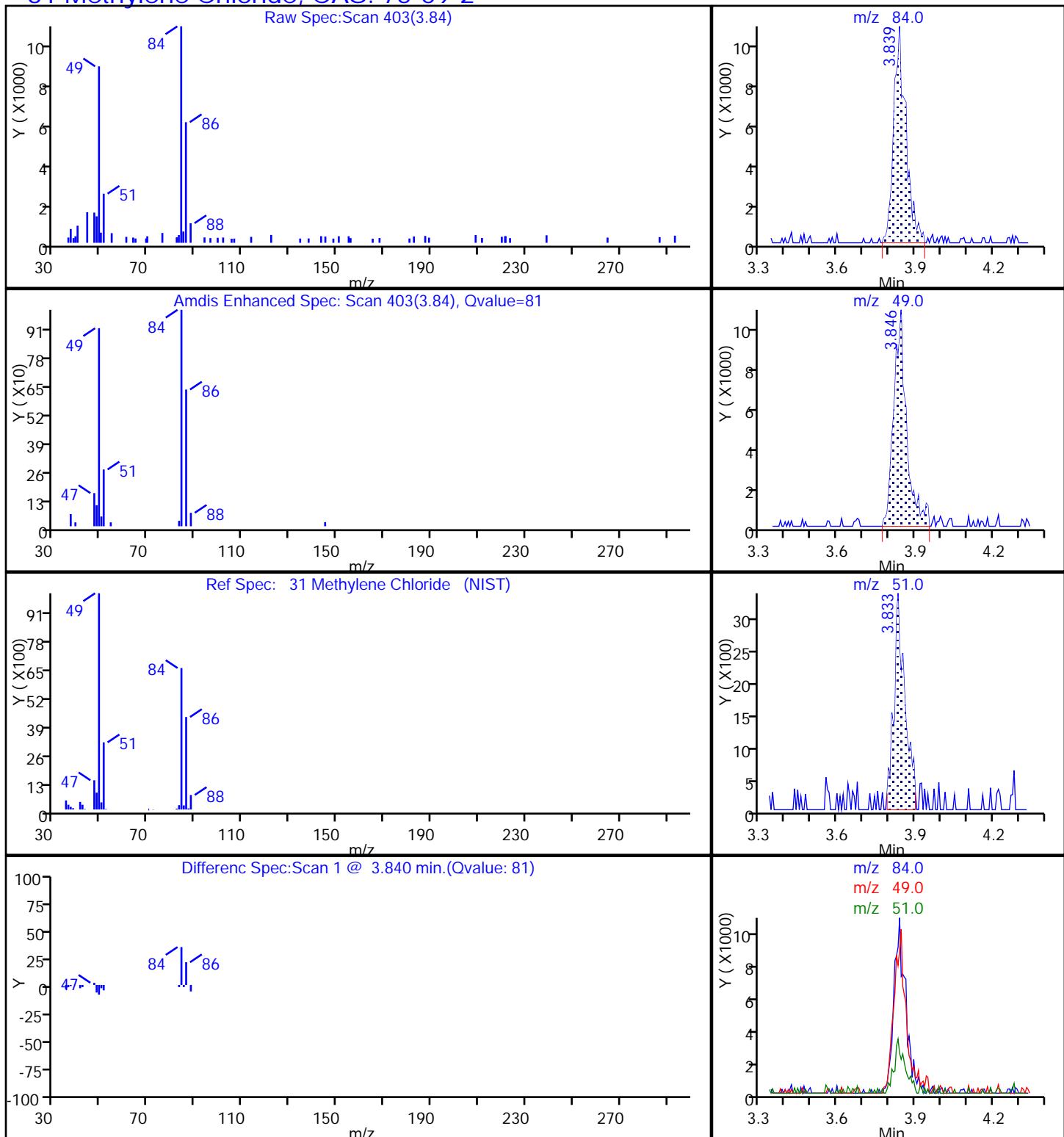
Data File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\6101819D.D
 Lims ID: 180-71131-B-3
 Client ID: HD-QC6-01/1-2
 Sample Type: Client
 Inject. Date: 18-Oct-2017 07:45:30 ALS Bottle#: 28 Worklist Smp#: 19
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018914-019
 Misc. Info.: 180-71131-B-3
 Operator ID: 034635 Instrument ID: CHHP6
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 18-Oct-2017 20:30:42 Calib Date: 28-Sep-2017 15:13:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170928-18631.b\60928P06.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK031

First Level Reviewer: bungardf Date: 18-Oct-2017 20:25:23

| Compound | Amount Added | Amount Recovered | % Rec. |
|-----------------------------------|--------------|------------------|--------|
| \$ 5 Dibromofluoromethane (Surr) | 50.0 | 47.6 | 95.24 |
| \$ 6 1,2-Dichloroethane-d4 (Surr) | 50.0 | 47.3 | 94.68 |
| \$ 7 Toluene-d8 (Surr) | 50.0 | 43.3 | 86.63 |
| \$ 8 4-Bromofluorobenzene (Surr) | 50.0 | 44.7 | 89.38 |

TestAmerica Pittsburgh
 Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101819D.D
 Injection Date: 18-Oct-2017 07:45:30 Instrument ID: CHHP6
 Lims ID: 180-71131-B-3 Lab Sample ID: 180-71131-3
 Client ID: HD-QC6-0/1-2
 Operator ID: 034635 ALS Bottle#: 28 Worklist Smp#: 19
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 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 BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Pittsburgh Job No.: 180-71131-1 Analy Batch No.: 217861

SDG No.: _____

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---------|-------------------|--------------|
| Level 1 | IC 180-217861/3 | 60724D03.D |
| Level 2 | IC 180-217861/4 | 60724D04.D |
| Level 3 | ICIS 180-217861/5 | 60724D05.D |
| Level 4 | IC 180-217861/6 | 60724D06.D |
| Level 5 | IC 180-217861/7 | 60724D07.D |
| Level 6 | IC 180-217861/8 | 60724D08.D |
| Level 7 | IC 180-217861/9 | 60724D09.D |
| Level 8 | IC 180-217861/10 | 60724D10.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 | 0.3450 0.3059 | 0.3450 0.2869 | 0.3353 0.3121 | 0.3123 | 0.3101 | Ave | | 0.3191 | | | 0.1000 | 6.5 | | 20.0 | | | |
| Chloromethane | 0.3003 0.2886 | 0.3121 0.2599 | 0.2915 0.2724 | 0.2904 | 0.2802 | Ave | | 0.2869 | | | 0.1000 | 5.6 | | 20.0 | | | |
| Vinyl chloride | 0.3294 0.2874 | 0.3184 0.2708 | 0.3073 0.2897 | 0.3106 | 0.2927 | Ave | | 0.3008 | | | 0.1000 | 6.3 | | 20.0 | | | |
| 1,3-Butadiene | 0.3092 0.2283 | 0.2618 0.2122 | 0.2620 0.2355 | 0.2508 | 0.2350 | Ave | | 0.2494 | | | 0.0100 | 11.9 | | 20.0 | | | |
| Bromomethane | 0.1553 0.1421 | 0.1665 0.1212 | 0.1437 0.1124 | 0.1420 | 0.1381 | Ave | | 0.1402 | | | 0.0500 | 12.3 | | 20.0 | | | |
| Chloroethane | 0.1804 0.1621 | 0.1823 0.1426 | 0.1613 0.1385 | 0.1692 | 0.1635 | Ave | | 0.1625 | | | 0.0500 | 9.7 | | 20.0 | | | |
| Trichlorofluoromethane | 0.3064 0.2991 | 0.3069 0.2753 | 0.2862 0.2942 | 0.3037 | 0.2975 | Ave | | 0.2961 | | | 0.1000 | 3.7 | | 20.0 | | | |
| Ethyl ether | 0.2704 0.2624 | 0.2948 0.2348 | 0.2431 0.2337 | 0.2503 | 0.2380 | Ave | | 0.2534 | | | 0.0100 | 8.4 | | 20.0 | | | |
| Acrolein | 0.0585 0.0578 | 0.0586 0.0511 | 0.0493 0.0540 | 0.0552 | 0.0498 | Ave | | 0.0543 | | | 0.0100 | 7.2 | | 20.0 | | | |
| 1,1-Dichloroethene | 0.2867 0.2623 | 0.2569 0.2445 | 0.2480 0.2685 | 0.2599 | 0.2523 | Ave | | 0.2599 | | | 0.1000 | 5.1 | | 20.0 | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.2563 0.2479 | 0.2553 0.2297 | 0.2505 0.2557 | 0.2551 | 0.2415 | Ave | | 0.2490 | | | 0.1000 | 3.7 | | 20.0 | | | |
| Acetone | 0.1203 0.1200 | 0.1188 0.0908 | 0.1059 0.0850 | 0.1178 | 0.0996 | Ave | | 0.1073 | | | 0.0500 | 13.2 | | 20.0 | | | |
| Iodomethane | 0.3702 0.3809 | 0.3779 0.3532 | 0.3472 0.3674 | 0.3706 | 0.3553 | Ave | | 0.3654 | | | 0.0100 | 3.3 | | 20.0 | | | |
| Carbon disulfide | 0.5160 0.6411 | 0.5084 0.6051 | 0.5286 0.6607 | 0.5745 | 0.5948 | Ave | | 0.5787 | | | 0.1000 | 9.9 | | 20.0 | | | |

Note: The M1 coefficient is the same as Ave RRF for an Ave curve type.

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

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

Analy Batch No.: 217861

SDG No.: _____

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

| 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 | | | | | | | | |
| Allyl chloride | 0.1400 0.1667 | 0.1427 0.1552 | 0.1408 0.1669 | 0.1545 | 0.1506 | Ave | | 0.1522 | | | 0.0100 | 7.1 | | 20.0 | | | |
| Methyl acetate | 0.2511 0.2511 | 0.2629 0.2169 | 0.2237 0.2238 | 0.2281 | 0.2235 | Ave | | 0.2351 | | | 0.1000 | 7.3 | | 20.0 | | | |
| Methylene Chloride | 0.4304 0.3426 | 0.3720 0.3106 | 0.3253 0.3196 | 0.3333 | 0.3214 | Ave | | 0.3444 | | | 0.1000 | 11.5 | | 20.0 | | | |
| tert-Butyl alcohol | 1.1832 1.1169 | 1.0643 1.1780 | 1.0783 1.0503 | 1.0936 | 1.0903 | Ave | | 1.1069 | | | 0.0100 | 4.5 | | 20.0 | | | |
| Acrylonitrile | 0.1433 0.1282 | 0.1504 0.1099 | 0.1193 0.1099 | 0.1260 | 0.1177 | Ave | | 0.1256 | | | 0.0100 | 11.8 | | 20.0 | | | |
| trans-1,2-Dichloroethene | 0.3079 0.2977 | 0.3084 0.2782 | 0.2862 0.2973 | 0.2967 | 0.2863 | Ave | | 0.2948 | | | 0.1000 | 3.6 | | 20.0 | | | |
| Methyl tert-butyl ether | 0.9852 0.9798 | 1.0649 0.8774 | 0.9152 0.8866 | 0.9403 | 0.8941 | Ave | | 0.9429 | | | 0.1000 | 6.8 | | 20.0 | | | |
| Hexane | 0.3830 0.3427 | 0.3628 0.3171 | 0.3558 0.3531 | 0.3486 | 0.3306 | Ave | | 0.3492 | | | 0.0100 | 5.7 | | 20.0 | | | |
| 1,1-Dichloroethane | 0.3424 0.5206 | 0.5334 0.4688 | 0.4893 0.4858 | 0.5109 | 0.4862 | Ave | | 0.4797 | | | 0.2000 | 12.4 | | 20.0 | | | |
| Vinyl acetate | 0.5164 0.5873 | 0.6180 0.5638 | 0.5095 0.5801 | 0.5598 | 0.5341 | Ave | | 0.5586 | | | 0.0100 | 6.6 | | 20.0 | | | |
| 2,2-Dichloropropane | 0.0453 0.0537 | 0.0486 0.0484 | 0.0465 0.0533 | 0.0511 | 0.0505 | Ave | | 0.0497 | | | 0.0100 | 6.1 | | 20.0 | | | |
| cis-1,2-Dichloroethene | 0.3357 0.3634 | 0.3627 0.3286 | 0.3396 0.3403 | 0.3473 | 0.3378 | Ave | | 0.3444 | | | 0.1000 | 3.7 | | 20.0 | | | |
| 2-Butanone (MEK) | 0.1549 0.1656 | 0.1778 0.1399 | 0.1558 0.1417 | 0.1537 | 0.1408 | Ave | | 0.1538 | | | 0.0500 | 8.6 | | 20.0 | | | |
| Bromochloromethane | 0.1516 0.1607 | 0.1622 0.1458 | 0.1383 0.1496 | 0.1503 | 0.1438 | Ave | | 0.1503 | | | 0.0100 | 5.4 | | 20.0 | | | |
| Tetrahydrofuran | 0.1403 0.1029 | 0.1208 0.0905 | 0.0949 0.0946 | 0.0965 | 0.0924 | Ave | | 0.1041 | | | 0.0100 | 16.8 | | 20.0 | | | |
| Chloroform | 0.5046 0.5414 | 0.5408 0.4888 | 0.5225 0.5010 | 0.5258 | 0.5131 | Ave | | 0.5173 | | | 0.2000 | 3.6 | | 20.0 | | | |
| 1,1,1-Trichloroethane | 0.3059 0.3476 | 0.3251 0.3197 | 0.3211 0.3440 | 0.3302 | 0.3364 | Ave | | 0.3287 | | | 0.1000 | 4.2 | | 20.0 | | | |
| Cyclohexane | 0.5176 0.4515 | 0.4846 0.4234 | 0.4744 0.4585 | 0.4759 | 0.4508 | Ave | | 0.4671 | | | 0.1000 | 6.0 | | 20.0 | | | |
| Carbon tetrachloride | 0.1966 0.2579 | 0.2202 0.2439 | 0.2226 0.2730 | 0.2389 | 0.2388 | Ave | | 0.2365 | | | 0.1000 | 10.0 | | 20.0 | | | |
| 1,1-Dichloropropene | 0.3981 0.4024 | 0.4008 0.3732 | 0.3987 0.3992 | 0.4036 | 0.3921 | Ave | | 0.3960 | | | 0.0100 | 2.5 | | 20.0 | | | |

Note: The M1 coefficient is the same as Ave RRF for an Ave curve type.

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

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

Analy Batch No.: 217861

SDG No.:

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

| 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 | | | | | | | | |
| Benzene | 1.2659 1.1411 | 1.3110 1.0204 | 1.1612 1.0175 | 1.1979 | 1.1046 | Ave | | 1.1524 | | | 0.5000 | 9.2 | | 20.0 | | | |
| Isobutyl alcohol | 0.0068 0.0078 | 0.0074 0.0073 | 0.0062 0.0074 | 0.0067 | 0.0061 | Ave | | 0.0069 | | * | 0.0100 | 9.0 | | 20.0 | | | |
| 1,2-Dichloroethane | 0.4414 0.4413 | 0.4873 0.3948 | 0.4087 0.4001 | 0.4303 | 0.4024 | Ave | | 0.4258 | | | 0.1000 | 7.3 | | 20.0 | | | |
| n-Heptane | 0.3046 0.2655 | 0.2715 0.2397 | 0.2583 0.2675 | 0.2685 | 0.2506 | Ave | | 0.2658 | | | 0.0100 | 7.1 | | 20.0 | | | |
| Trichloroethene | 0.2857 0.2913 | 0.2895 0.2663 | 0.2730 0.2779 | 0.2828 | 0.2736 | Ave | | 0.2800 | | | 0.2000 | 3.1 | | 20.0 | | | |
| Methylcyclohexane | 0.5506 0.4736 | 0.5082 0.4397 | 0.4903 0.4745 | 0.5019 | 0.4725 | Ave | | 0.4889 | | | 0.1000 | 6.7 | | 20.0 | | | |
| 1,2-Dichloropropane | 0.2899 0.3108 | 0.2949 0.2788 | 0.2747 0.2913 | 0.2876 | 0.2769 | Ave | | 0.2881 | | | 0.1000 | 4.1 | | 20.0 | | | |
| 1,4-Dioxane | 0.0030 0.0026 | 0.0033 0.0022 | 0.0027 0.0024 | 0.0026 | 0.0023 | Ave | | 0.0026 | | * | 0.0100 | 13.6 | | 20.0 | | | |
| Dibromomethane | 0.1974 0.2025 | 0.1996 0.1829 | 0.1756 0.1868 | 0.1860 | 0.1767 | Ave | | 0.1884 | | | 0.0100 | 5.5 | | 20.0 | | | |
| Bromodichloromethane | 0.2608 0.3692 | 0.3120 0.3396 | 0.2889 0.3543 | 0.3275 | 0.3140 | Ave | | 0.3208 | | | 0.2000 | 10.9 | | 20.0 | | | |
| 2-Chloroethyl vinyl ether | 0.1868 0.2074 | 0.2167 0.1868 | 0.1781 0.1938 | 0.1904 | 0.1674 | Ave | | 0.1909 | | | 0.0100 | 8.2 | | 20.0 | | | |
| cis-1,3-Dichloropropene | 0.2705 0.4410 | 0.3316 0.3997 | 0.3266 0.4142 | 0.3665 | 0.3611 | Ave | | 0.3639 | | | 0.2000 | 15.0 | | 20.0 | | | |
| 4-Methyl-2-pentanone (MIBK) | 1.3134 1.2837 | 1.4428 1.1619 | 1.3220 1.0694 | 1.3558 | 1.2661 | Ave | | 1.2769 | | | 0.1000 | 9.1 | | 20.0 | | | |
| Toluene | 6.0508 4.5624 | 5.5576 4.1666 | 5.1577 3.8337 | 5.2430 | 4.8472 | Ave | | 4.9274 | | | 0.4000 | 14.8 | | 20.0 | | | |
| trans-1,3-Dichloropropene | 1.0419 1.5300 | 1.2118 1.4618 | 1.1943 1.4145 | 1.3692 | 1.3334 | Ave | | 1.3196 | | | 0.1000 | 12.2 | | 20.0 | | | |
| Ethyl methacrylate | 1.5098 1.7677 | 1.7528 1.6625 | 1.6297 1.5641 | 1.7656 | 1.6495 | Ave | | 1.6627 | | | 0.0100 | 5.8 | | 20.0 | | | |
| 1,1,2-Trichloroethane | 1.1786 1.1625 | 1.2819 1.0786 | 1.1275 1.0374 | 1.1670 | 1.0865 | Ave | | 1.1400 | | | 0.1000 | 6.6 | | 20.0 | | | |
| Tetrachloroethene | 1.0463 0.8489 | 0.8922 0.8062 | 0.8845 0.7981 | 0.9220 | 0.8591 | Ave | | 0.8822 | | | 0.2000 | 8.9 | | 20.0 | | | |
| 1,3-Dichloropropane | 2.2848 2.0510 | 2.3788 1.8937 | 2.0325 1.7922 | 2.1291 | 1.9759 | Ave | | 2.0672 | | | 0.0100 | 9.4 | | 20.0 | | | |
| 2-Hexanone | 0.8467 0.8290 | 0.9552 0.7643 | 0.8185 0.7213 | 0.8612 | 0.7781 | Ave | | 0.8218 | | | 0.1000 | 8.6 | | 20.0 | | | |

Note: The M1 coefficient is the same as Ave RRF for an Ave curve type.

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

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

Analy Batch No.: 217861

SDG No.:

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

| 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 | 0.6235 0.9071 | 0.7065 0.8567 | 0.6998 0.8436 | 0.7772 | 0.7651 | Ave | | 0.7724 | | | 0.1000 | 12.2 | | 20.0 | | | |
| 1,2-Dibromoethane (EDB) | 1.0866 1.1527 | 1.1708 1.0816 | 1.0670 1.0267 | 1.1262 | 1.0622 | Ave | | 1.0967 | | | 0.1000 | 4.5 | | 20.0 | | | |
| 3-Chlorobenzotrifluoride | 1.5158 1.3678 | 1.5396 1.3015 | 1.4427 1.2972 | 1.5937 | 1.3990 | Ave | | 1.4322 | | | 0.0100 | 7.7 | | 20.0 | | | |
| Chlorobenzene | 3.8282 3.0865 | 3.6360 2.8090 | 3.2721 2.6244 | 3.3129 | 3.0652 | Ave | | 3.2043 | | | 0.5000 | 12.5 | | 20.0 | | | |
| 4-Chlorobenzotrifluoride | 1.3979 1.2982 | 1.3616 1.2351 | 1.3127 1.2373 | 1.5001 | 1.2773 | Ave | | 1.3275 | | | 0.0100 | 6.7 | | 20.0 | | | |
| 1,1,1,2-Tetrachloroethane | 0.7632 1.0328 | 0.8610 0.9619 | 0.8342 0.9319 | 0.9709 | 0.9327 | Ave | | 0.9111 | | | 0.0100 | 9.5 | | 20.0 | | | |
| Ethylbenzene | 2.0557 1.8003 | 1.9366 1.6726 | 1.8780 1.5959 | 1.9373 | 1.7866 | Ave | | 1.8329 | | | 0.1000 | 8.2 | | 20.0 | | | |
| m-Xylene & p-Xylene | 2.5501 2.2147 | 2.4209 2.0393 | 2.2209 1.9570 | 2.3806 | 2.1868 | Ave | | 2.2463 | | | 0.1000 | 8.8 | | 20.0 | | | |
| o-Xylene | 2.4843 2.2465 | 2.4921 2.0108 | 2.2993 1.9118 | 2.3599 | 2.2034 | Ave | | 2.2510 | | | 0.3000 | 9.2 | | 20.0 | | | |
| Styrene | 4.0891 3.5907 | 4.0950 3.2108 | 3.7484 2.9710 | 3.9219 | 3.4981 | Ave | | 3.6406 | | | 0.3000 | 11.2 | | 20.0 | | | |
| Bromoform | 0.3370 0.5330 | 0.3789 0.5010 | 0.3818 0.4964 | 0.4305 | 0.4190 | Ave | | 0.4347 | | | 0.1000 | 15.9 | | 20.0 | | | |
| 2-Chlorobenzotrifluoride | 1.4533 1.4424 | 1.6153 1.3223 | 1.4179 1.3036 | 1.6403 | 1.4193 | Ave | | 1.4518 | | | 0.0100 | 8.4 | | 20.0 | | | |
| Isopropylbenzene | 6.3984 4.6371 | 5.8487 4.0713 | 5.3090 3.7512 | 5.4219 | 4.8709 | Ave | | 5.0386 | | | 0.1000 | 17.6 | | 20.0 | | | |
| Bromobenzene | 0.9410 0.9394 | 0.9810 0.9122 | 0.9383 0.9038 | 0.9005 | 0.9180 | Ave | | 0.9293 | | | 0.0100 | 2.8 | | 20.0 | | | |
| 1,1,2,2-Tetrachloroethane | 1.6867 1.6395 | 1.8800 1.4633 | 1.6064 1.3686 | 1.6823 | 1.5069 | Ave | | 1.6042 | | | 0.3000 | 9.9 | | 20.0 | | | |
| trans-1,4-Dichloro-2-butene | 0.2724 0.2884 | 0.2827 0.2833 | 0.2498 0.2904 | 0.2507 | 0.2582 | Ave | | 0.2720 | | | 0.0100 | 6.2 | | 20.0 | | | |
| 1,2,3-Trichloropropane | 0.3844 0.3903 | 0.4180 0.3743 | 0.3879 0.3799 | 0.3764 | 0.3765 | Ave | | 0.3860 | | | 0.0100 | 3.7 | | 20.0 | | | |
| N-Propylbenzene | 1.0896 0.9901 | 1.0404 0.9508 | 1.0060 0.9746 | 1.0147 | 1.0247 | Ave | | 1.0114 | | | 0.0100 | 4.2 | | 20.0 | | | |
| 2-Chlorotoluene | 0.9244 0.8870 | 0.8858 0.8323 | 0.8682 0.8579 | 0.8757 | 0.8785 | Ave | | 0.8762 | | | 0.0100 | 3.0 | | 20.0 | | | |
| 3-Chlorotoluene | 0.9488 0.9026 | 0.9481 0.8733 | 0.9020 0.9002 | 0.9547 | 0.9252 | Ave | | 0.9194 | | | 0.0100 | 3.2 | | 20.0 | | | |

Note: The M1 coefficient is the same as Ave RRF for an Ave curve type.

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

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

Analy Batch No.: 217861

SDG No.: _____

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

| 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 | 3.2098 2.7572 | 3.1610 2.5370 | 2.9762 2.4836 | 2.9682 | 2.8830 | Ave | | 2.8720 | | | 0.0100 | 9.3 | | 20.0 | | | |
| 4-Chlorotoluene | 1.0814 0.9671 | 0.9434 0.9047 | 0.9555 0.9161 | 0.9436 | 0.9398 | Ave | | 0.9565 | | | 0.0100 | 5.7 | | 20.0 | | | |
| tert-Butylbenzene | 2.5322 2.2077 | 2.3962 2.0211 | 2.3742 2.0481 | 2.3326 | 2.3286 | Ave | | 2.2801 | | | 0.0100 | 7.7 | | 20.0 | | | |
| 1,2,4-Trimethylbenzene | 3.5127 2.8439 | 3.3123 2.5942 | 3.1049 2.5435 | 3.0350 | 3.0086 | Ave | | 2.9944 | | | 0.0100 | 11.1 | | 20.0 | | | |
| 3,4-Dichlorobenzotrifluoride | 0.7362 0.6508 | 0.6581 0.6193 | 0.6198 0.6784 | 0.6940 | 0.6587 | Ave | | 0.6644 | | | 0.0100 | 5.8 | | 20.0 | | | |
| sec-Butylbenzene | 3.9837 2.9724 | 3.4720 2.6959 | 3.2967 2.6895 | 3.2643 | 3.2125 | Ave | | 3.1984 | | | 0.0100 | 13.3 | | 20.0 | | | |
| 1,3-Dichlorobenzene | 1.9656 1.6199 | 1.7934 1.5145 | 1.6495 1.5107 | 1.6503 | 1.6466 | Ave | | 1.6688 | | | 0.6000 | 8.9 | | 20.0 | | | |
| 4-Isopropyltoluene | 3.0764 2.5117 | 2.9764 2.2679 | 2.6994 2.2836 | 2.6864 | 2.6251 | Ave | | 2.6409 | | | 0.0100 | 11.0 | | 20.0 | | | |
| 1,4-Dichlorobenzene | 1.9865 1.6860 | 1.8621 1.5565 | 1.7487 1.5598 | 1.7034 | 1.7043 | Ave | | 1.7259 | | | 0.5000 | 8.4 | | 20.0 | | | |
| 2,4-Dichlorobenzotrifluoride | 0.6454 0.6171 | 0.6327 0.5625 | 0.6090 0.6636 | 0.6702 | 0.6085 | Ave | | 0.6261 | | | 0.0100 | 5.6 | | 20.0 | | | |
| 2,5-Dichlorobenzotrifluoride | 0.6732 0.7026 | 0.7481 0.6866 | 0.6523 0.6767 | 0.7060 | 0.6830 | Ave | | 0.6911 | | | 0.0100 | 4.1 | | 20.0 | | | |
| n-Butylbenzene | 2.8298 2.2864 | 2.6464 2.0727 | 2.4207 2.1235 | 2.4803 | 2.4098 | Ave | | 2.4087 | | | 0.0100 | 10.5 | | 20.0 | | | |
| 1,2-Dichlorobenzene | 1.7708 1.5647 | 1.7956 1.4356 | 1.5665 1.4544 | 1.5674 | 1.5442 | Ave | | 1.5874 | | | 0.4000 | 8.3 | | 20.0 | | | |
| 1,2-Dibromo-3-Chloropropane | 0.1216 0.1606 | 0.1391 0.1568 | 0.1266 0.1669 | 0.1326 | 0.1278 | Ave | | 0.1415 | | | 0.0500 | 12.4 | | 20.0 | | | |
| 2,4- & 2,5- & 2,6- Dichlorotoluene | 1.0715 0.9570 | 1.1692 0.9265 | 0.9793 0.9637 | 1.0382 | 0.9824 | Ave | | 1.0110 | | | 0.0100 | 7.8 | | 20.0 | | | |
| 2,3- & 3,4- Dichlorotoluene | 1.1353 1.0683 | 1.2293 1.0765 | 1.0355 1.1269 | 1.0901 | 1.0726 | Ave | | 1.1043 | | | 0.0100 | 5.4 | | 20.0 | | | |
| 1,2,4-Trichlorobenzene | 0.8306 0.8182 | 0.9355 0.8486 | 0.7503 0.8953 | 0.7691 | 0.7875 | Ave | | 0.8294 | | | 0.2000 | 7.6 | | 20.0 | | | |
| Hexachlorobutadiene | 0.3307 0.2462 | 0.2635 0.2582 | 0.2213 0.2957 | 0.2402 | 0.2448 | Ave | | 0.2626 | | | 0.0100 | 13.3 | | 20.0 | | | |
| Naphthalene | 3.1019 2.3472 | 2.8951 2.4151 | 2.3638 2.3957 | 2.3524 | 2.3536 | Ave | | 2.5281 | | | 0.0100 | 11.7 | | 20.0 | | | |
| 1,2,3-Trichlorobenzene | 0.7804 0.7270 | 0.8087 0.8154 | 0.6385 0.8565 | 0.6517 | 0.6792 | Ave | | 0.7447 | | | 0.0100 | 11.0 | | 20.0 | | | |

Note: The M1 coefficient is the same as Ave RRF for an Ave curve type.

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

Lab Name: TestAmerica Pittsburgh Job No.: 180-71131-1 Analy Batch No.: 217861
SDG No.: _____
Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N
Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

| 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 | | | | | | | | |
| 2,4,5-Trichlorotoluene | 0.4043 0.4250 | 0.4160 0.5255 | 0.3407 0.5717 | 0.3454 | 0.3978 | Ave | | 0.4283 | | | 0.0100 | 19.0 | | 20.0 | | | |
| 2,3,6-Trichlorotoluene | 0.3645 0.3663 | 0.3955 0.4558 | 0.3167 0.5003 | 0.3075 | 0.3648 | Ave | | 0.3839 | | | 0.0100 | 17.1 | | 20.0 | | | |
| Dibromofluoromethane (Surr) | 0.3105 0.2657 | 0.2699 0.2422 | 0.2388 0.2506 | 0.2560 | 0.2449 | Ave | | 0.2598 | | | | 8.9 | | 20.0 | | | |
| 1,2-Dichloroethane-d4 (Surr) | 0.4947 0.3692 | 0.4116 0.3263 | 0.3458 0.3341 | 0.3589 | 0.3302 | Ave | | 0.3713 | | | | 15.4 | | 20.0 | | | |
| Toluene-d8 (Surr) | 5.9733 3.6566 | 4.7106 3.3970 | 4.0457 ++++ | 4.1808 | 3.8728 | Lin2 | 11.440 | 3.7819 | | | | | | | 0.9930 | | 0.9900 |
| 4-Bromofluorobenzene (Surr) | 2.3906 1.6487 | 1.9549 1.4705 | 1.6323 1.4101 | 1.7133 | 1.5905 | Ave | | 1.7264 | | | | 18.2 | | 20.0 | | | |

Note: The M1 coefficient is the same as Ave RRF for an Ave curve type.

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

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

Analy Batch No.: 217861

SDG No.: _____

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---------|-------------------|--------------|
| Level 1 | IC 180-217861/3 | 60724D03.D |
| Level 2 | IC 180-217861/4 | 60724D04.D |
| Level 3 | ICIS 180-217861/5 | 60724D05.D |
| Level 4 | IC 180-217861/6 | 60724D06.D |
| Level 5 | IC 180-217861/7 | 60724D07.D |
| Level 6 | IC 180-217861/8 | 60724D08.D |
| Level 7 | IC 180-217861/9 | 60724D09.D |
| Level 8 | IC 180-217861/10 | 60724D10.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 | 27479 839726 | 121547 998454 | 286699 1341098 | 392309 | 537841 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Chloromethane | FB | Ave | 23920 792135 | 109949 904618 | 249199 1170186 | 364709 | 485997 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Vinyl chloride | FB | Ave | 26240 788826 | 112168 942517 | 262771 1244722 | 390082 | 507676 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,3-Butadiene | FB | Ave | 24629 626705 | 92248 738623 | 223993 1011679 | 315046 | 407662 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Bromomethane | FB | Ave | 12374 390167 | 58670 421777 | 122895 482936 | 178416 | 239488 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Chloroethane | FB | Ave | 14369 445022 | 64213 496292 | 137952 595090 | 212582 | 283541 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Trichlorofluoromethane | FB | Ave | 24407 820933 | 108123 958134 | 244680 1263842 | 381467 | 515987 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Ethyl ether | FB | Ave | 21539 720418 | 103875 817030 | 207890 1004152 | 314417 | 412759 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Acrolein | FB | Ave | 93239 203936 | 103236 222222 | 126353 255428 | 161845 | 172660 | 100 225 | 125 250 | 150 275 | 175 | 200 |
| 1,1-Dichloroethene | FB | Ave | 22840 719926 | 90501 850942 | 212019 1153420 | 326499 | 437661 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | FB | Ave | 20417 680574 | 89949 799492 | 214148 1098540 | 320427 | 418931 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Acetone | FB | Ave | 47909 658887 | 83742 631699 | 181114 730103 | 295809 | 345529 | 25.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| Iodomethane | FB | Ave | 29488 1045605 | 133154 1229211 | 296892 1578693 | 465530 | 616342 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Carbon disulfide | FB | Ave | 41100 1759930 | 1779135 2106147 | 451961 2838844 | 721571 | 1031794 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Allyl chloride | FB | Ave | 11154 457646 | 50272 540192 | 120354 717090 | 194002 | 261163 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |

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

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

Analy Batch No.: 217861

SDG No.:

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

| 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 |
| Methyl acetate | FB | Ave | 40003 1378375 | 185280 1509491 | 382502 1923255 | 572896 | 775230 | 10.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| Methylene Chloride | FB | Ave | 34286 940505 | 131057 1081026 | 278118 1372983 | 418660 | 557470 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| tert-Butyl alcohol | TBAd 9 | Ave | 20227 572173 | 71644 544523 | 151604 673044 | 251071 | 297303 | 50.0 1750 | 250 2000 | 500 2500 | 750 | 1000 |
| Acrylonitrile | FB | Ave | 114140 3519987 | 530059 3825550 | 1020211 4723140 | 1582844 | 2040960 | 50.0 1750 | 250 2000 | 500 2500 | 750 | 1000 |
| trans-1,2-Dichloroethene | FB | Ave | 24529 817284 | 108647 968185 | 244707 1277192 | 372663 | 496677 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Methyl tert-butyl ether | FB | Ave | 78478 2689634 | 375187 3053893 | 782498 3809189 | 1181045 | 1550808 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Hexane | FB | Ave | 30509 940765 | 127818 1103506 | 304164 1517069 | 437873 | 573395 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,1-Dichloroethane | FB | Ave | 27271 1429058 | 187923 1631648 | 418314 2087088 | 641771 | 843415 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Vinyl acetate | FB | Ave | 41136 1612184 | 217747 1962237 | 435645 2492452 | 703120 | 926426 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2,2-Dichloropropane | FB | Ave | 3609 147487 | 17106 168555 | 39797 228817 | 64241 | 87515 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| cis-1,2-Dichloroethene | FB | Ave | 26745 997518 | 127777 1143732 | 290355 1462208 | 436220 | 585924 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2-Butanone (MEK) | FB | Ave | 61711 909301 | 125277 973759 | 266460 1217733 | 385993 | 488342 | 25.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| Bromochloromethane | FB | Ave | 12079 441109 | 57129 507483 | 118276 642539 | 188762 | 249508 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Tetrahydrofuran | FB | Ave | 22350 565011 | 85090 629732 | 162324 812537 | 242295 | 320475 | 10.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| Chloroform | FB | Ave | 40199 1486297 | 190538 1701079 | 446745 2152497 | 660420 | 889966 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,1,1-Trichloroethane | FB | Ave | 24367 954279 | 114539 112586 | 274535 1477890 | 414781 | 583435 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Cyclohexane | FB | Ave | 41233 1239342 | 170750 1473582 | 405642 1969875 | 597815 | 781868 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Carbon tetrachloride | FB | Ave | 15662 707925 | 77574 848801 | 190304 1172757 | 300016 | 414150 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,1-Dichloropropene | FB | Ave | 31712 1104618 | 141227 1298950 | 340889 1715254 | 506906 | 680030 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Benzene | FB | Ave | 100837 3132477 | 461882 3551507 | 992835 4371437 | 1504594 | 1915947 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Isobutyl alcohol | FB | Ave | 13528 537905 | 64845 631529 | 132080 789622 | 208868 | 262400 | 125 4375 | 625 5000 | 1250 6250 | 1875 | 2500 |

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

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

Analy Batch No.: 217861

SDG No.:

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

| 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 |
| 1,2-Dichloroethane | FB | Ave | 35160 1211453 | 171694 1374193 | 349409 1719102 | 540487 | 697923 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| n-Heptane | FB | Ave | 24265 728795 | 95666 834225 | 220812 1149322 | 337209 | 434669 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Trichloroethene | FB | Ave | 22756 799778 | 102011 926685 | 233389 1194165 | 355153 | 474544 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Methylcyclohexane | FB | Ave | 43857 1300055 | 179050 1530213 | 419227 2038808 | 630450 | 819514 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2-Dichloropropane | FB | Ave | 23091 853109 | 103896 970394 | 234878 1251517 | 361293 | 480287 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,4-Dioxane | FB | Ave | 4766 141286 | 23040 152489 | 45919 203123 | 65610 | 80878 | 100 3500 | 500 4000 | 1000 5000 | 1500 | 2000 |
| Dibromomethane | FB | Ave | 15724 555898 | 70312 636427 | 150124 802657 | 233620 | 306535 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Bromodichloromethane | FB | Ave | 20777 1013623 | 109927 1181877 | 246980 1522204 | 411363 | 544627 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2-Chloroethyl vinyl ether | FB | Ave | 29761 1138561 | 152667 1300177 | 304468 1665721 | 478201 | 580737 | 10.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| cis-1,3-Dichloropropene | FB | Ave | 21547 1210517 | 116838 1391254 | 279255 1779441 | 460333 | 626305 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 4-Methyl-2-pentanone (MIBK) | CBNZ d5 | Ave | 113074 1737974 | 233783 1932325 | 495598 2362456 | 755882 | 970750 | 25.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| Toluene | CBNZ d5 | Ave | 104189 3088570 | 450260 3464609 | 966776 4234419 | 1461492 | 1858285 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| trans-1,3-Dichloropropene | CBNZ d5 | Ave | 17940 1035772 | 98174 1215519 | 223869 1562404 | 381658 | 511188 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Ethyl methacrylate | CBNZ d5 | Ave | 25997 1196636 | 142004 1382390 | 305467 1727540 | 492154 | 632371 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,1,2-Trichloroethane | CBNZ d5 | Ave | 20294 786936 | 103854 896840 | 211345 1145832 | 325307 | 416541 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Tetrachloroethene | CBNZ d5 | Ave | 18016 574638 | 72287 670325 | 165794 881532 | 257015 | 329342 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,3-Dichloropropane | CBNZ d5 | Ave | 39343 1388403 | 192724 1574600 | 380984 1979514 | 593488 | 757496 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2-Hexanone | CBNZ d5 | Ave | 72896 1122326 | 154768 1271094 | 306856 1593297 | 480105 | 596567 | 25.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| Dibromochloromethane | CBNZ d5 | Ave | 10736 614068 | 57240 712324 | 131179 931753 | 216633 | 293309 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2-Dibromoethane (EDB) | CBNZ d5 | Ave | 18711 780325 | 94851 899338 | 200005 1134020 | 313940 | 407201 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 3-Chlorobenzotrifluoride | CBNZ d5 | Ave | 26101 925971 | 124733 1082251 | 270429 1432766 | 444234 | 536353 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |

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

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

Analy Batch No.: 217861

SDG No.: _____

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

| 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 | CBNZ d5 | Ave | 65919 2089428 | 294577 2335758 | 613324 2898680 | 923461 | 1175123 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 4-Chlorobenzotrifluoride | CBNZ d5 | Ave | 24071 878791 | 110312 1026977 | 246049 1366587 | 418159 | 489696 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,1,1,2-Tetrachloroethane | CBNZ d5 | Ave | 13142 699176 | 69755 799872 | 156373 1029362 | 270626 | 357566 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Ethylbenzene | CBNZ d5 | Ave | 35398 1218751 | 156900 1390812 | 352019 1762716 | 540015 | 684943 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| m-Xylene & p-Xylene | CBNZ d5 | Ave | 43910 1499275 | 196132 1695741 | 416283 2161557 | 663604 | 838371 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| o-Xylene | CBNZ d5 | Ave | 42778 1520782 | 201906 1672026 | 430994 2111655 | 657817 | 844701 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Styrene | CBNZ d5 | Ave | 70410 2430770 | 331761 2669824 | 702614 3281557 | 1093228 | 1341052 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Bromoform | CBNZ d5 | Ave | 5803 360803 | 30696 416604 | 71560 548248 | 120009 | 160632 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2-Chlorobenzotrifluoride | CBNZ d5 | Ave | 25025 976474 | 130865 1099473 | 265784 1439864 | 457223 | 544128 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Isopropylbenzene | CBNZ d5 | Ave | 110174 3139141 | 473843 3385367 | 995129 4143279 | 1511361 | 1867348 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Bromobenzene | DCBd 4 | Ave | 25122 913147 | 122999 1008153 | 249243 1259105 | 380643 | 474286 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,1,2,2-Tetrachloroethane | CBNZ d5 | Ave | 29044 1109880 | 152313 1216769 | 301111 1511708 | 468952 | 577708 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| trans-1,4-Dichloro-2-butene | DCBd 4 | Ave | 7272 280292 | 35451 313051 | 66355 404523 | 105972 | 133396 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2,3-Trichloropropane | DCBd 4 | Ave | 10261 379404 | 52406 413676 | 103048 529322 | 159086 | 194531 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| N-Propylbenzene | DCBd 4 | Ave | 29088 962443 | 130443 1050858 | 267230 1357847 | 428911 | 529403 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2-Chlorotoluene | DCBd 4 | Ave | 24679 862193 | 111064 919859 | 230624 1195161 | 370145 | 453885 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 3-Chlorotoluene | DCBd 4 | Ave | 25329 877299 | 118876 965183 | 239610 1254125 | 403537 | 478005 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,3,5-Trimethylbenzene | DCBd 4 | Ave | 85688 2680034 | 396334 2803848 | 790599 3460126 | 1254643 | 1489442 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 4-Chlorotoluene | DCBd 4 | Ave | 28869 940075 | 118288 999834 | 253826 1276343 | 398838 | 485508 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| tert-Butylbenzene | DCBd 4 | Ave | 67600 2145902 | 300432 2233746 | 630674 2853353 | 985989 | 1203013 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2,4-Trimethylbenzene | DCBd 4 | Ave | 93775 2764298 | 415297 2867121 | 824782 3543615 | 1282863 | 1554360 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |

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

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

Analy Batch No.: 217861

SDG No.: _____

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

| 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 | DCBd 4 | Ave | 19655 632603 | 82510 684446 | 164643 945109 | 293347 | 340323 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| sec-Butylbenzene | DCBd 4 | Ave | 106349 2889256 | 435324 2979536 | 875734 3747062 | 1379795 | 1659704 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,3-Dichlorobenzene | DCBd 4 | Ave | 52474 1574554 | 224858 1673821 | 438172 2104721 | 697551 | 850676 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 4-Isopropyltoluene | DCBd 4 | Ave | 82129 2441419 | 373177 2506437 | 717050 3181497 | 1135538 | 1356230 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,4-Dichlorobenzene | DCBd 4 | Ave | 53033 1638813 | 233465 1720245 | 464508 2173124 | 720028 | 880492 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2,4-Dichlorobenzotrifluoride | DCBd 4 | Ave | 17229 599851 | 79327 621645 | 161786 924522 | 283275 | 314370 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2,5-Dichlorobenzotrifluoride | DCBd 4 | Ave | 17972 682948 | 93800 758780 | 173267 942783 | 298436 | 352876 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| n-Butylbenzene | DCBd 4 | Ave | 75545 2222409 | 331802 2290785 | 643029 2958420 | 1048415 | 1244987 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2-Dichlorobenzene | DCBd 4 | Ave | 47273 1520936 | 225127 1586590 | 416117 2026312 | 662537 | 797769 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2-Dibromo-3-Chloropropane | DCBd 4 | Ave | 3245 156142 | 17440 173322 | 33625 232535 | 56051 | 66004 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2,4- & 2,5- & 2,6- Dichlorotoluene | DCBd 4 | Ave | 85812 2790747 | 439793 3071993 | 780454 4027755 | 1316529 | 1522613 | 15.0 525 | 75.0 600 | 150 750 | 225 | 300 |
| 2,3- & 3,4- Dichlorotoluene | DCBd 4 | Ave | 60616 2076832 | 308264 2379503 | 550121 3139946 | 921595 | 1108272 | 10.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| 1,2,4-Trichlorobenzene | DCBd 4 | Ave | 22175 795349 | 117288 937825 | 199297 1247374 | 325076 | 406868 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Hexachlorobutadiene | DCBd 4 | Ave | 8829 239351 | 33039 285364 | 58792 411971 | 101525 | 126465 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Naphthalene | DCBd 4 | Ave | 82809 2281539 | 362983 2669188 | 627907 3337709 | 994327 | 1215966 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2,3-Trichlorobenzene | DCBd 4 | Ave | 20833 706689 | 101393 901210 | 169607 1193234 | 275471 | 350907 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2,4,5-Trichlorotoluene | DCBd 4 | Ave | 10792 413111 | 52163 580730 | 90501 796492 | 146009 | 205495 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2,3,6-Trichlorotoluene | DCBd 4 | Ave | 9731 356014 | 49582 503740 | 84129 697018 | 129996 | 188457 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Dibromofluoromethane (Surr) | FB | Ave | 24732 729396 | 95101 842973 | 204197 1076618 | 321543 | 424756 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2-Dichloroethane-d4 (Surr) | FB | Ave | 39405 1013596 | 145003 1135703 | 295614 1435595 | 450831 | 572691 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Toluene-d8 (Surr) | CBNZ d5 | Lin2 | 102855 2475360 | 381639 2824683 | 758339 +++++ | 1165400 | 1484720 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |

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

Lab Name: TestAmerica Pittsburgh Job No.: 180-71131-1 Analy Batch No.: 217861

SDG No.: _____

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

| 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 |
| 4-Bromofluorobenzene (Surr) | CBNZ d5 | Ave | 41164 1116127 | 158379 1222775 | 305961 1557524 | 477594 | 609762 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |

Curve Type Legend:

Ave = Average ISTD

Lin2 = Linear 1/conc^2 ISTD

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
READBACK PERCENT ERROR

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

Analy Batch No.: 217861

SDG No.: _____

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---------|-------------------|--------------|
| Level 1 | IC 180-217861/3 | 60724D03.D |
| Level 2 | IC 180-217861/4 | 60724D04.D |
| Level 3 | ICIS 180-217861/5 | 60724D05.D |
| Level 4 | IC 180-217861/6 | 60724D06.D |
| Level 5 | IC 180-217861/7 | 60724D07.D |
| Level 6 | IC 180-217861/8 | 60724D08.D |
| Level 7 | IC 180-217861/9 | 60724D09.D |
| Level 8 | IC 180-217861/10 | 60724D10.D |

| ANALYTE | PERCENT ERROR | | | | | | PERCENT ERROR LIMIT | | | | | |
|---------------------------------------|--------------------|--------------------|---------|---------|---------|---------|---------------------|----------------|-------|-------|-------|-------|
| | LVL 1 # LVL 7 # | LVL 2 # LVL 8 # | LVL 3 # | LVL 4 # | LVL 5 # | LVL 6 # | LVL 1 LVL 7 | LVL 2 LVL 8 | LVL 3 | LVL 4 | LVL 5 | LVL 6 |
| Dichlorodifluoromethane | 8.1 -10.1 | 8.1 -2.2 | 5.1 | -2.1 | -2.8 | -4.1 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Chloromethane | 4.7 -9.4 | 8.8 -5.1 | 1.6 | 1.2 | -2.3 | 0.6 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Vinyl chloride | 9.5 -10.0 | 5.8 -3.7 | 2.2 | 3.3 | -2.7 | -4.5 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,3-Butadiene | 24.0 -14.9 | 5.0 -5.6 | 5.1 | 0.6 | -5.7 | -8.4 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Bromomethane | 10.8 -13.5 | 18.8 -19.8 | 2.5 | 1.3 | -1.5 | 1.4 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Chloroethane | 11.0 -12.2 | 12.2 -14.8 | -0.7 | 4.2 | 0.6 | -0.2 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Trichlorofluoromethane | 3.5 -7.0 | 3.6 -0.7 | -3.4 | 2.6 | 0.4 | 1.0 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Ethyl ether | 6.7 -7.4 | 16.3 -7.8 | -4.1 | -1.2 | -6.1 | 3.5 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Acrolein | 7.8 -5.9 | 8.0 -0.4 | -9.3 | 1.7 | -8.3 | 6.4 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,1-Dichloroethene | 10.3 -5.9 | -1.2 3.3 | -4.6 | 0.0 | -2.9 | 0.9 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 2.9 -7.7 | 2.5 2.7 | 0.6 | 2.5 | -3.0 | -0.4 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Acetone | 12.1 -15.4 | 10.8 -20.8 | -1.3 | 9.8 | -7.1 | 11.9 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Iodomethane | 1.3 -3.3 | 3.4 0.6 | -5.0 | 1.4 | -2.7 | 4.3 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Carbon disulfide | -10.8 4.6 | -12.1 14.2 | -8.6 | -0.7 | 2.8 | 10.8 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Allyl chloride | -8.0 2.0 | -6.2 9.7 | -7.5 | 1.5 | -1.1 | 9.6 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
READBACK PERCENT ERROR

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

Analy Batch No.: 217861

SDG No.:

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

| ANALYTE | PERCENT ERROR | | | | | | PERCENT ERROR LIMIT | | | | | |
|--------------------------|--------------------|--------------------|---------|---------|---------|---------|---------------------|----------------|-------|-------|-------|-------|
| | LVL 1 # LVL 7 # | LVL 2 # LVL 8 # | LVL 3 # | LVL 4 # | LVL 5 # | LVL 6 # | LVL 1 LVL 7 | LVL 2 LVL 8 | LVL 3 | LVL 4 | LVL 5 | LVL 6 |
| Methyl acetate | 6.8 -7.8 | 11.8 -4.8 | -4.9 | -3.0 | -5.0 | 6.8 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Methylene Chloride | 25.0 -9.8 | 8.0 -7.2 | -5.5 | -3.2 | -6.7 | -0.5 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| tert-Butyl alcohol | 6.9 6.4 | -3.8 -5.1 | -2.6 | -1.2 | -1.5 | 0.9 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| Acrylonitrile | 14.1 -12.5 | 19.8 -12.5 | -5.0 | 0.3 | -6.3 | 2.1 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| trans-1,2-Dichloroethene | 4.4 -5.7 | 4.6 0.8 | -2.9 | 0.6 | -2.9 | 1.0 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| Methyl tert-butyl ether | 4.5 -6.9 | 12.9 -6.0 | -2.9 | -0.3 | -5.2 | 3.9 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| Hexane | 9.7 -9.2 | 3.9 1.1 | 1.9 | -0.2 | -5.3 | -1.9 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,1-Dichloroethane | -28.6 -2.3 | 11.2 1.3 | 2.0 | 6.5 | 1.4 | 8.5 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| Vinyl acetate | -7.6 0.9 | 10.6 3.8 | -8.8 | 0.2 | -4.4 | 5.1 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2,2-Dichloropropane | -8.8 -2.5 | -2.3 7.2 | -6.3 | 3.0 | 1.6 | 8.2 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| cis-1,2-Dichloroethene | -2.5 -4.6 | 5.3 -1.2 | -1.4 | 0.8 | -1.9 | 5.5 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2-Butanone (MEK) | 0.8 -9.0 | 15.6 -7.8 | 1.3 | -0.1 | -8.5 | 7.7 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| Bromochloromethane | 0.9 -3.0 | 7.9 -0.5 | -8.0 | 0.0 | -4.3 | 6.9 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| Tetrahydrofuran | 34.8 -13.1 | 16.0 -9.2 | -8.8 | -7.3 | -11.3 | -1.1 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| Chloroform | -2.4 -5.5 | 4.6 -3.1 | 1.0 | 1.7 | -0.8 | 4.7 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,1,1-Trichloroethane | -7.0 -2.8 | -1.1 4.6 | -2.3 | 0.5 | 2.3 | 5.7 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| Cyclohexane | 10.8 -9.4 | 3.8 -1.8 | 1.6 | 1.9 | -3.5 | -3.3 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| Carbon tetrachloride | -16.9 3.1 | -6.9 15.4 | -5.9 | 1.0 | 1.0 | 9.1 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,1-Dichloropropene | 0.5 -5.8 | 1.2 0.8 | 0.7 | 1.9 | -1.0 | 1.6 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| Benzene | 9.8 -11.5 | 13.8 -11.7 | 0.8 | 3.9 | -4.2 | -1.0 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |
| Isobutyl alcohol | -2.1 4.6 | 6.1 6.0 | -10.9 | -4.1 | -12.8 | 13.0 | 50 30 30 | 30 30 | 30 | 30 | 30 | 30 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
READBACK PERCENT ERROR

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

Analy Batch No.: 217861

SDG No.: _____

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

| ANALYTE | PERCENT ERROR | | | | | | PERCENT ERROR LIMIT | | | | | |
|-----------------------------|--------------------|--------------------|---------|---------|---------|---------|---------------------|----------------|-------|-------|-------|-------|
| | LVL 1 # LVL 7 # | LVL 2 # LVL 8 # | LVL 3 # | LVL 4 # | LVL 5 # | LVL 6 # | LVL 1 LVL 7 | LVL 2 LVL 8 | LVL 3 | LVL 4 | LVL 5 | LVL 6 |
| 1,2-Dichloroethane | 3.7 -7.3 | 14.5 -6.0 | -4.0 | 1.1 | -5.5 | 3.6 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| n-Heptane | 14.6 -9.8 | 2.2 0.7 | -2.8 | 1.0 | -5.7 | -0.1 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Trichloroethene | 2.0 -4.9 | 3.4 -0.7 | -2.5 | 1.0 | -2.3 | 4.0 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Methylcyclohexane | 12.6 -10.1 | 3.9 -2.9 | 0.3 | 2.7 | -3.4 | -3.1 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,2-Dichloropropane | 0.6 -3.2 | 2.4 1.1 | -4.7 | -0.2 | -3.9 | 7.9 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,4-Dioxane | 13.9 -16.6 | 24.5 -10.0 | 2.2 | -0.6 | -11.3 | -2.0 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Dibromomethane | 4.8 -3.0 | 5.9 -0.9 | -6.8 | -1.3 | -6.2 | 7.5 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Bromodichloromethane | -18.7 5.9 | -2.7 10.4 | -10.0 | 2.1 | -2.1 | 15.1 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2-Chloroethyl vinyl ether | -2.2 -2.2 | 13.5 1.5 | -6.7 | -0.3 | -12.3 | 8.6 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| cis-1,3-Dichloropropene | -25.7 9.8 | -8.9 13.8 | -10.2 | 0.7 | -0.8 | 21.2 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 4-Methyl-2-pentanone (MIBK) | 2.9 -9.0 | 13.0 -16.2 | 3.5 | 6.2 | -0.8 | 0.5 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Toluene | 22.8 -15.4 | 12.8 -22.2 | 4.7 | 6.4 | -1.6 | -7.4 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| trans-1,3-Dichloropropene | -21.0 10.8 | -8.2 7.2 | -9.5 | 3.8 | 1.0 | 15.9 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Ethyl methacrylate | -9.2 0.0 | 5.4 -5.9 | -2.0 | 6.2 | -0.8 | 6.3 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,1,2-Trichloroethane | 3.4 -5.4 | 12.4 -9.0 | -1.1 | 2.4 | -4.7 | 2.0 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Tetrachloroethene | 18.6 -8.6 | 1.1 -9.5 | 0.3 | 4.5 | -2.6 | -3.8 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,3-Dichloropropane | 10.5 -8.4 | 15.1 -13.3 | -1.7 | 3.0 | -4.4 | -0.8 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2-Hexanone | 3.0 -7.0 | 16.2 -12.2 | -0.4 | 4.8 | -5.3 | 0.9 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Dibromochloromethane | -19.3 10.9 | -8.5 9.2 | -9.4 | 0.6 | -1.0 | 17.4 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,2-Dibromoethane (EDB) | -0.9 -1.4 | 6.8 -6.4 | -2.7 | 2.7 | -3.2 | 5.1 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 3-Chlorobenzotrifluoride | 5.8 -9.1 | 7.5 -9.4 | 0.7 | 11.3 | -2.3 | -4.5 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
READBACK PERCENT ERROR

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

Analy Batch No.: 217861

SDG No.:

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

| ANALYTE | PERCENT ERROR | | | | | | PERCENT ERROR LIMIT | | | | | |
|-----------------------------|--------------------|--------------------|---------|---------|---------|---------|---------------------|----------------|-------|-------|-------|-------|
| | LVL 1 # LVL 7 # | LVL 2 # LVL 8 # | LVL 3 # | LVL 4 # | LVL 5 # | LVL 6 # | LVL 1 LVL 7 | LVL 2 LVL 8 | LVL 3 | LVL 4 | LVL 5 | LVL 6 |
| Chlorobenzene | 19.5 -12.3 | 13.5 -18.1 | 2.1 | 3.4 | -4.3 | -3.7 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 4-Chlorobenzotrifluoride | 5.3 -7.0 | 2.6 -6.8 | -1.1 | 13.0 | -3.8 | -2.2 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,1,1,2-Tetrachloroethane | -16.2 5.6 | -5.5 2.3 | -8.4 | 6.6 | 2.4 | 13.4 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Ethylbenzene | 12.2 -8.7 | 5.7 -12.9 | 2.5 | 5.7 | -2.5 | -1.8 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| m-Xylene & p-Xylene | 13.5 -9.2 | 7.8 -12.9 | -1.1 | 6.0 | -2.6 | -1.4 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| o-Xylene | 10.4 -10.7 | 10.7 -15.1 | 2.1 | 4.8 | -2.1 | -0.2 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Styrene | 12.3 -11.8 | 12.5 -18.4 | 3.0 | 7.7 | -3.9 | -1.4 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Bromoform | -22.5 15.3 | -12.8 14.2 | -12.2 | -1.0 | -3.6 | 22.6 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2-Chlorobenzotrifluoride | 0.1 -8.9 | 11.3 -10.2 | -2.3 | 13.0 | -2.2 | -0.6 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Isopropylbenzene | 27.0 -19.2 | 16.1 -25.6 | 5.4 | 7.6 | -3.3 | -8.0 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Bromobenzene | 1.3 -1.8 | 5.6 -2.7 | 1.0 | -3.1 | -1.2 | 1.1 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,1,2,2-Tetrachloroethane | 5.1 -8.8 | 17.2 -14.7 | 0.1 | 4.9 | -6.1 | 2.2 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| trans-1,4-Dichloro-2-butene | 0.2 4.1 | 4.0 6.8 | -8.2 | -7.8 | -5.1 | 6.0 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,2,3-Trichloropropane | -0.4 -3.0 | 8.3 -1.6 | 0.5 | -2.5 | -2.4 | 1.1 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| N-Propylbenzene | 7.7 -6.0 | 2.9 -3.6 | -0.5 | 0.3 | 1.3 | -2.1 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2-Chlorotoluene | 5.5 -5.0 | 1.1 -2.1 | -0.9 | -0.1 | 0.3 | 1.2 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 3-Chlorotoluene | 3.2 -5.0 | 3.1 -2.1 | -1.9 | 3.8 | 0.6 | -1.8 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,3,5-Trimethylbenzene | 11.8 -11.7 | 10.1 -13.5 | 3.6 | 3.4 | 0.4 | -4.0 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 4-Chlorotoluene | 13.1 -5.4 | -1.4 -4.2 | -0.1 | -1.3 | -1.7 | 1.1 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| tert-Butylbenzene | 11.1 -11.4 | 5.1 -10.2 | 4.1 | 2.3 | 2.1 | -3.2 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,2,4-Trimethylbenzene | 17.3 -13.4 | 10.6 -15.1 | 3.7 | 1.4 | 0.5 | -5.0 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
READBACK PERCENT ERROR

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

Analy Batch No.: 217861

SDG No.: _____

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

| ANALYTE | PERCENT ERROR | | | | | | PERCENT ERROR LIMIT | | | | | |
|------------------------------------|--------------------|--------------------|---------|---------|---------|---------|---------------------|----------------|-------|-------|-------|-------|
| | LVL 1 # LVL 7 # | LVL 2 # LVL 8 # | LVL 3 # | LVL 4 # | LVL 5 # | LVL 6 # | LVL 1 LVL 7 | LVL 2 LVL 8 | LVL 3 | LVL 4 | LVL 5 | LVL 6 |
| 3,4-Dichlorobenzotrifluoride | 10.8 -6.8 | -1.0 2.1 | -6.7 | 4.5 | -0.9 | -2.0 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| sec-Butylbenzene | 24.6 -15.7 | 8.6 -15.9 | 3.1 | 2.1 | 0.4 | -7.1 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,3-Dichlorobenzene | 17.8 -9.2 | 7.5 -9.5 | -1.2 | -1.1 | -1.3 | -2.9 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 4-Isopropyltoluene | 16.5 -14.1 | 12.7 -13.5 | 2.2 | 1.7 | -0.6 | -4.9 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,4-Dichlorobenzene | 15.1 -9.8 | 7.9 -9.6 | 1.3 | -1.3 | -1.3 | -2.3 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2,4-Dichlorobenzotrifluoride | 3.1 -10.2 | 1.0 6.0 | -2.7 | 7.0 | -2.8 | -1.4 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2,5-Dichlorobenzotrifluoride | -2.6 -0.7 | 8.3 -2.1 | -5.6 | 2.2 | -1.2 | 1.7 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| n-Butylbenzene | 17.5 -13.9 | 9.9 -11.8 | 0.5 | 3.0 | 0.0 | -5.1 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,2-Dichlorobenzene | 11.6 -9.6 | 13.1 -8.4 | -1.3 | -1.3 | -2.7 | -1.4 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,2-Dibromo-3-Chloropropane | -14.1 10.8 | -1.7 18.0 | -10.5 | -6.3 | -9.7 | 13.5 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2,4- & 2,5- & 2,6- Dichlorotoluene | 6.0 -8.4 | 15.7 -4.7 | -3.1 | 2.7 | -2.8 | -5.3 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2,3- & 3,4- Dichlorotoluene | 2.8 -2.5 | 11.3 2.0 | -6.2 | -1.3 | -2.9 | -3.3 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,2,4-Trichlorobenzene | 0.2 2.3 | 12.8 8.0 | -9.5 | -7.3 | -5.0 | -1.3 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Hexachlorobutadiene | 25.9 -1.7 | 0.4 12.6 | -15.7 | -8.5 | -6.8 | -6.2 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Naphthalene | 22.7 -4.5 | 14.5 -5.2 | -6.5 | -7.0 | -6.9 | -7.2 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,2,3-Trichlorobenzene | 4.8 9.5 | 8.6 15.0 | -14.3 | -12.5 | -8.8 | -2.4 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2,4,5-Trichlorotoluene | -5.6 22.7 | -2.9 33.5 * | -20.5 | -19.3 | -7.1 | -0.8 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2,3,6-Trichlorotoluene | -5.1 18.7 | 3.0 30.3 * | -17.5 | -19.9 | -5.0 | -4.6 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Dibromofluoromethane (Surr) | 19.5 -6.8 | 3.9 -3.6 | -8.1 | -1.5 | -5.8 | 2.3 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,2-Dichloroethane-d4 (Surr) | 33.2 -12.1 | 10.8 -10.0 | -6.9 | -3.3 | -11.1 | -0.6 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Toluene-d8 (Surr) | -2.6 -11.7 | 12.5 +++++ | 0.9 | 6.5 | -0.6 | -5.0 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
READBACK PERCENT ERROR

Lab Name: TestAmerica Pittsburgh Job No.: 180-71131-1 Analy Batch No.: 217861

SDG No.: _____

Instrument ID: CHHP6 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 07/24/2017 06:39 Calibration End Date: 07/24/2017 09:28 Calibration ID: 35029

| ANALYTE | PERCENT ERROR | | | | | | PERCENT ERROR LIMIT | | | | | |
|-----------------------------|---------------|---------|---------|---------|---------|---------|---------------------|-------|-------|-------|-------|-------|
| | LVL 1 # | LVL 2 # | LVL 3 # | LVL 4 # | LVL 5 # | LVL 6 # | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 6 |
| 4-Bromofluorobenzene (Surr) | 38.5 | 13.2 | -5.5 | -0.8 | -7.9 | -4.5 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -14.8 | -18.3 | | | | | 30 | 30 | | | | |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D03.D
 Lims ID: IC VSTD1
 Client ID:
 Sample Type: IC Calib Level: 1
 Inject. Date: 24-Jul-2017 06:39:30 ALS Bottle#: 3 Worklist Smp#: 3
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017705-003
 Misc. Info.: IC VSTD1
 Operator ID: 034635 Instrument ID: CHHP6
 Sublist: chrom-MSVOA_LL_CHHP6*sub10
 Method: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 25-Jul-2017 01:44:30 Calib Date: 24-Jul-2017 09:28:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D10.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK021

First Level Reviewer: bungardf Date: 24-Jul-2017 07:18:57

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 3.968 | 3.973 | -0.005 | 93 | 341897 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.052 | 7.052 | 0.000 | 98 | 796580 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.173 | 10.172 | 0.001 | 88 | 172191 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.515 | 12.515 | 0.001 | 97 | 266961 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr) | 113 | 6.316 | 6.315 | 0.001 | 92 | 24732 | 5.00 | 5.97 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur) | 65 | 6.693 | 6.693 | 0.000 | 62 | 39405 | 5.00 | 6.66 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.719 | 8.712 | 0.007 | 93 | 102855 | 5.00 | 4.87 | |
| \$ 8 4-Bromofluorobenzene (Surr) | 95 | 11.353 | 11.353 | 0.000 | 84 | 41164 | 5.00 | 6.92 | |
| 11 Dichlorodifluoromethane | 85 | 1.486 | 1.479 | 0.007 | 98 | 27479 | 5.00 | 5.41 | |
| 12 Chloromethane | 50 | 1.632 | 1.631 | 0.001 | 97 | 23920 | 5.00 | 5.23 | |
| 13 Vinyl chloride | 62 | 1.747 | 1.753 | -0.006 | 96 | 26240 | 5.00 | 5.48 | |
| 14 Butadiene | 39 | 1.790 | 1.789 | 0.001 | 89 | 24629 | 5.00 | 6.20 | |
| 15 Bromomethane | 94 | 2.076 | 2.075 | 0.001 | 72 | 12374 | 5.00 | 5.54 | |
| 16 Chloroethane | 64 | 2.210 | 2.203 | 0.007 | 82 | 14369 | 5.00 | 5.55 | |
| 17 Dichlorofluoromethane | 67 | 2.459 | 2.459 | 0.000 | 95 | 29354 | 5.00 | 5.26 | |
| 18 Trichlorofluoromethane | 101 | 2.496 | 2.501 | -0.005 | 85 | 24407 | 5.00 | 5.17 | M |
| 20 Ethyl ether | 59 | 2.818 | 2.824 | -0.006 | 87 | 21539 | 5.00 | 5.33 | |
| 21 Acrolein | 56 | 2.989 | 2.994 | -0.005 | 100 | 93239 | 100.0 | 107.8 | |
| 22 1,1-Dichloroethene | 96 | 3.110 | 3.109 | 0.001 | 96 | 22840 | 5.00 | 5.52 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.153 | 3.158 | -0.005 | 88 | 20417 | 5.00 | 5.15 | |
| 24 Acetone | 43 | 3.183 | 3.195 | -0.012 | 99 | 47909 | 25.0 | 28.0 | |
| 25 Iodomethane | 142 | 3.287 | 3.292 | -0.005 | 92 | 29488 | 5.00 | 5.07 | |
| 26 Carbon disulfide | 76 | 3.360 | 3.365 | -0.005 | 98 | 41100 | 5.00 | 4.46 | |
| 29 3-Chloro-1-propene | 76 | 3.633 | 3.633 | 0.000 | 80 | 11154 | 5.00 | 4.60 | |
| 30 Methyl acetate | 43 | 3.646 | 3.651 | -0.005 | 95 | 40003 | 10.0 | 10.7 | |
| 31 Methylene Chloride | 84 | 3.828 | 3.846 | -0.018 | 93 | 34286 | 5.00 | 6.25 | |
| 32 2-Methyl-2-propanol | 59 | 4.108 | 4.113 | -0.005 | 89 | 20227 | 50.0 | 53.4 | |
| 33 Acrylonitrile | 53 | 4.236 | 4.235 | 0.001 | 99 | 114140 | 50.0 | 57.0 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.272 | 4.271 | 0.001 | 96 | 24529 | 5.00 | 5.22 | |
| 35 Methyl tert-butyl ether | 73 | 4.284 | 4.290 | -0.006 | 95 | 78478 | 5.00 | 5.22 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 36 Hexane | 57 | 4.704 | 4.709 | -0.005 | 88 | 30509 | 5.00 | 5.48 | |
| 37 1,1-Dichloroethane | 63 | 4.923 | 4.922 | 0.001 | 73 | 27271 | 5.00 | 3.57 | |
| 38 Vinyl acetate | 43 | 4.978 | 4.983 | -0.005 | 97 | 41136 | 5.00 | 4.62 | |
| 42 2,2-Dichloropropane | 97 | 5.683 | 5.689 | -0.006 | 51 | 3609 | 5.00 | 4.56 | |
| 43 cis-1,2-Dichloroethene | 96 | 5.690 | 5.695 | -0.005 | 75 | 26745 | 5.00 | 4.87 | |
| 44 2-Butanone (MEK) | 43 | 5.708 | 5.713 | -0.005 | 98 | 61711 | 25.0 | 25.2 | |
| 48 Chlorobromomethane | 128 | 5.982 | 5.981 | 0.001 | 98 | 12079 | 5.00 | 5.04 | |
| 49 Tetrahydrofuran | 42 | 5.994 | 5.987 | 0.007 | 91 | 22350 | 10.0 | 13.5 | |
| 50 Chloroform | 83 | 6.140 | 6.133 | 0.007 | 92 | 40199 | 5.00 | 4.88 | |
| 51 1,1,1-Trichloroethane | 97 | 6.280 | 6.285 | -0.005 | 95 | 24367 | 5.00 | 4.65 | |
| 52 Cyclohexane | 56 | 6.359 | 6.364 | -0.005 | 91 | 41233 | 5.00 | 5.54 | |
| 53 Carbon tetrachloride | 117 | 6.468 | 6.468 | 0.000 | 93 | 15662 | 5.00 | 4.16 | |
| 54 1,1-Dichloropropene | 75 | 6.480 | 6.486 | -0.006 | 96 | 31712 | 5.00 | 5.03 | |
| 56 Benzene | 78 | 6.693 | 6.699 | -0.006 | 96 | 100837 | 5.00 | 5.49 | |
| 55 Isobutyl alcohol | 41 | 6.693 | 7.076 | -0.383 | 37 | 13528 | 125.0 | 122.4 | |
| 57 1,2-Dichloroethane | 62 | 6.779 | 6.778 | 0.001 | 97 | 35160 | 5.00 | 5.18 | |
| 59 n-Heptane | 43 | 7.077 | 7.076 | 0.001 | 48 | 24265 | 5.00 | 5.73 | |
| 61 Trichloroethene | 130 | 7.442 | 7.447 | -0.005 | 98 | 22756 | 5.00 | 5.10 | |
| 63 Methylcyclohexane | 83 | 7.679 | 7.678 | 0.001 | 84 | 43857 | 5.00 | 5.63 | |
| 64 1,2-Dichloropropane | 63 | 7.715 | 7.715 | 0.000 | 89 | 23091 | 5.00 | 5.03 | |
| 65 1,4-Dioxane | 88 | 7.807 | 7.885 | -0.078 | 42 | 4766 | 100.0 | 113.9 | |
| 67 Dibromomethane | 93 | 7.801 | 7.806 | -0.005 | 93 | 15724 | 5.00 | 5.24 | |
| 68 Dichlorobromomethane | 83 | 8.001 | 8.007 | -0.006 | 94 | 20777 | 5.00 | 4.07 | |
| 70 2-Chloroethyl vinyl ether | 63 | 8.312 | 8.311 | 0.001 | 91 | 29761 | 10.0 | 9.78 | |
| 71 cis-1,3-Dichloropropene | 75 | 8.451 | 8.457 | -0.006 | 94 | 21547 | 5.00 | 3.72 | |
| 72 4-Methyl-2-pentanone (MIBK) | 43 | 8.616 | 8.615 | 0.001 | 94 | 113074 | 25.0 | 25.7 | |
| 73 Toluene | 91 | 8.786 | 8.779 | 0.007 | 99 | 104189 | 5.00 | 6.14 | |
| 74 trans-1,3-Dichloropropene | 75 | 9.035 | 9.035 | 0.000 | 93 | 17940 | 5.00 | 3.95 | |
| 75 Ethyl methacrylate | 69 | 9.102 | 9.102 | 0.000 | 90 | 25997 | 5.00 | 4.54 | |
| 76 1,1,2-Trichloroethane | 97 | 9.230 | 9.229 | 0.001 | 90 | 20294 | 5.00 | 5.17 | |
| 77 Tetrachloroethene | 164 | 9.291 | 9.296 | -0.005 | 96 | 18016 | 5.00 | 5.93 | |
| 78 1,3-Dichloropropane | 76 | 9.382 | 9.388 | -0.006 | 88 | 39343 | 5.00 | 5.53 | |
| 79 2-Hexanone | 43 | 9.449 | 9.448 | 0.001 | 95 | 72896 | 25.0 | 25.8 | |
| 81 Chlorodibromomethane | 129 | 9.607 | 9.601 | 0.006 | 90 | 10736 | 5.00 | 4.04 | |
| 82 Ethylene Dibromide | 107 | 9.705 | 9.710 | -0.005 | 97 | 18711 | 5.00 | 4.95 | |
| 83 3-Chlorobenzotrifluoride | 180 | 10.185 | 10.185 | 0.000 | 61 | 26101 | 5.00 | 5.29 | |
| 84 Chlorobenzene | 112 | 10.204 | 10.197 | 0.007 | 95 | 65919 | 5.00 | 5.97 | |
| 85 4-Chlorobenzotrifluoride | 180 | 10.270 | 10.270 | 0.000 | 95 | 24071 | 5.00 | 5.27 | |
| 86 1,1,1,2-Tetrachloroethane | 131 | 10.289 | 10.294 | -0.005 | 40 | 13142 | 5.00 | 4.19 | |
| 87 Ethylbenzene | 106 | 10.301 | 10.300 | 0.001 | 98 | 35398 | 5.00 | 5.61 | |
| 88 m-Xylene & p-Xylene | 106 | 10.435 | 10.434 | 0.001 | 99 | 43910 | 5.00 | 5.68 | |
| 89 o-Xylene | 106 | 10.812 | 10.811 | 0.001 | 97 | 42778 | 5.00 | 5.52 | |
| 90 Styrene | 104 | 10.836 | 10.835 | 0.001 | 95 | 70410 | 5.00 | 5.62 | |
| 91 Bromoform | 173 | 11.007 | 11.018 | -0.011 | 82 | 5803 | 5.00 | 3.88 | |
| 92 2-Chlorobenzotrifluoride | 180 | 11.092 | 11.091 | 0.001 | 94 | 25025 | 5.00 | 5.01 | |
| 93 Isopropylbenzene | 105 | 11.183 | 11.182 | 0.001 | 95 | 110174 | 5.00 | 6.35 | |
| 95 Bromobenzene | 156 | 11.493 | 11.492 | 0.001 | 94 | 25122 | 5.00 | 5.06 | |
| 96 1,1,2,2-Tetrachloroethane | 83 | 11.499 | 11.499 | 0.000 | 94 | 29044 | 5.00 | 5.26 | |
| 97 trans-1,4-Dichloro-2-butene | 53 | 11.530 | 11.535 | -0.005 | 70 | 7272 | 5.00 | 5.01 | |
| 98 1,2,3-Trichloropropane | 110 | 11.548 | 11.553 | -0.005 | 88 | 10261 | 5.00 | 4.98 | |
| 99 N-Propylbenzene | 120 | 11.597 | 11.596 | 0.001 | 99 | 29088 | 5.00 | 5.39 | |
| 100 2-Chlorotoluene | 126 | 11.682 | 11.681 | 0.001 | 95 | 24679 | 5.00 | 5.28 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 101 3-Chlorotoluene | 126 | 11.749 | 11.748 | 0.001 | 97 | 25329 | 5.00 | 5.16 | |
| 102 1,3,5-Trimethylbenzene | 105 | 11.785 | 11.784 | 0.001 | 92 | 85688 | 5.00 | 5.59 | |
| 103 4-Chlorotoluene | 126 | 11.810 | 11.809 | 0.001 | 97 | 28869 | 5.00 | 5.65 | |
| 104 tert-Butylbenzene | 119 | 12.095 | 12.095 | 0.000 | 90 | 67600 | 5.00 | 5.55 | |
| 106 1,2,4-Trimethylbenzene | 105 | 12.156 | 12.156 | 0.000 | 95 | 93775 | 5.00 | 5.87 | |
| 107 1,2-dichloro-4-(trifluoromethyl) | 214 | 12.199 | 12.204 | -0.005 | 96 | 19655 | 5.00 | 5.54 | |
| 108 sec-Butylbenzene | 105 | 12.321 | 12.320 | 0.001 | 94 | 106349 | 5.00 | 6.23 | |
| 109 1,3-Dichlorobenzene | 146 | 12.436 | 12.435 | 0.001 | 95 | 52474 | 5.00 | 5.89 | |
| 110 4-Isopropyltoluene | 119 | 12.473 | 12.478 | -0.005 | 95 | 82129 | 5.00 | 5.82 | |
| 111 1,4-Dichlorobenzene | 146 | 12.533 | 12.539 | -0.006 | 93 | 53033 | 5.00 | 5.76 | |
| 113 2,4-Dichloro-1-(trifluoromethyl) | 214 | 12.570 | 12.569 | 0.001 | 94 | 17229 | 5.00 | 5.15 | |
| 114 2,5-Dichlorobenzotrifluoride | 214 | 12.613 | 12.612 | 0.001 | 96 | 17972 | 5.00 | 4.87 | |
| 116 n-Butylbenzene | 91 | 12.880 | 12.886 | -0.006 | 98 | 75545 | 5.00 | 5.87 | |
| 117 1,2-Dichlorobenzene | 146 | 12.892 | 12.892 | 0.000 | 94 | 47273 | 5.00 | 5.58 | |
| 118 1,2-Dibromo-3-Chloropropan | 75 | 13.677 | 13.683 | -0.006 | 66 | 3245 | 5.00 | 4.30 | |
| 119 2,4- & 2,5- & 2,6- Dichlorobenzene | 125 | 13.823 | 13.816 | 0.007 | 99 | 85812 | 15.0 | 15.9 | |
| 121 2,3- & 3,4- Dichlorotoluene | 125 | 14.237 | 14.236 | 0.001 | 98 | 60616 | 10.0 | 10.3 | |
| 122 1,2,4-Trichlorobenzene | 180 | 14.505 | 14.498 | 0.007 | 92 | 22175 | 5.00 | 5.01 | |
| 123 Hexachlorobutadiene | 225 | 14.644 | 14.650 | -0.006 | 89 | 8829 | 5.00 | 6.30 | |
| 124 Naphthalene | 128 | 14.760 | 14.759 | 0.001 | 97 | 82809 | 5.00 | 6.13 | |
| 125 1,2,3-Trichlorobenzene | 180 | 14.985 | 14.984 | 0.001 | 94 | 20833 | 5.00 | 5.24 | |
| 126 2,4,5-Trichlorotoluene | 159 | 15.776 | 15.775 | 0.001 | 0 | 10792 | 5.00 | 4.72 | |
| 127 2,3,6-Trichlorotoluene | 159 | 15.879 | 15.885 | -0.006 | 91 | 9731 | 5.00 | 4.75 | |
| 146 3,4-Dichlorotoluene | 1 | 0.000 | | | | | ND | ND | |
| S 131 Xylenes, Total | 106 | | | | 0 | | 10.0 | 11.2 | |
| S 130 1,2-Dichloroethene, Total | 96 | | | | 0 | | 10.0 | 10.1 | |
| S 132 1,3-Dichloropropene, Total | 1 | | | | 0 | | 10.0 | 7.66 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

Reagents:

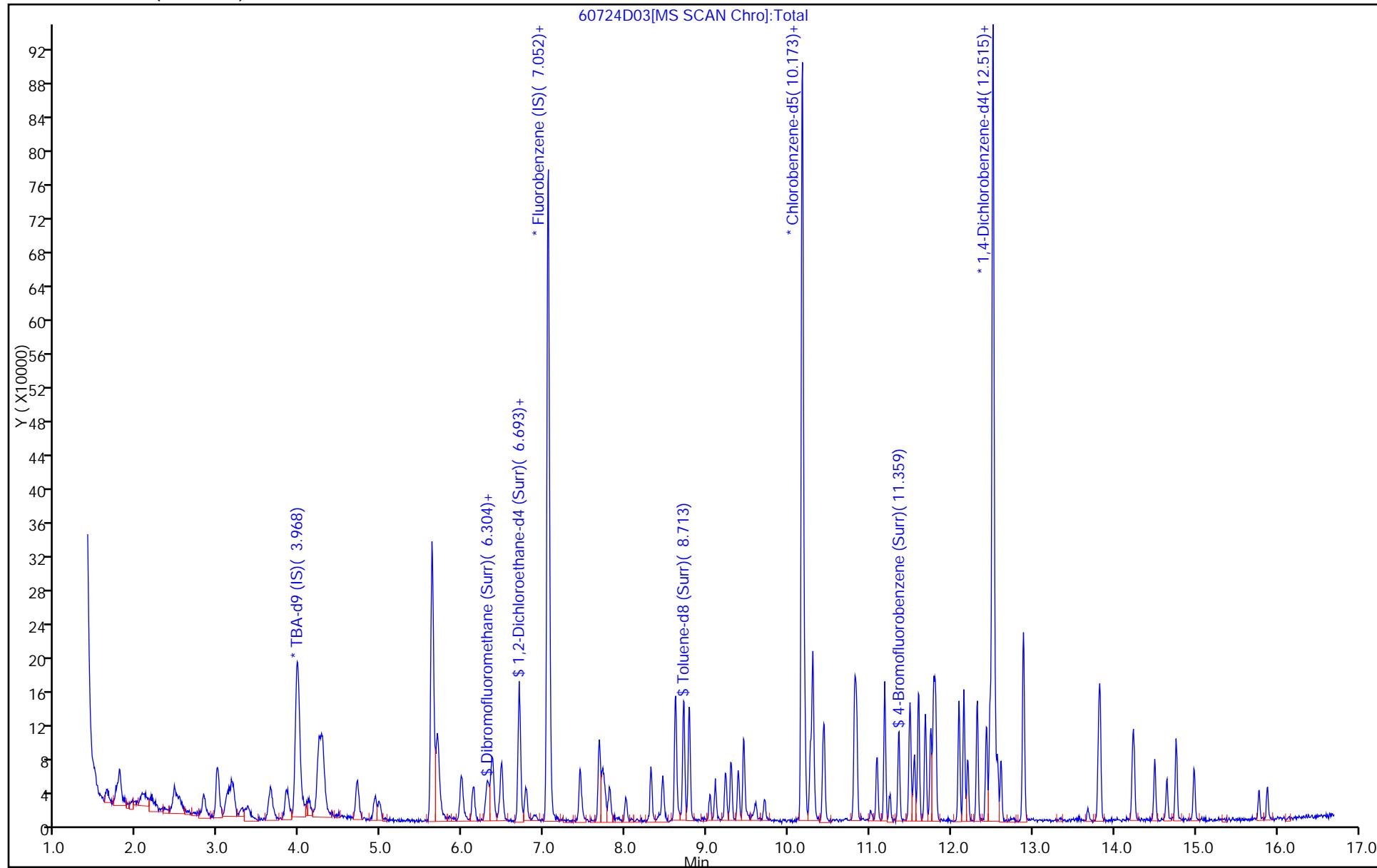
| | | |
|---------------------|--------------------|-----------|
| voaWEEmix1stR_00009 | Amount Added: 0.20 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 0.80 | Units: uL |
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 0.20 | Units: uL |
| voaWVA1stRest_00016 | Amount Added: 0.20 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 4.00 | Units: uL |
| VOA8260VOAPRI_00263 | Amount Added: 0.20 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 0.20 | Units: uL |

Report Date: 25-Jul-2017 01:44:31

Chrom Revision: 2.2 20-Jun-2017 07:42:38

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D03.D
Injection Date: 24-Jul-2017 06:39:30 Instrument ID: CHHP6
Lims ID: IC VSTD1 Operator ID: 034635
Client ID:
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 3
Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh

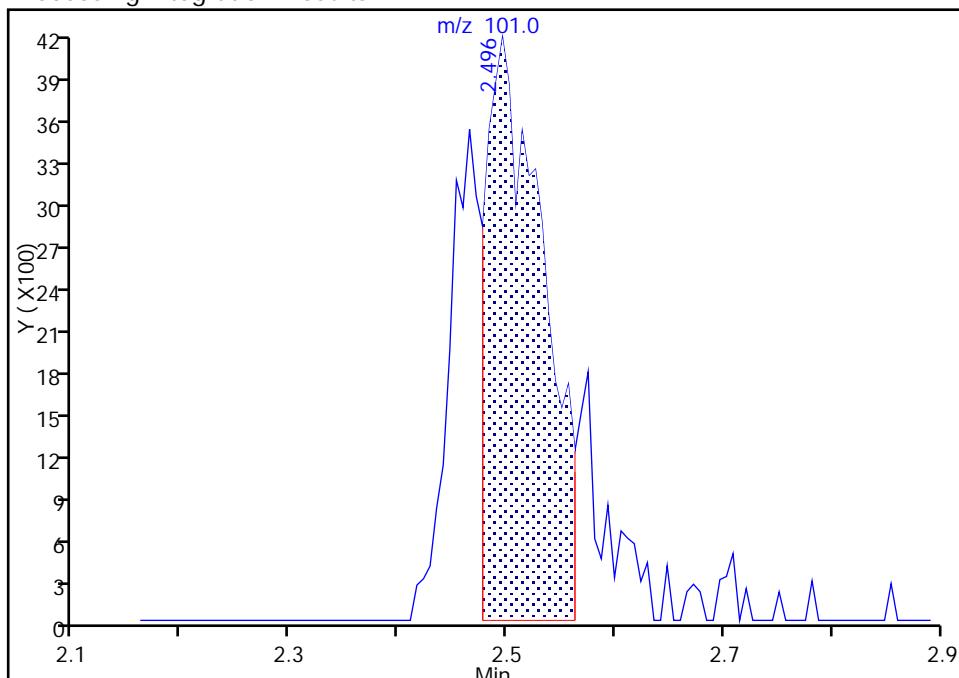
Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D03.D
 Injection Date: 24-Jul-2017 06:39:30 Instrument ID: CHHP6
 Lims ID: IC VSTD1
 Client ID:
 Operator ID: 034635 ALS Bottle#: 3 Worklist Smp#: 3
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
 Column: DB-624 (0.18 mm) Detector: MS SCAN

18 Trichlorofluoromethane, CAS: 75-69-4

Signal: 1

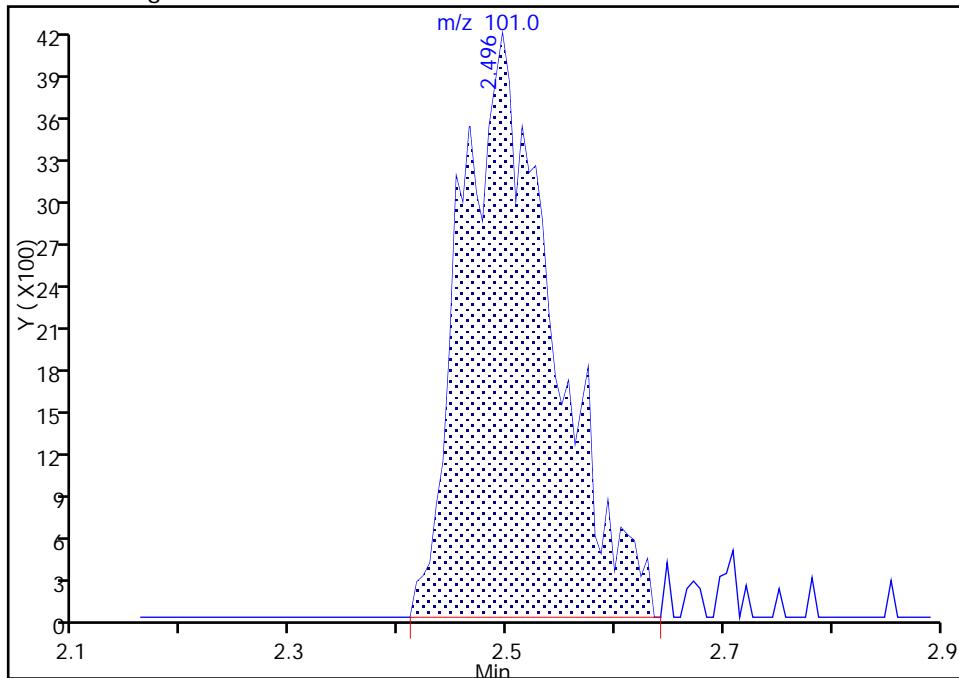
RT: 2.50
 Area: 15241
 Amount: 5.000000
 Amount Units: ng

Processing Integration Results



RT: 2.50
 Area: 24407
 Amount: 5.173113
 Amount Units: ng

Manual Integration Results



Reviewer: bungardf, 24-Jul-2017 07:17:36

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D04.D
 Lims ID: IC VSTD5
 Client ID:
 Sample Type: IC Calib Level: 2
 Inject. Date: 24-Jul-2017 07:03:30 ALS Bottle#: 4 Worklist Smp#: 4
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017705-004
 Misc. Info.: IC VSTD5
 Operator ID: 034635 Instrument ID: CHHP6
 Sublist: chrom-MSVOA_LL_CHHP6*sub10
 Method: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 25-Jul-2017 01:44:32 Calib Date: 24-Jul-2017 09:28:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D10.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK021

First Level Reviewer: bungardf Date: 24-Jul-2017 07:25:58

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 3.968 | 3.973 | -0.005 | 93 | 269256 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.052 | 7.052 | 0.000 | 98 | 704642 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.173 | 10.172 | 0.001 | 88 | 162034 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.509 | 12.515 | -0.005 | 97 | 250761 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr) | 113 | 6.316 | 6.315 | 0.001 | 93 | 95101 | 25.0 | 26.0 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur) | 65 | 6.693 | 6.693 | 0.000 | 59 | 145003 | 25.0 | 27.7 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.713 | 8.712 | 0.001 | 93 | 381639 | 25.0 | 28.1 | |
| \$ 8 4-Bromofluorobenzene (Surr) | 95 | 11.353 | 11.353 | 0.000 | 83 | 158379 | 25.0 | 28.3 | |
| 11 Dichlorodifluoromethane | 85 | 1.480 | 1.479 | 0.001 | 98 | 121547 | 25.0 | 27.0 | |
| 12 Chloromethane | 50 | 1.626 | 1.631 | -0.005 | 99 | 109949 | 25.0 | 27.2 | |
| 13 Vinyl chloride | 62 | 1.753 | 1.753 | 0.000 | 97 | 112168 | 25.0 | 26.5 | |
| 14 Butadiene | 39 | 1.790 | 1.789 | 0.001 | 89 | 92248 | 25.0 | 26.3 | |
| 15 Bromomethane | 94 | 2.082 | 2.075 | 0.007 | 88 | 58670 | 25.0 | 29.7 | |
| 16 Chloroethane | 64 | 2.204 | 2.203 | 0.001 | 98 | 64213 | 25.0 | 28.0 | |
| 17 Dichlorofluoromethane | 67 | 2.459 | 2.459 | 0.000 | 97 | 133871 | 25.0 | 27.1 | |
| 18 Trichlorofluoromethane | 101 | 2.489 | 2.501 | -0.012 | 98 | 108123 | 25.0 | 25.9 | |
| 20 Ethyl ether | 59 | 2.824 | 2.824 | 0.000 | 88 | 103875 | 25.0 | 29.1 | |
| 21 Acrolein | 56 | 3.000 | 2.994 | 0.006 | 98 | 103236 | 125.0 | 134.9 | |
| 22 1,1-Dichloroethene | 96 | 3.104 | 3.109 | -0.005 | 98 | 90501 | 25.0 | 24.7 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.159 | 3.158 | 0.001 | 94 | 89949 | 25.0 | 25.6 | |
| 24 Acetone | 43 | 3.183 | 3.195 | -0.012 | 100 | 83742 | 50.0 | 55.4 | |
| 25 Iodomethane | 142 | 3.286 | 3.292 | -0.006 | 96 | 133154 | 25.0 | 25.9 | |
| 26 Carbon disulfide | 76 | 3.372 | 3.365 | 0.007 | 98 | 179135 | 25.0 | 22.0 | |
| 29 3-Chloro-1-propene | 76 | 3.639 | 3.633 | 0.006 | 90 | 50272 | 25.0 | 23.4 | |
| 30 Methyl acetate | 43 | 3.651 | 3.651 | 0.000 | 96 | 185280 | 50.0 | 55.9 | |
| 31 Methylene Chloride | 84 | 3.846 | 3.846 | 0.000 | 90 | 131057 | 25.0 | 27.0 | |
| 32 2-Methyl-2-propanol | 59 | 4.108 | 4.113 | -0.005 | 91 | 71644 | 250.0 | 240.4 | |
| 33 Acrylonitrile | 53 | 4.235 | 4.235 | 0.000 | 100 | 530059 | 250.0 | 299.5 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.272 | 4.271 | 0.001 | 96 | 108647 | 25.0 | 26.1 | |
| 35 Methyl tert-butyl ether | 73 | 4.284 | 4.290 | -0.006 | 96 | 375187 | 25.0 | 28.2 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 36 Hexane | 57 | 4.704 | 4.709 | -0.005 | 90 | 127818 | 25.0 | 26.0 | |
| 37 1,1-Dichloroethane | 63 | 4.929 | 4.922 | 0.007 | 96 | 187923 | 25.0 | 27.8 | |
| 38 Vinyl acetate | 43 | 4.984 | 4.983 | 0.001 | 97 | 217747 | 25.0 | 27.7 | |
| 42 2,2-Dichloropropane | 97 | 5.683 | 5.689 | -0.006 | 54 | 17106 | 25.0 | 24.4 | |
| 43 cis-1,2-Dichloroethene | 96 | 5.689 | 5.695 | -0.006 | 80 | 127777 | 25.0 | 26.3 | |
| 44 2-Butanone (MEK) | 43 | 5.702 | 5.713 | -0.011 | 73 | 125277 | 50.0 | 57.8 | |
| 48 Chlorobromomethane | 128 | 5.981 | 5.981 | 0.000 | 98 | 57129 | 25.0 | 27.0 | |
| 49 Tetrahydrofuran | 42 | 5.994 | 5.987 | 0.007 | 87 | 85090 | 50.0 | 58.0 | |
| 50 Chloroform | 83 | 6.133 | 6.133 | 0.000 | 93 | 190538 | 25.0 | 26.1 | |
| 51 1,1,1-Trichloroethane | 97 | 6.292 | 6.285 | 0.007 | 97 | 114539 | 25.0 | 24.7 | |
| 52 Cyclohexane | 56 | 6.359 | 6.364 | -0.005 | 91 | 170750 | 25.0 | 25.9 | |
| 53 Carbon tetrachloride | 117 | 6.468 | 6.468 | 0.000 | 95 | 77574 | 25.0 | 23.3 | |
| 54 1,1-Dichloropropene | 75 | 6.480 | 6.486 | -0.006 | 96 | 141227 | 25.0 | 25.3 | |
| 56 Benzene | 78 | 6.699 | 6.699 | 0.000 | 97 | 461882 | 25.0 | 28.4 | |
| 57 1,2-Dichloroethane | 62 | 6.778 | 6.778 | 0.000 | 98 | 171694 | 25.0 | 28.6 | |
| 55 Isobutyl alcohol | 41 | 6.699 | 7.076 | -0.377 | 40 | 64845 | 625.0 | 663.4 | |
| 59 n-Heptane | 43 | 7.076 | 7.076 | 0.000 | 80 | 95666 | 25.0 | 25.5 | |
| 61 Trichloroethene | 130 | 7.447 | 7.447 | 0.000 | 98 | 102011 | 25.0 | 25.9 | |
| 63 Methylcyclohexane | 83 | 7.679 | 7.678 | 0.001 | 87 | 179050 | 25.0 | 26.0 | |
| 64 1,2-Dichloropropane | 63 | 7.715 | 7.715 | 0.000 | 92 | 103896 | 25.0 | 25.6 | |
| 67 Dibromomethane | 93 | 7.806 | 7.806 | 0.000 | 96 | 70312 | 25.0 | 26.5 | |
| 65 1,4-Dioxane | 88 | 7.800 | 7.885 | -0.085 | 38 | 23040 | 500.0 | 622.3 | M |
| 68 Dichlorobromomethane | 83 | 8.007 | 8.007 | 0.000 | 98 | 109927 | 25.0 | 24.3 | |
| 70 2-Chloroethyl vinyl ether | 63 | 8.317 | 8.311 | 0.006 | 92 | 152667 | 50.0 | 56.7 | |
| 71 cis-1,3-Dichloropropene | 75 | 8.457 | 8.457 | 0.000 | 95 | 116838 | 25.0 | 22.8 | |
| 72 4-Methyl-2-pentanone (MIBK) | 43 | 8.615 | 8.615 | 0.000 | 96 | 233783 | 50.0 | 56.5 | |
| 73 Toluene | 91 | 8.786 | 8.779 | 0.007 | 99 | 450260 | 25.0 | 28.2 | |
| 74 trans-1,3-Dichloropropene | 75 | 9.035 | 9.035 | 0.000 | 92 | 98174 | 25.0 | 23.0 | |
| 75 Ethyl methacrylate | 69 | 9.102 | 9.102 | 0.000 | 87 | 142004 | 25.0 | 26.4 | |
| 76 1,1,2-Trichloroethane | 97 | 9.230 | 9.229 | 0.001 | 92 | 103854 | 25.0 | 28.1 | |
| 77 Tetrachloroethene | 164 | 9.297 | 9.296 | 0.001 | 96 | 72287 | 25.0 | 25.3 | |
| 78 1,3-Dichloropropane | 76 | 9.388 | 9.388 | 0.000 | 88 | 192724 | 25.0 | 28.8 | |
| 79 2-Hexanone | 43 | 9.449 | 9.448 | 0.001 | 95 | 154768 | 50.0 | 58.1 | |
| 81 Chlorodibromomethane | 129 | 9.601 | 9.601 | 0.000 | 87 | 57240 | 25.0 | 22.9 | |
| 82 Ethylene Dibromide | 107 | 9.711 | 9.710 | 0.000 | 98 | 94851 | 25.0 | 26.7 | |
| 83 3-Chlorobenzotrifluoride | 180 | 10.185 | 10.185 | 0.000 | 91 | 124733 | 25.0 | 26.9 | |
| 84 Chlorobenzene | 112 | 10.203 | 10.197 | 0.006 | 94 | 294577 | 25.0 | 28.4 | |
| 85 4-Chlorobenzotrifluoride | 180 | 10.270 | 10.270 | 0.000 | 96 | 110312 | 25.0 | 25.6 | |
| 86 1,1,1,2-Tetrachloroethane | 131 | 10.295 | 10.294 | 0.001 | 87 | 69755 | 25.0 | 23.6 | |
| 87 Ethylbenzene | 106 | 10.301 | 10.300 | 0.001 | 98 | 156900 | 25.0 | 26.4 | |
| 88 m-Xylene & p-Xylene | 106 | 10.434 | 10.434 | 0.000 | 99 | 196132 | 25.0 | 26.9 | |
| 89 o-Xylene | 106 | 10.818 | 10.811 | 0.007 | 96 | 201906 | 25.0 | 27.7 | |
| 90 Styrene | 104 | 10.836 | 10.835 | 0.001 | 94 | 331761 | 25.0 | 28.1 | |
| 91 Bromoform | 173 | 11.012 | 11.018 | -0.006 | 94 | 30696 | 25.0 | 21.8 | |
| 92 2-Chlorobenzotrifluoride | 180 | 11.091 | 11.091 | 0.000 | 97 | 130865 | 25.0 | 27.8 | |
| 93 Isopropylbenzene | 105 | 11.183 | 11.182 | 0.001 | 96 | 473843 | 25.0 | 29.0 | |
| 95 Bromobenzene | 156 | 11.487 | 11.492 | -0.005 | 97 | 122999 | 25.0 | 26.4 | |
| 96 1,1,2,2-Tetrachloroethane | 83 | 11.499 | 11.499 | 0.000 | 95 | 152313 | 25.0 | 29.3 | |
| 97 trans-1,4-Dichloro-2-butene | 53 | 11.536 | 11.535 | 0.001 | 69 | 35451 | 25.0 | 26.0 | |
| 98 1,2,3-Trichloropropane | 110 | 11.548 | 11.553 | -0.005 | 86 | 52406 | 25.0 | 27.1 | |
| 99 N-Propylbenzene | 120 | 11.596 | 11.596 | 0.000 | 98 | 130443 | 25.0 | 25.7 | |
| 100 2-Chlorotoluene | 126 | 11.682 | 11.681 | 0.001 | 95 | 111064 | 25.0 | 25.3 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 101 3-Chlorotoluene | 126 | 11.748 | 11.748 | 0.000 | 97 | 118876 | 25.0 | 25.8 | |
| 102 1,3,5-Trimethylbenzene | 105 | 11.785 | 11.784 | 0.001 | 93 | 396334 | 25.0 | 27.5 | |
| 103 4-Chlorotoluene | 126 | 11.809 | 11.809 | 0.000 | 99 | 118288 | 25.0 | 24.7 | |
| 104 tert-Butylbenzene | 119 | 12.095 | 12.095 | 0.000 | 91 | 300432 | 25.0 | 26.3 | |
| 106 1,2,4-Trimethylbenzene | 105 | 12.156 | 12.156 | 0.000 | 98 | 415297 | 25.0 | 27.7 | |
| 107 1,2-dichloro-4-(trifluoromethyl) | 214 | 12.205 | 12.204 | 0.001 | 96 | 82510 | 25.0 | 24.8 | |
| 108 sec-Butylbenzene | 105 | 12.320 | 12.320 | 0.000 | 94 | 435324 | 25.0 | 27.1 | |
| 109 1,3-Dichlorobenzene | 146 | 12.430 | 12.435 | -0.005 | 96 | 224858 | 25.0 | 26.9 | |
| 110 4-Isopropyltoluene | 119 | 12.478 | 12.478 | 0.000 | 96 | 373177 | 25.0 | 28.2 | |
| 111 1,4-Dichlorobenzene | 146 | 12.539 | 12.539 | 0.000 | 95 | 233465 | 25.0 | 27.0 | |
| 113 2,4-Dichloro-1-(trifluoromethyl) | 214 | 12.570 | 12.569 | 0.001 | 97 | 79327 | 25.0 | 25.3 | |
| 114 2,5-Dichlorobenzotrifluoride | 214 | 12.612 | 12.612 | 0.000 | 98 | 93800 | 25.0 | 27.1 | |
| 116 n-Butylbenzene | 91 | 12.886 | 12.886 | 0.000 | 98 | 331802 | 25.0 | 27.5 | |
| 117 1,2-Dichlorobenzene | 146 | 12.892 | 12.892 | 0.000 | 96 | 225127 | 25.0 | 28.3 | |
| 118 1,2-Dibromo-3-Chloropropan | 75 | 13.677 | 13.683 | -0.006 | 73 | 17440 | 25.0 | 24.6 | |
| 119 2,4- & 2,5- & 2,6- Dichlorobenzene | 125 | 13.823 | 13.816 | 0.007 | 99 | 439793 | 75.0 | 86.7 | |
| 121 2,3- & 3,4- Dichlorotoluene | 125 | 14.237 | 14.236 | 0.001 | 99 | 308264 | 50.0 | 55.7 | |
| 122 1,2,4-Trichlorobenzene | 180 | 14.498 | 14.498 | 0.000 | 95 | 117288 | 25.0 | 28.2 | |
| 123 Hexachlorobutadiene | 225 | 14.644 | 14.650 | -0.006 | 93 | 33039 | 25.0 | 25.1 | |
| 124 Naphthalene | 128 | 14.760 | 14.759 | 0.001 | 98 | 362983 | 25.0 | 28.6 | |
| 125 1,2,3-Trichlorobenzene | 180 | 14.985 | 14.984 | 0.001 | 96 | 101393 | 25.0 | 27.1 | |
| 126 2,4,5-Trichlorotoluene | 159 | 15.776 | 15.775 | 0.001 | 0 | 52163 | 25.0 | 24.3 | |
| 127 2,3,6-Trichlorotoluene | 159 | 15.879 | 15.885 | -0.006 | 98 | 49582 | 25.0 | 25.8 | |
| 146 3,4-Dichlorotoluene | 1 | 0.000 | | | | | ND | ND | |
| S 130 1,2-Dichloroethene, Total | 96 | | | | 0 | | 50.0 | 52.5 | |
| S 131 Xylenes, Total | 106 | | | | 0 | | 50.0 | 54.6 | |
| S 132 1,3-Dichloropropene, Total | 1 | | | | 0 | | 50.0 | 45.7 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

Reagents:

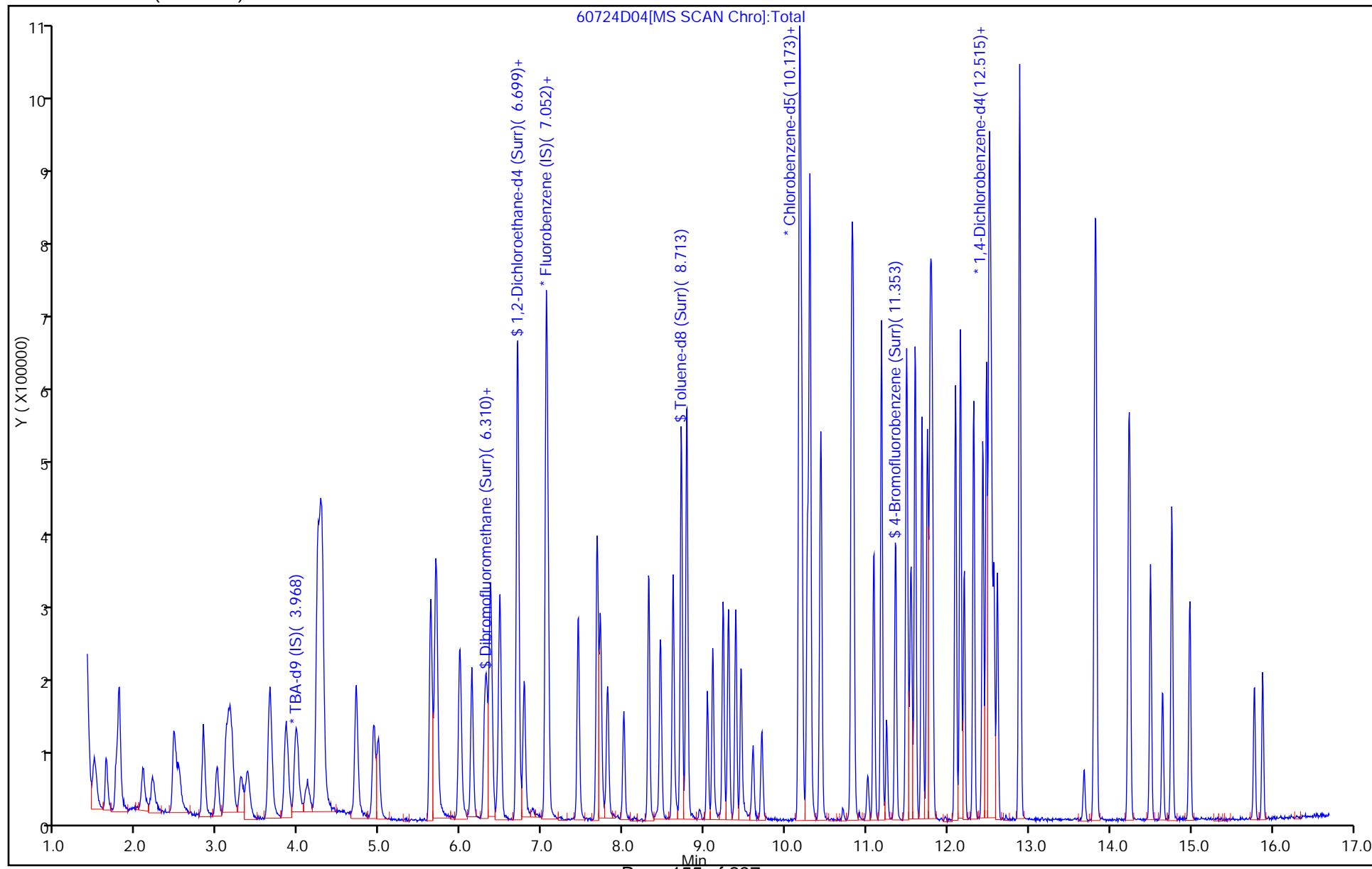
| | | |
|---------------------|--------------------|-----------|
| voaWEEmix1stR_00009 | Amount Added: 1.00 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 1.00 | Units: uL |
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 1.00 | Units: uL |
| voaWVA1stRest_00016 | Amount Added: 1.00 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 5.00 | Units: uL |
| VOA8260VOAPRI_00263 | Amount Added: 1.00 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 1.00 | Units: uL |

Report Date: 25-Jul-2017 01:44:33

Chrom Revision: 2.2 20-Jun-2017 07:42:38

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D04.D
Injection Date: 24-Jul-2017 07:03:30 Instrument ID: CHHP6
Lims ID: IC VSTD5 Operator ID: 034635
Client ID:
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 4
Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh

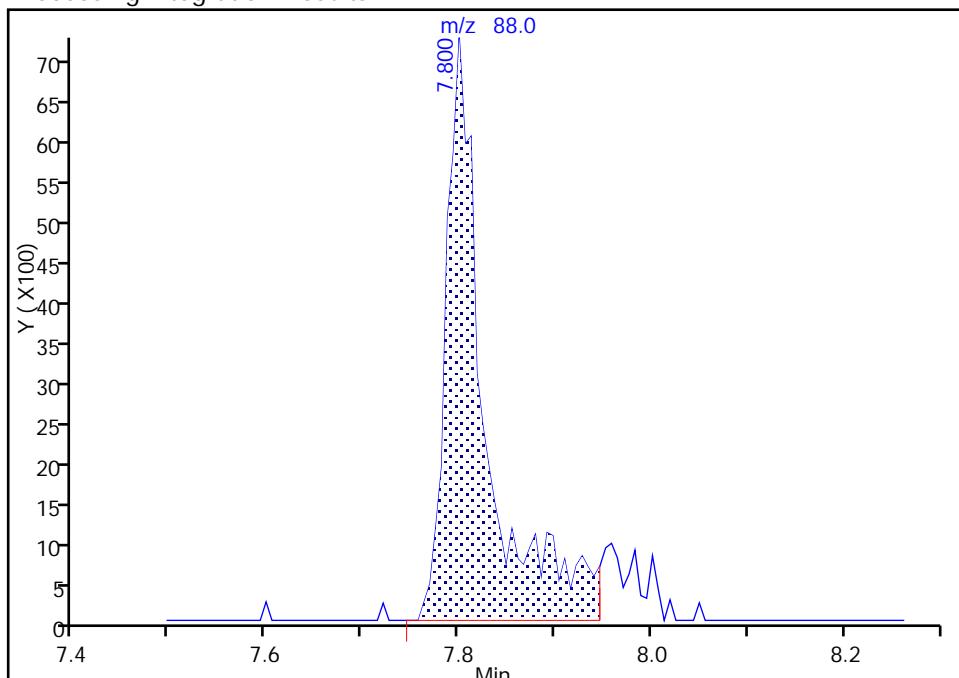
Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D04.D
 Injection Date: 24-Jul-2017 07:03:30 Instrument ID: CHHP6
 Lims ID: IC VSTD5
 Client ID:
 Operator ID: 034635 ALS Bottle#: 4 Worklist Smp#: 4
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
 Column: DB-624 (0.18 mm) Detector: MS SCAN

65 1,4-Dioxane, CAS: 123-91-1

Signal: 1

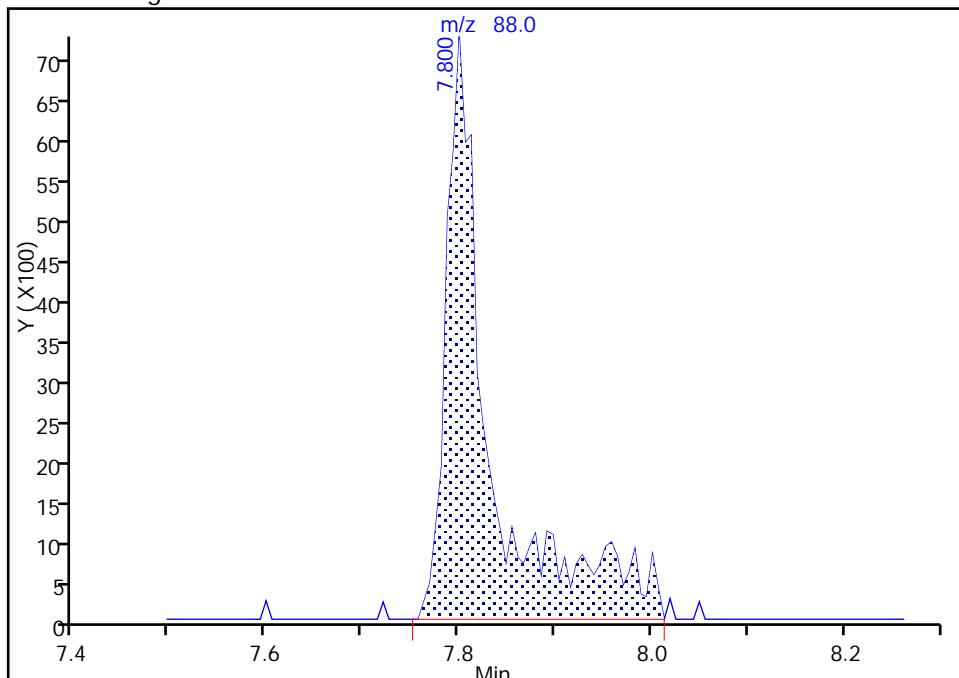
RT: 7.80
 Area: 20744
 Amount: 495.9874
 Amount Units: ng

Processing Integration Results



RT: 7.80
 Area: 23040
 Amount: 622.2830
 Amount Units: ng

Manual Integration Results



Reviewer: bungardf, 24-Jul-2017 07:27:10

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D05.D
 Lims ID: ICIS VSTD10
 Client ID:
 Sample Type: ICIS Calib Level: 3
 Inject. Date: 24-Jul-2017 07:27:30 ALS Bottle#: 5 Worklist Smp#: 5
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017705-005
 Misc. Info.: ICIS VSTD10
 Operator ID: 034635 Instrument ID: CHHP6
 Sublist: chrom-MSVOA_LL_CHHP6*sub10
 Method: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 25-Jul-2017 02:13:16 Calib Date: 24-Jul-2017 09:28:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D10.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK021

First Level Reviewer: bungardf Date: 25-Jul-2017 01:51:12

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 3.978 | 3.978 | 0.000 | 95 | 281180 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.050 | 7.050 | 0.000 | 99 | 854988 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.171 | 10.171 | 0.000 | 87 | 187443 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.513 | 12.513 | 0.000 | 94 | 265638 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr) | 113 | 6.314 | 6.314 | 0.000 | 93 | 204197 | 50.0 | 46.0 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur) | 65 | 6.691 | 6.691 | 0.000 | 69 | 295614 | 50.0 | 46.6 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.717 | 8.717 | 0.000 | 92 | 758339 | 50.0 | 50.5 | |
| \$ 8 4-Bromofluorobenzene (Surr) | 95 | 11.357 | 11.357 | 0.000 | 84 | 305961 | 50.0 | 47.3 | |
| 11 Dichlorodifluoromethane | 85 | 1.484 | 1.484 | 0.000 | 99 | 286699 | 50.0 | 52.5 | |
| 12 Chloromethane | 50 | 1.630 | 1.630 | 0.000 | 100 | 249199 | 50.0 | 50.8 | |
| 13 Vinyl chloride | 62 | 1.758 | 1.758 | 0.000 | 98 | 262771 | 50.0 | 51.1 | |
| 14 Butadiene | 39 | 1.788 | 1.788 | 0.000 | 92 | 223993 | 50.0 | 52.5 | |
| 15 Bromomethane | 94 | 2.086 | 2.086 | 0.000 | 87 | 122895 | 50.0 | 51.3 | |
| 16 Chloroethane | 64 | 2.202 | 2.202 | 0.000 | 98 | 137952 | 50.0 | 49.6 | |
| 17 Dichlorofluoromethane | 67 | 2.463 | 2.463 | 0.000 | 96 | 297904 | 50.0 | 49.7 | |
| 18 Trichlorofluoromethane | 101 | 2.506 | 2.506 | 0.000 | 95 | 244680 | 50.0 | 48.3 | |
| 20 Ethyl ether | 59 | 2.822 | 2.822 | 0.000 | 88 | 207890 | 50.0 | 48.0 | |
| 21 Acrolein | 56 | 2.993 | 2.993 | 0.000 | 100 | 126353 | 150.0 | 136.1 | |
| 22 1,1-Dichloroethene | 96 | 3.114 | 3.114 | 0.000 | 98 | 212019 | 50.0 | 47.7 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.157 | 3.157 | 0.000 | 94 | 214148 | 50.0 | 50.3 | |
| 24 Acetone | 43 | 3.187 | 3.187 | 0.000 | 100 | 181114 | 100.0 | 98.7 | |
| 25 Iodomethane | 142 | 3.291 | 3.291 | 0.000 | 99 | 296892 | 50.0 | 47.5 | |
| 26 Carbon disulfide | 76 | 3.370 | 3.370 | 0.000 | 98 | 451961 | 50.0 | 45.7 | |
| 29 3-Chloro-1-propene | 76 | 3.644 | 3.644 | 0.000 | 92 | 120354 | 50.0 | 46.3 | |
| 30 Methyl acetate | 43 | 3.656 | 3.656 | 0.000 | 96 | 382502 | 100.0 | 95.1 | |
| 31 Methylene Chloride | 84 | 3.844 | 3.844 | 0.000 | 89 | 278118 | 50.0 | 47.2 | |
| 32 2-Methyl-2-propanol | 59 | 4.112 | 4.112 | 0.000 | 93 | 151604 | 500.0 | 487.1 | |
| 33 Acrylonitrile | 53 | 4.234 | 4.234 | 0.000 | 100 | 1020211 | 500.0 | 475.0 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.276 | 4.276 | 0.000 | 99 | 244707 | 50.0 | 48.5 | |
| 35 Methyl tert-butyl ether | 73 | 4.295 | 4.295 | 0.000 | 95 | 782498 | 50.0 | 48.5 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|--------------|------------------|------------------|-----|----------|---------------|-----------------|-------|
| 36 Hexane | 57 | 4.708 | 4.708 | 0.000 | 91 | 304164 | 50.0 | 50.9 | |
| 37 1,1-Dichloroethane | 63 | 4.933 | 4.933 | 0.000 | 97 | 418314 | 50.0 | 51.0 | |
| 38 Vinyl acetate | 43 | 4.982 | 4.982 | 0.000 | 97 | 435645 | 50.0 | 45.6 | |
| 42 2,2-Dichloropropane | 97 | 5.688 | 5.688 | 0.000 | 61 | 39797 | 50.0 | 46.8 | |
| 43 cis-1,2-Dichloroethene | 96 | 5.694 | 5.694 | 0.000 | 79 | 290355 | 50.0 | 49.3 | |
| 44 2-Butanone (MEK) | 43 | 5.712 | 5.712 | 0.000 | 99 | 266460 | 100.0 | 101.3 | |
| 48 Chlorobromomethane | 128 | 5.986 | 5.986 | 0.000 | 98 | 118276 | 50.0 | 46.0 | |
| 49 Tetrahydrofuran | 42 | 5.998 | 5.998 | 0.000 | 90 | 162324 | 100.0 | 91.2 | |
| 50 Chloroform | 83 | 6.138 | 6.138 | 0.000 | 93 | 446745 | 50.0 | 50.5 | |
| 51 1,1,1-Trichloroethane | 97 | 6.296 | 6.296 | 0.000 | 97 | 274535 | 50.0 | 48.8 | |
| 52 Cyclohexane | 56 | 6.363 | 6.363 | 0.000 | 90 | 405642 | 50.0 | 50.8 | |
| 53 Carbon tetrachloride | 117 | 6.466 | 6.466 | 0.000 | 95 | 190304 | 50.0 | 47.1 | |
| 54 1,1-Dichloropropene | 75 | 6.485 | 6.485 | 0.000 | 98 | 340889 | 50.0 | 50.3 | |
| 56 Benzene | 78 | 6.698 | 6.698 | 0.000 | 97 | 992835 | 50.0 | 50.4 | |
| 55 Isobutyl alcohol | 41 | 6.704 | 6.704 | 0.000 | 41 | 132080 | 1250.0 | 1113.7 | |
| 57 1,2-Dichloroethane | 62 | 6.777 | 6.777 | 0.000 | 98 | 349409 | 50.0 | 48.0 | |
| 59 n-Heptane | 43 | 7.075 | 7.075 | 0.000 | 86 | 220812 | 50.0 | 48.6 | |
| 61 Trichloroethene | 130 | 7.446 | 7.446 | 0.000 | 97 | 233389 | 50.0 | 48.7 | |
| 63 Methylcyclohexane | 83 | 7.677 | 7.677 | 0.000 | 86 | 419227 | 50.0 | 50.1 | |
| 64 1,2-Dichloropropane | 63 | 7.720 | 7.720 | 0.000 | 95 | 234878 | 50.0 | 47.7 | |
| 65 1,4-Dioxane | 88 | 7.805 | 7.805 | 0.000 | 31 | 45919 | 1000.0 | 1022.1 | M |
| 67 Dibromomethane | 93 | 7.811 | 7.811 | 0.000 | 95 | 150124 | 50.0 | 46.6 | |
| 68 Dichlorobromomethane | 83 | 8.005 | 8.005 | 0.000 | 99 | 246980 | 50.0 | 45.0 | |
| 70 2-Chloroethyl vinyl ether | 63 | 8.316 | 8.316 | 0.000 | 91 | 304468 | 100.0 | 93.3 | |
| 71 cis-1,3-Dichloropropene | 75 | 8.456 | 8.456 | 0.000 | 95 | 279255 | 50.0 | 44.9 | |
| 72 4-Methyl-2-pentanone (MIBK) | 43 | 8.614 | 8.614 | 0.000 | 94 | 495598 | 100.0 | 103.5 | |
| 73 Toluene | 91 | 8.784 | 8.784 | 0.000 | 99 | 966776 | 50.0 | 52.3 | |
| 74 trans-1,3-Dichloropropene | 75 | 9.040 | 9.040 | 0.000 | 92 | 223869 | 50.0 | 45.3 | |
| 75 Ethyl methacrylate | 69 | 9.107 | 9.107 | 0.000 | 88 | 305467 | 50.0 | 49.0 | |
| 76 1,1,2-Trichloroethane | 97 | 9.228 | 9.228 | 0.000 | 91 | 211345 | 50.0 | 49.5 | |
| 77 Tetrachloroethene | 164 | 9.295 | 9.295 | 0.000 | 95 | 165794 | 50.0 | 50.1 | |
| 78 1,3-Dichloropropane | 76 | 9.386 | 9.386 | 0.000 | 89 | 380984 | 50.0 | 49.2 | |
| 79 2-Hexanone | 43 | 9.453 | 9.453 | 0.000 | 94 | 306856 | 100.0 | 99.6 | |
| 81 Chlorodibromomethane | 129 | 9.599 | 9.599 | 0.000 | 89 | 131179 | 50.0 | 45.3 | |
| 82 Ethylene Dibromide | 107 | 9.709 | 9.709 | 0.000 | 99 | 200005 | 50.0 | 48.6 | |
| 83 3-Chlorobenzotrifluoride | 180 | 10.183 | 10.183 | 0.000 | 93 | 270429 | 50.0 | 50.4 | |
| 84 Chlorobenzene | 112 | 10.202 | 10.202 | 0.000 | 94 | 613324 | 50.0 | 51.1 | |
| 85 4-Chlorobenzotrifluoride | 180 | 10.269 | 10.269 | 0.000 | 97 | 246049 | 50.0 | 49.4 | |
| 86 1,1,1,2-Tetrachloroethane | 131 | 10.293 | 10.293 | 0.000 | 87 | 156373 | 50.0 | 45.8 | |
| 87 Ethylbenzene | 106 | 10.305 | 10.305 | 0.000 | 98 | 352019 | 50.0 | 51.2 | |
| 88 m-Xylene & p-Xylene | 106 | 10.433 | 10.433 | 0.000 | 100 | 416283 | 50.0 | 49.4 | |
| 89 o-Xylene | 106 | 10.816 | 10.816 | 0.000 | 96 | 430994 | 50.0 | 51.1 | |
| 90 Styrene | 104 | 10.834 | 10.834 | 0.000 | 94 | 702614 | 50.0 | 51.5 | |
| 91 Bromoform | 173 | 11.017 | 11.017 | 0.000 | 94 | 71560 | 50.0 | 43.9 | |
| 92 2-Chlorobenzotrifluoride | 180 | 11.090 | 11.090 | 0.000 | 95 | 265784 | 50.0 | 48.8 | |
| 93 Isopropylbenzene | 105 | 11.181 | 11.181 | 0.000 | 96 | 995129 | 50.0 | 52.7 | |
| 95 Bromobenzene | 156 | 11.491 | 11.491 | 0.000 | 95 | 249243 | 50.0 | 50.5 | |
| 96 1,1,2,2-Tetrachloroethane | 83 | 11.497 | 11.497 | 0.000 | 95 | 301111 | 50.0 | 50.1 | |
| 97 trans-1,4-Dichloro-2-butene | 53 | 11.534 | 11.534 | 0.000 | 75 | 66355 | 50.0 | 45.9 | |
| 98 1,2,3-Trichloropropane | 110 | 11.552 | 11.552 | 0.000 | 89 | 103048 | 50.0 | 50.3 | |
| 99 N-Propylbenzene | 120 | 11.601 | 11.601 | 0.000 | 98 | 267230 | 50.0 | 49.7 | |
| 100 2-Chlorotoluene | 126 | 11.680 | 11.680 | 0.000 | 95 | 230624 | 50.0 | 49.5 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 101 3-Chlorotoluene | 126 | 11.747 | 11.747 | 0.000 | 96 | 239610 | 50.0 | 49.1 | |
| 102 1,3,5-Trimethylbenzene | 105 | 11.783 | 11.783 | 0.000 | 93 | 790599 | 50.0 | 51.8 | |
| 103 4-Chlorotoluene | 126 | 11.808 | 11.808 | 0.000 | 99 | 253826 | 50.0 | 50.0 | |
| 104 tert-Butylbenzene | 119 | 12.094 | 12.094 | 0.000 | 91 | 630674 | 50.0 | 52.1 | |
| 106 1,2,4-Trimethylbenzene | 105 | 12.154 | 12.154 | 0.000 | 98 | 824782 | 50.0 | 51.8 | |
| 107 1,2-dichloro-4-(trifluoromethyl) | 214 | 12.209 | 12.209 | 0.000 | 97 | 164643 | 50.0 | 46.6 | |
| 108 sec-Butylbenzene | 105 | 12.319 | 12.319 | 0.000 | 95 | 875734 | 50.0 | 51.5 | |
| 109 1,3-Dichlorobenzene | 146 | 12.434 | 12.434 | 0.000 | 96 | 438172 | 50.0 | 49.4 | |
| 110 4-Isopropyltoluene | 119 | 12.477 | 12.477 | 0.000 | 96 | 717050 | 50.0 | 51.1 | |
| 111 1,4-Dichlorobenzene | 146 | 12.538 | 12.538 | 0.000 | 94 | 464508 | 50.0 | 50.7 | |
| 113 2,4-Dichloro-1-(trifluoromethyl) | 214 | 12.568 | 12.568 | 0.000 | 95 | 161786 | 50.0 | 48.6 | |
| 114 2,5-Dichlorobenzotrifluoride | 214 | 12.611 | 12.611 | 0.000 | 97 | 173267 | 50.0 | 47.2 | |
| 116 n-Butylbenzene | 91 | 12.884 | 12.884 | 0.000 | 97 | 643029 | 50.0 | 50.2 | |
| 117 1,2-Dichlorobenzene | 146 | 12.890 | 12.890 | 0.000 | 96 | 416117 | 50.0 | 49.3 | |
| 118 1,2-Dibromo-3-Chloropropan | 75 | 13.681 | 13.681 | 0.000 | 77 | 33625 | 50.0 | 44.7 | |
| 119 2,4- & 2,5- & 2,6- Dichlorobenzene | 125 | 13.821 | 13.821 | 0.000 | 99 | 780454 | 150.0 | 145.3 | |
| 121 2,3- & 3,4- Dichlorotoluene | 125 | 14.235 | 14.235 | 0.000 | 99 | 550121 | 100.0 | 93.8 | |
| 122 1,2,4-Trichlorobenzene | 180 | 14.497 | 14.497 | 0.000 | 94 | 199297 | 50.0 | 45.2 | |
| 123 Hexachlorobutadiene | 225 | 14.649 | 14.649 | 0.000 | 97 | 58792 | 50.0 | 42.1 | |
| 124 Naphthalene | 128 | 14.764 | 14.764 | 0.000 | 98 | 627907 | 50.0 | 46.7 | |
| 125 1,2,3-Trichlorobenzene | 180 | 14.983 | 14.983 | 0.000 | 95 | 169607 | 50.0 | 42.9 | |
| 126 2,4,5-Trichlorotoluene | 159 | 15.774 | 15.774 | 0.000 | 0 | 90501 | 50.0 | 39.8 | |
| 127 2,3,6-Trichlorotoluene | 159 | 15.884 | 15.884 | 0.000 | 96 | 84129 | 50.0 | 41.2 | |
| 146 3,4-Dichlorotoluene | 1 | 0.000 | | | | | ND | ND | |
| S 130 1,2-Dichloroethene, Total | 96 | | | | 0 | | 100.0 | 97.8 | |
| S 131 Xylenes, Total | 106 | | | | 0 | | 100.0 | 100.5 | |
| S 132 1,3-Dichloropropene, Total | 1 | | | | 0 | | 100.0 | 90.1 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

Reagents:

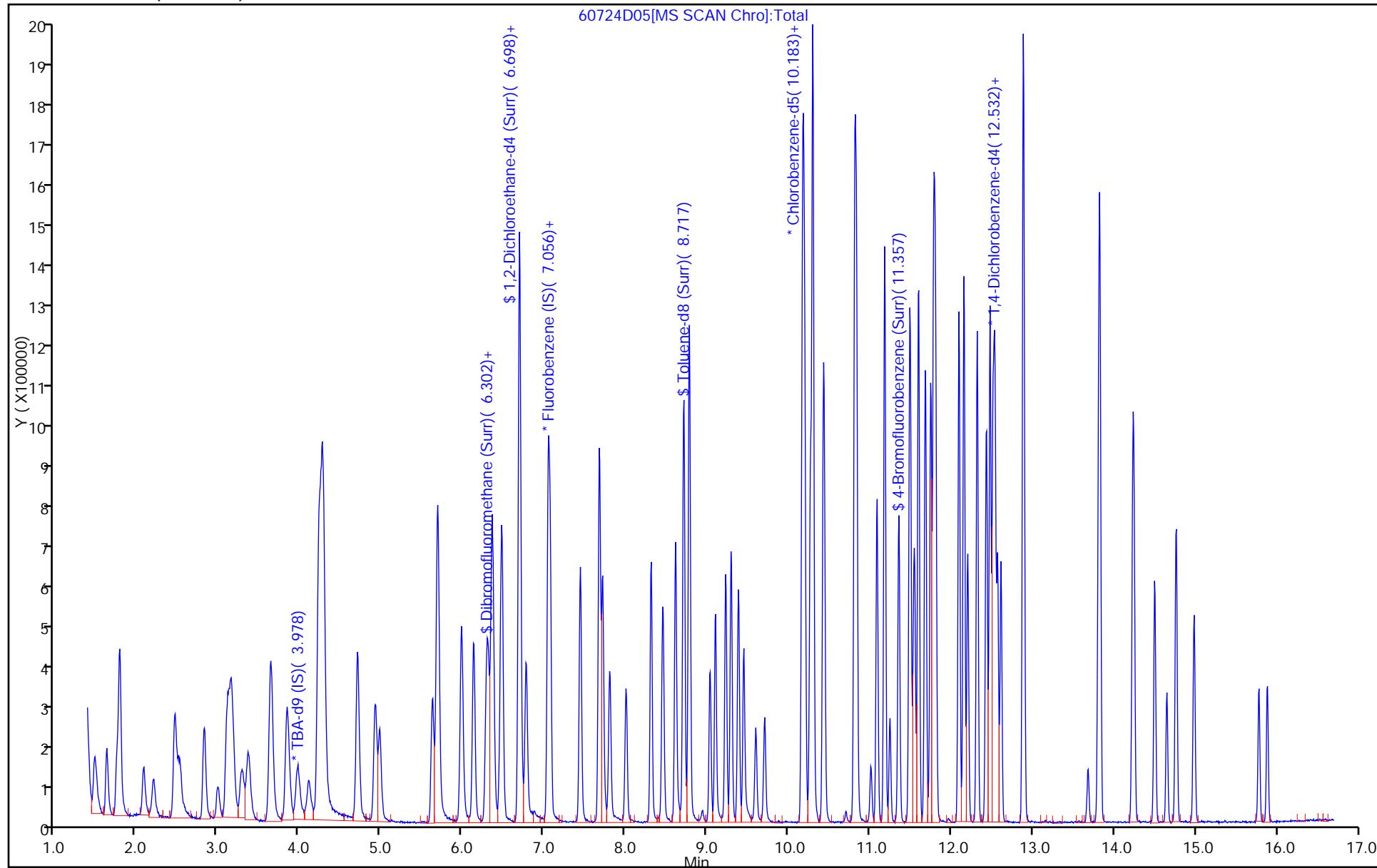
| | | |
|---------------------|--------------------|-----------|
| voaWEEmix1stR_00009 | Amount Added: 2.00 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 2.00 | Units: uL |
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 2.00 | Units: uL |
| voaWVA1stRest_00016 | Amount Added: 2.00 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 6.00 | Units: uL |
| VOA8260VOAPRI_00263 | Amount Added: 2.00 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 2.00 | Units: uL |

Report Date: 25-Jul-2017 02:13:18

Chrom Revision: 2.2 20-Jun-2017 07:42:38

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D05.D
 Injection Date: 24-Jul-2017 07:27:30 Instrument ID: CHHP6
 Lims ID: ICIS VSTD10 Operator ID: 034635
 Client ID:
 Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 5
 Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
 Column: DB-624 (0.18 mm)



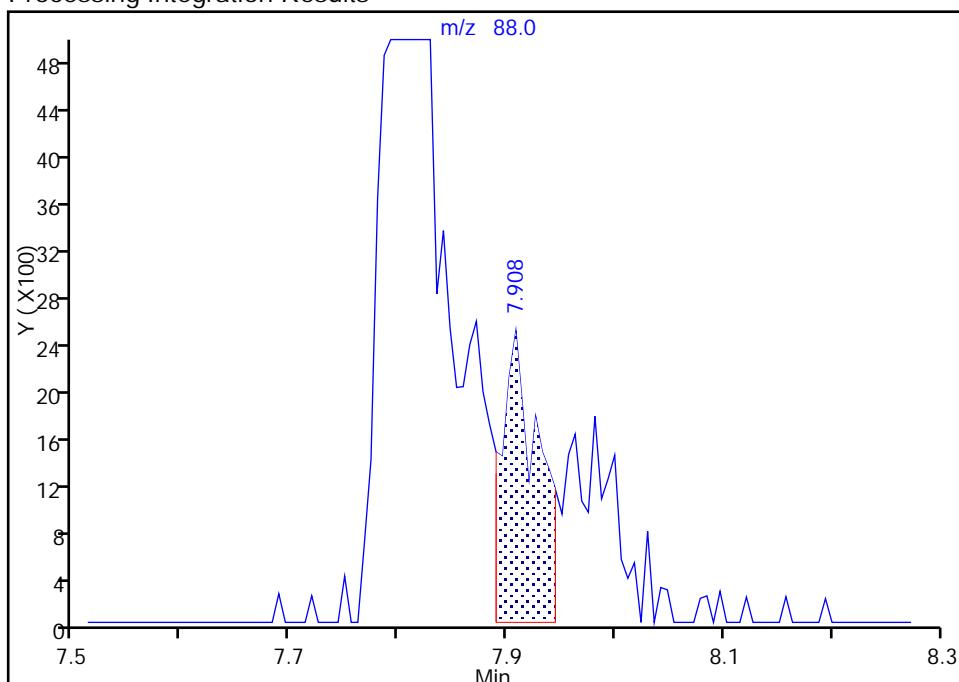
TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D05.D
 Injection Date: 24-Jul-2017 07:27:30 Instrument ID: CHHP6
 Lims ID: ICIS VSTD10
 Client ID:
 Operator ID: 034635 ALS Bottle#: 5 Worklist Smp#: 5
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
 Column: DB-624 (0.18 mm) Detector: MS SCAN

65 1,4-Dioxane, CAS: 123-91-1
Signal: 1

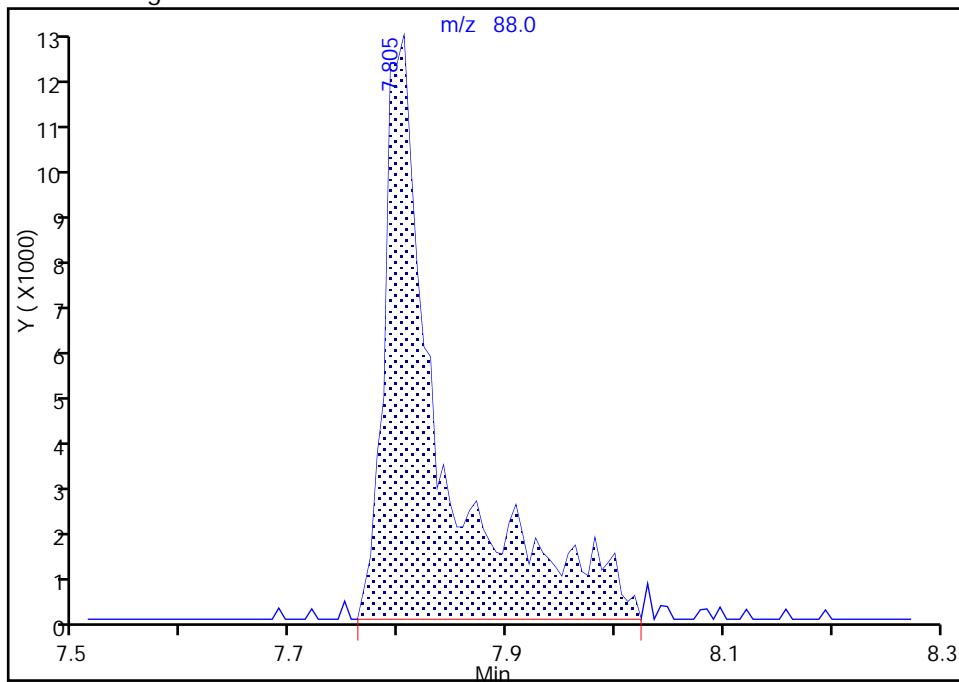
RT: 7.91
 Area: 5831
 Amount: 154.9459
 Amount Units: ng

Processing Integration Results



RT: 7.80
 Area: 45919
 Amount: 1022.1306
 Amount Units: ng

Manual Integration Results



Reviewer: bungardf, 24-Jul-2017 07:58:04

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D06.D
 Lims ID: IC VSTD15
 Client ID:
 Sample Type: IC Calib Level: 4
 Inject. Date: 24-Jul-2017 07:52:30 ALS Bottle#: 6 Worklist Smp#: 6
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017705-006
 Misc. Info.: IC VSTD15
 Operator ID: 034635 Instrument ID: CHHP6
 Sublist: chrom-MSVOA_LL_CHHP6*sub10
 Method: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 25-Jul-2017 01:44:34 Calib Date: 24-Jul-2017 09:28:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D10.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK021

First Level Reviewer: bungardf Date: 24-Jul-2017 08:44:01

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 3.972 | 3.972 | 0.000 | 94 | 306123 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.051 | 7.051 | 0.000 | 98 | 837369 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.172 | 10.172 | 0.000 | 87 | 185833 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.514 | 12.514 | 0.000 | 93 | 281796 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr) | 113 | 6.315 | 6.315 | 0.000 | 93 | 321543 | 75.0 | 73.9 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur) | 65 | 6.692 | 6.692 | 0.000 | 69 | 450831 | 75.0 | 72.5 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.718 | 8.718 | 0.000 | 93 | 1165400 | 75.0 | 79.9 | |
| \$ 8 4-Bromofluorobenzene (Surr) | 95 | 11.352 | 11.352 | 0.000 | 85 | 477594 | 75.0 | 74.4 | |
| 11 Dichlorodifluoromethane | 85 | 1.484 | 1.484 | 0.000 | 98 | 392309 | 75.0 | 73.4 | |
| 12 Chloromethane | 50 | 1.630 | 1.630 | 0.000 | 98 | 364709 | 75.0 | 75.9 | |
| 13 Vinyl chloride | 62 | 1.758 | 1.758 | 0.000 | 98 | 390082 | 75.0 | 77.4 | |
| 14 Butadiene | 39 | 1.788 | 1.788 | 0.000 | 89 | 315046 | 75.0 | 75.4 | |
| 15 Bromomethane | 94 | 2.087 | 2.087 | 0.000 | 90 | 178416 | 75.0 | 76.0 | |
| 16 Chloroethane | 64 | 2.202 | 2.202 | 0.000 | 98 | 212582 | 75.0 | 78.1 | |
| 17 Dichlorofluoromethane | 67 | 2.470 | 2.470 | 0.000 | 96 | 452413 | 75.0 | 77.1 | |
| 18 Trichlorofluoromethane | 101 | 2.506 | 2.506 | 0.000 | 98 | 381467 | 75.0 | 76.9 | |
| 20 Ethyl ether | 59 | 2.823 | 2.823 | 0.000 | 87 | 314417 | 75.0 | 74.1 | |
| 21 Acrolein | 56 | 2.999 | 2.999 | 0.000 | 100 | 161845 | 175.0 | 178.0 | |
| 22 1,1-Dichloroethene | 96 | 3.115 | 3.115 | 0.000 | 98 | 326499 | 75.0 | 75.0 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.157 | 3.157 | 0.000 | 95 | 320427 | 75.0 | 76.8 | |
| 24 Acetone | 43 | 3.194 | 3.194 | 0.000 | 100 | 295809 | 150.0 | 164.7 | |
| 25 Iodomethane | 142 | 3.291 | 3.291 | 0.000 | 98 | 465530 | 75.0 | 76.1 | |
| 26 Carbon disulfide | 76 | 3.364 | 3.364 | 0.000 | 99 | 721571 | 75.0 | 74.5 | |
| 29 3-Chloro-1-propene | 76 | 3.638 | 3.638 | 0.000 | 92 | 194002 | 75.0 | 76.1 | |
| 30 Methyl acetate | 43 | 3.662 | 3.662 | 0.000 | 97 | 572896 | 150.0 | 145.5 | |
| 31 Methylene Chloride | 84 | 3.839 | 3.839 | 0.000 | 89 | 418660 | 75.0 | 72.6 | |
| 32 2-Methyl-2-propanol | 59 | 4.112 | 4.112 | 0.000 | 93 | 251071 | 750.0 | 741.0 | |
| 33 Acrylonitrile | 53 | 4.240 | 4.240 | 0.000 | 99 | 1582844 | 750.0 | 752.5 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.277 | 4.277 | 0.000 | 98 | 372663 | 75.0 | 75.5 | |
| 35 Methyl tert-butyl ether | 73 | 4.295 | 4.295 | 0.000 | 95 | 1181045 | 75.0 | 74.8 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|--------------|------------------|------------------|-----|----------|---------------|-----------------|-------|
| 36 Hexane | 57 | 4.709 | 4.709 | 0.000 | 91 | 437873 | 75.0 | 74.9 | |
| 37 1,1-Dichloroethane | 63 | 4.928 | 4.928 | 0.000 | 95 | 641771 | 75.0 | 79.9 | |
| 38 Vinyl acetate | 43 | 4.982 | 4.982 | 0.000 | 97 | 703120 | 75.0 | 75.2 | |
| 42 2,2-Dichloropropane | 97 | 5.682 | 5.682 | 0.000 | 56 | 64241 | 75.0 | 77.2 | |
| 43 cis-1,2-Dichloroethene | 96 | 5.694 | 5.694 | 0.000 | 80 | 436220 | 75.0 | 75.6 | |
| 44 2-Butanone (MEK) | 43 | 5.706 | 5.706 | 0.000 | 97 | 385993 | 150.0 | 149.9 | |
| 48 Chlorobromomethane | 128 | 5.986 | 5.986 | 0.000 | 98 | 188762 | 75.0 | 75.0 | |
| 49 Tetrahydrofuran | 42 | 5.992 | 5.992 | 0.000 | 85 | 242295 | 150.0 | 139.0 | |
| 50 Chloroform | 83 | 6.138 | 6.138 | 0.000 | 93 | 660420 | 75.0 | 76.2 | |
| 51 1,1,1-Trichloroethane | 97 | 6.290 | 6.290 | 0.000 | 98 | 414781 | 75.0 | 75.3 | |
| 52 Cyclohexane | 56 | 6.363 | 6.363 | 0.000 | 90 | 597815 | 75.0 | 76.4 | |
| 53 Carbon tetrachloride | 117 | 6.467 | 6.467 | 0.000 | 96 | 300016 | 75.0 | 75.8 | |
| 54 1,1-Dichloropropene | 75 | 6.485 | 6.485 | 0.000 | 98 | 506906 | 75.0 | 76.4 | |
| 56 Benzene | 78 | 6.704 | 6.704 | 0.000 | 97 | 1504594 | 75.0 | 78.0 | |
| 57 1,2-Dichloroethane | 62 | 6.783 | 6.783 | 0.000 | 97 | 540487 | 75.0 | 75.8 | |
| 59 n-Heptane | 43 | 7.075 | 7.075 | 0.000 | 89 | 337209 | 75.0 | 75.8 | |
| 55 Isobutyl alcohol | 41 | 6.698 | 6.698 | 0.000 | 42 | 208868 | 1875.0 | 1798.2 | |
| 61 Trichloroethene | 130 | 7.446 | 7.446 | 0.000 | 98 | 355153 | 75.0 | 75.7 | |
| 63 Methylcyclohexane | 83 | 7.677 | 7.677 | 0.000 | 87 | 630450 | 75.0 | 77.0 | |
| 64 1,2-Dichloropropane | 63 | 7.720 | 7.720 | 0.000 | 93 | 361293 | 75.0 | 74.9 | |
| 67 Dibromomethane | 93 | 7.811 | 7.811 | 0.000 | 96 | 233620 | 75.0 | 74.0 | |
| 65 1,4-Dioxane | 88 | 7.805 | 7.805 | 0.000 | 61 | 65610 | 1500.0 | 1491.2 | M |
| 68 Dichlorobromomethane | 83 | 8.006 | 8.006 | 0.000 | 100 | 411363 | 75.0 | 76.6 | |
| 70 2-Chloroethyl vinyl ether | 63 | 8.316 | 8.316 | 0.000 | 92 | 478201 | 150.0 | 149.6 | |
| 71 cis-1,3-Dichloropropene | 75 | 8.456 | 8.456 | 0.000 | 96 | 460333 | 75.0 | 75.5 | |
| 72 4-Methyl-2-pentanone (MIBK) | 43 | 8.614 | 8.614 | 0.000 | 94 | 755882 | 150.0 | 159.3 | |
| 73 Toluene | 91 | 8.784 | 8.784 | 0.000 | 98 | 1461492 | 75.0 | 79.8 | |
| 74 trans-1,3-Dichloropropene | 75 | 9.040 | 9.040 | 0.000 | 92 | 381658 | 75.0 | 77.8 | |
| 75 Ethyl methacrylate | 69 | 9.101 | 9.101 | 0.000 | 87 | 492154 | 75.0 | 79.6 | |
| 76 1,1,2-Trichloroethane | 97 | 9.229 | 9.229 | 0.000 | 91 | 325307 | 75.0 | 76.8 | |
| 77 Tetrachloroethene | 164 | 9.295 | 9.295 | 0.000 | 97 | 257015 | 75.0 | 78.4 | |
| 78 1,3-Dichloropropane | 76 | 9.387 | 9.387 | 0.000 | 89 | 593488 | 75.0 | 77.2 | |
| 79 2-Hexanone | 43 | 9.454 | 9.454 | 0.000 | 94 | 480105 | 150.0 | 157.2 | |
| 81 Chlorodibromomethane | 129 | 9.600 | 9.600 | 0.000 | 89 | 216633 | 75.0 | 75.5 | |
| 82 Ethylene Dibromide | 107 | 9.709 | 9.709 | 0.000 | 98 | 313940 | 75.0 | 77.0 | |
| 83 3-Chlorobenzotrifluoride | 180 | 10.184 | 10.184 | 0.000 | 92 | 444234 | 75.0 | 83.5 | |
| 84 Chlorobenzene | 112 | 10.202 | 10.202 | 0.000 | 95 | 923461 | 75.0 | 77.5 | |
| 85 4-Chlorobenzotrifluoride | 180 | 10.269 | 10.269 | 0.000 | 97 | 418159 | 75.0 | 84.8 | |
| 86 1,1,1,2-Tetrachloroethane | 131 | 10.293 | 10.293 | 0.000 | 88 | 270626 | 75.0 | 79.9 | |
| 87 Ethylbenzene | 106 | 10.299 | 10.299 | 0.000 | 98 | 540015 | 75.0 | 79.3 | |
| 88 m-Xylene & p-Xylene | 106 | 10.433 | 10.433 | 0.000 | 99 | 663604 | 75.0 | 79.5 | |
| 89 o-Xylene | 106 | 10.816 | 10.816 | 0.000 | 95 | 657817 | 75.0 | 78.6 | |
| 90 Styrene | 104 | 10.835 | 10.835 | 0.000 | 94 | 1093228 | 75.0 | 80.8 | |
| 91 Bromoform | 173 | 11.017 | 11.017 | 0.000 | 94 | 120009 | 75.0 | 74.3 | |
| 92 2-Chlorobenzotrifluoride | 180 | 11.090 | 11.090 | 0.000 | 96 | 457223 | 75.0 | 84.7 | |
| 93 Isopropylbenzene | 105 | 11.181 | 11.181 | 0.000 | 97 | 1511361 | 75.0 | 80.7 | |
| 95 Bromobenzene | 156 | 11.492 | 11.492 | 0.000 | 95 | 380643 | 75.0 | 72.7 | |
| 96 1,1,2,2-Tetrachloroethane | 83 | 11.498 | 11.498 | 0.000 | 95 | 468952 | 75.0 | 78.7 | |
| 97 trans-1,4-Dichloro-2-butene | 53 | 11.534 | 11.534 | 0.000 | 76 | 105972 | 75.0 | 69.1 | |
| 98 1,2,3-Trichloropropane | 110 | 11.552 | 11.552 | 0.000 | 87 | 159086 | 75.0 | 73.1 | |
| 99 N-Propylbenzene | 120 | 11.601 | 11.601 | 0.000 | 98 | 428911 | 75.0 | 75.2 | |
| 100 2-Chlorotoluene | 126 | 11.680 | 11.680 | 0.000 | 95 | 370145 | 75.0 | 75.0 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 101 3-Chlorotoluene | 126 | 11.747 | 11.747 | 0.000 | 95 | 403537 | 75.0 | 77.9 | |
| 102 1,3,5-Trimethylbenzene | 105 | 11.784 | 11.784 | 0.000 | 94 | 1254643 | 75.0 | 77.5 | |
| 103 4-Chlorotoluene | 126 | 11.808 | 11.808 | 0.000 | 99 | 398838 | 75.0 | 74.0 | |
| 104 tert-Butylbenzene | 119 | 12.094 | 12.094 | 0.000 | 91 | 985989 | 75.0 | 76.7 | |
| 106 1,2,4-Trimethylbenzene | 105 | 12.155 | 12.155 | 0.000 | 98 | 1282863 | 75.0 | 76.0 | |
| 107 1,2-dichloro-4-(trifluoromethyl) | 214 | 12.203 | 12.203 | 0.000 | 96 | 293347 | 75.0 | 78.3 | |
| 108 sec-Butylbenzene | 105 | 12.319 | 12.319 | 0.000 | 95 | 1379795 | 75.0 | 76.5 | |
| 109 1,3-Dichlorobenzene | 146 | 12.435 | 12.435 | 0.000 | 95 | 697551 | 75.0 | 74.2 | |
| 110 4-Isopropyltoluene | 119 | 12.477 | 12.477 | 0.000 | 96 | 1135538 | 75.0 | 76.3 | |
| 111 1,4-Dichlorobenzene | 146 | 12.538 | 12.538 | 0.000 | 93 | 720028 | 75.0 | 74.0 | |
| 113 2,4-Dichloro-1-(trifluoromethyl) | 214 | 12.568 | 12.568 | 0.000 | 97 | 283275 | 75.0 | 80.3 | |
| 114 2,5-Dichlorobenzotrifluoride | 214 | 12.611 | 12.611 | 0.000 | 97 | 298436 | 75.0 | 76.6 | |
| 116 n-Butylbenzene | 91 | 12.885 | 12.885 | 0.000 | 97 | 1048415 | 75.0 | 77.2 | |
| 117 1,2-Dichlorobenzene | 146 | 12.891 | 12.891 | 0.000 | 96 | 662537 | 75.0 | 74.1 | |
| 118 1,2-Dibromo-3-Chloropropan | 75 | 13.682 | 13.682 | 0.000 | 79 | 56051 | 75.0 | 70.3 | |
| 119 2,4- & 2,5- & 2,6- Dichlorobenzene | 125 | 13.822 | 13.822 | 0.000 | 99 | 1316529 | 225.0 | 231.1 | |
| 121 2,3- & 3,4- Dichlorotoluene | 125 | 14.235 | 14.235 | 0.000 | 98 | 921595 | 150.0 | 148.1 | |
| 122 1,2,4-Trichlorobenzene | 180 | 14.503 | 14.503 | 0.000 | 93 | 325076 | 75.0 | 69.5 | |
| 123 Hexachlorobutadiene | 225 | 14.649 | 14.649 | 0.000 | 96 | 101525 | 75.0 | 68.6 | |
| 124 Naphthalene | 128 | 14.765 | 14.765 | 0.000 | 97 | 994327 | 75.0 | 69.8 | |
| 125 1,2,3-Trichlorobenzene | 180 | 14.984 | 14.984 | 0.000 | 95 | 275471 | 75.0 | 65.6 | |
| 126 2,4,5-Trichlorotoluene | 159 | 15.780 | 15.780 | 0.000 | 0 | 146009 | 75.0 | 60.5 | |
| 127 2,3,6-Trichlorotoluene | 159 | 15.884 | 15.884 | 0.000 | 96 | 129996 | 75.0 | 60.1 | |
| 146 3,4-Dichlorotoluene | 1 | 0.000 | | | | | ND | ND | |
| S 131 Xylenes, Total | 106 | | | | 0 | | 150.0 | 158.1 | |
| S 130 1,2-Dichloroethene, Total | 96 | | | | 0 | | 150.0 | 151.1 | |
| S 132 1,3-Dichloropropene, Total | 1 | | | | 0 | | 150.0 | 153.4 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

Reagents:

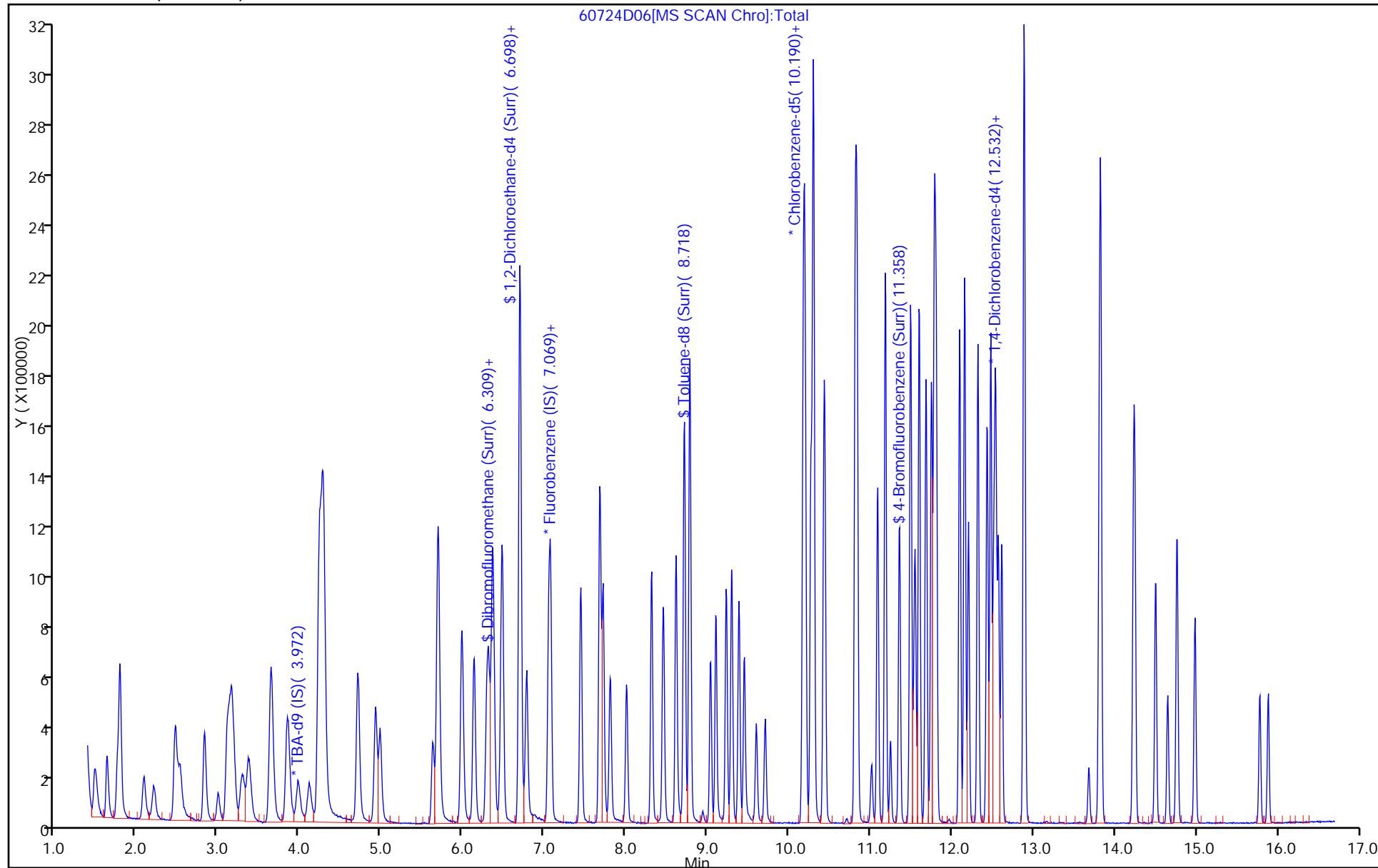
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|---------------------|--------------------|-----------|
| voaWEEmix1stR_00009 | Amount Added: 3.00 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 3.00 | Units: uL |
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 3.00 | Units: uL |
| voaWVA1stRest_00016 | Amount Added: 3.00 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 7.00 | Units: uL |
| VOA8260VOAPRI_00263 | Amount Added: 3.00 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 3.00 | Units: uL |

Report Date: 25-Jul-2017 01:44:35

Chrom Revision: 2.2 20-Jun-2017 07:42:38

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D06.D
Injection Date: 24-Jul-2017 07:52:30 Instrument ID: CHHP6
Lims ID: IC VSTD15 Operator ID: 034635
Client ID:
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 6
Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh

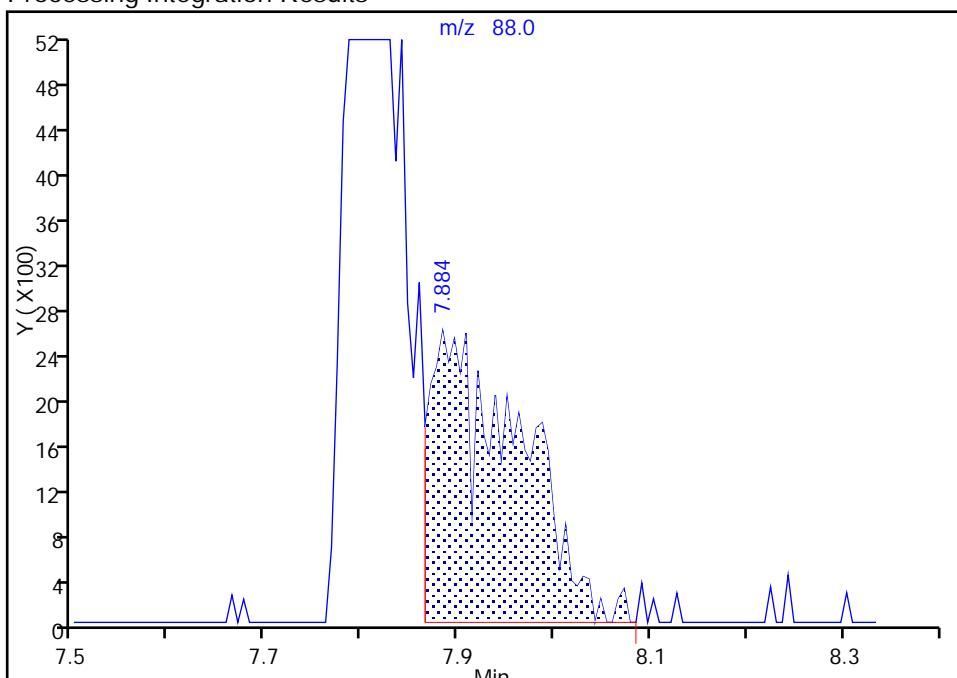
Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D06.D
 Injection Date: 24-Jul-2017 07:52:30 Instrument ID: CHHP6
 Lims ID: IC VSTD15
 Client ID:
 Operator ID: 034635 ALS Bottle#: 6 Worklist Smp#: 6
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
 Column: DB-624 (0.18 mm) Detector: MS SCAN

65 1,4-Dioxane, CAS: 123-91-1

Signal: 1

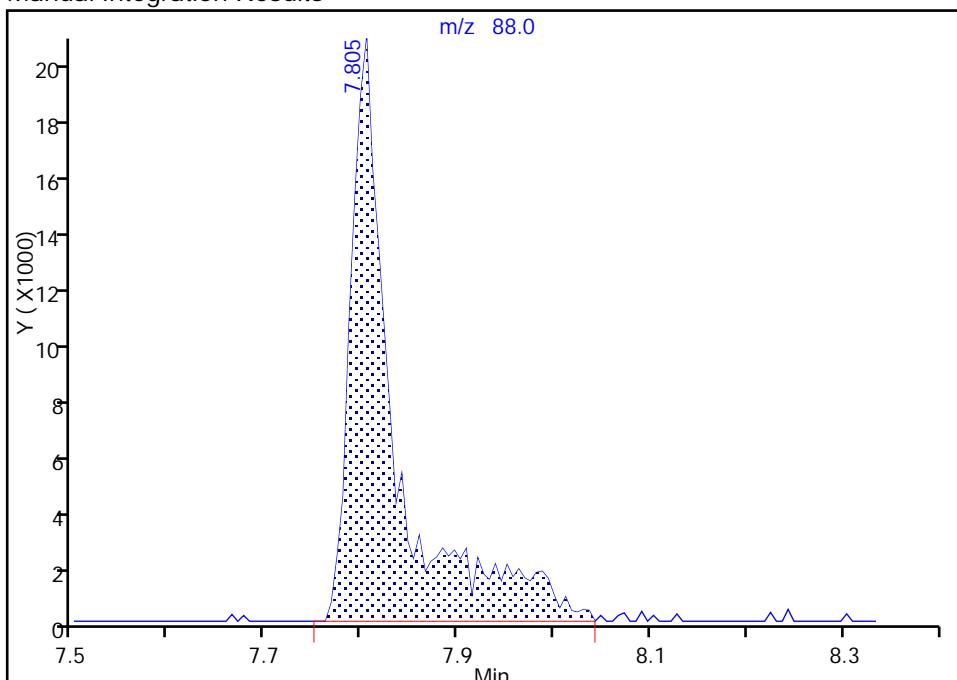
RT: 7.88
 Area: 16500
 Amount: 410.3640
 Amount Units: ng

Processing Integration Results



RT: 7.81
 Area: 65610
 Amount: 1491.1699
 Amount Units: ng

Manual Integration Results



Reviewer: bungardf, 24-Jul-2017 08:11:32

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170724-17705.b\60724D07.D
 Lims ID: IC VSTD20
 Client ID:
 Sample Type: IC Calib Level: 5
 Inject. Date: 24-Jul-2017 08:16:30 ALS Bottle#: 7 Worklist Smp#: 7
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017705-007
 Misc. Info.: IC VSTD20
 Operator ID: 034635 Instrument ID: CHHP6
 Sublist: chrom-MSVOA_LL_CHHP6*sub10
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170724-17705.b\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 25-Jul-2017 01:44:36 Calib Date: 24-Jul-2017 09:28:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170724-17705.b\60724D10.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK021

First Level Reviewer: bungardf Date: 24-Jul-2017 08:52:45

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 3.976 | 3.972 | 0.004 | 94 | 272680 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.048 | 7.051 | -0.003 | 99 | 867274 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.169 | 10.172 | -0.003 | 87 | 191685 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.511 | 12.514 | -0.003 | 96 | 258316 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr) | 113 | 6.318 | 6.315 | 0.003 | 93 | 424756 | 100.0 | 94.2 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur) | 65 | 6.689 | 6.692 | -0.003 | 70 | 572691 | 100.0 | 88.9 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.715 | 8.718 | -0.003 | 93 | 1484720 | 100.0 | 99.4 | |
| \$ 8 4-Bromofluorobenzene (Surr) | 95 | 11.355 | 11.352 | 0.003 | 83 | 609762 | 100.0 | 92.1 | |
| 11 Dichlorodifluoromethane | 85 | 1.482 | 1.484 | -0.002 | 98 | 537841 | 100.0 | 97.2 | |
| 12 Chloromethane | 50 | 1.628 | 1.630 | -0.002 | 99 | 485997 | 100.0 | 97.7 | |
| 13 Vinyl chloride | 62 | 1.755 | 1.758 | -0.003 | 98 | 507676 | 100.0 | 97.3 | |
| 14 Butadiene | 39 | 1.786 | 1.788 | -0.002 | 89 | 407662 | 100.0 | 94.3 | |
| 15 Bromomethane | 94 | 2.084 | 2.087 | -0.003 | 90 | 239488 | 100.0 | 98.5 | |
| 16 Chloroethane | 64 | 2.199 | 2.202 | -0.003 | 98 | 283541 | 100.0 | 100.6 | |
| 17 Dichlorofluoromethane | 67 | 2.467 | 2.470 | -0.003 | 97 | 601770 | 100.0 | 99.1 | |
| 18 Trichlorofluoromethane | 101 | 2.498 | 2.506 | -0.008 | 99 | 515987 | 100.0 | 100.4 | |
| 20 Ethyl ether | 59 | 2.820 | 2.823 | -0.003 | 88 | 412759 | 100.0 | 93.9 | |
| 21 Acrolein | 56 | 2.990 | 2.999 | -0.009 | 100 | 172660 | 200.0 | 183.4 | |
| 22 1,1-Dichloroethene | 96 | 3.112 | 3.115 | -0.003 | 97 | 437661 | 100.0 | 97.1 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.155 | 3.157 | -0.002 | 95 | 418931 | 100.0 | 97.0 | |
| 24 Acetone | 43 | 3.191 | 3.194 | -0.003 | 100 | 345529 | 200.0 | 185.7 | |
| 25 Iodomethane | 142 | 3.282 | 3.291 | -0.009 | 99 | 616342 | 100.0 | 97.3 | |
| 26 Carbon disulfide | 76 | 3.361 | 3.364 | -0.003 | 99 | 1031794 | 100.0 | 102.8 | |
| 29 3-Chloro-1-propene | 76 | 3.635 | 3.638 | -0.003 | 93 | 261163 | 100.0 | 98.9 | |
| 30 Methyl acetate | 43 | 3.653 | 3.662 | -0.009 | 97 | 775230 | 200.0 | 190.1 | |
| 31 Methylene Chloride | 84 | 3.842 | 3.839 | 0.003 | 88 | 557470 | 100.0 | 93.3 | |
| 32 2-Methyl-2-propanol | 59 | 4.104 | 4.112 | -0.008 | 94 | 297303 | 1000.0 | 985.0 | |
| 33 Acrylonitrile | 53 | 4.231 | 4.240 | -0.009 | 98 | 2040960 | 1000.0 | 936.8 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.274 | 4.277 | -0.003 | 99 | 496677 | 100.0 | 97.1 | |
| 35 Methyl tert-butyl ether | 73 | 4.286 | 4.295 | -0.009 | 96 | 1550808 | 100.0 | 94.8 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|--------------|------------------|------------------|-----|----------|---------------|-----------------|-------|
| 36 Hexane | 57 | 4.706 | 4.709 | -0.003 | 90 | 573395 | 100.0 | 94.7 | |
| 37 1,1-Dichloroethane | 63 | 4.925 | 4.928 | -0.003 | 95 | 843415 | 100.0 | 101.4 | |
| 38 Vinyl acetate | 43 | 4.980 | 4.982 | -0.002 | 97 | 926426 | 100.0 | 95.6 | |
| 42 2,2-Dichloropropane | 97 | 5.691 | 5.682 | 0.009 | 90 | 87515 | 100.0 | 101.6 | |
| 43 cis-1,2-Dichloroethene | 96 | 5.691 | 5.694 | -0.003 | 80 | 585924 | 100.0 | 98.1 | |
| 44 2-Butanone (MEK) | 43 | 5.704 | 5.706 | -0.002 | 98 | 488342 | 200.0 | 183.1 | |
| 48 Chlorobromomethane | 128 | 5.983 | 5.986 | -0.003 | 97 | 249508 | 100.0 | 95.7 | |
| 49 Tetrahydrofuran | 42 | 5.989 | 5.992 | -0.003 | 87 | 320475 | 200.0 | 177.5 | |
| 50 Chloroform | 83 | 6.135 | 6.138 | -0.003 | 93 | 889966 | 100.0 | 99.2 | |
| 51 1,1,1-Trichloroethane | 97 | 6.294 | 6.290 | 0.004 | 98 | 583435 | 100.0 | 102.3 | |
| 52 Cyclohexane | 56 | 6.367 | 6.363 | 0.004 | 89 | 781868 | 100.0 | 96.5 | |
| 53 Carbon tetrachloride | 117 | 6.470 | 6.467 | 0.003 | 96 | 414150 | 100.0 | 101.0 | |
| 54 1,1-Dichloropropene | 75 | 6.482 | 6.485 | -0.003 | 98 | 680030 | 100.0 | 99.0 | |
| 56 Benzene | 78 | 6.701 | 6.704 | -0.003 | 98 | 1915947 | 100.0 | 95.8 | |
| 57 1,2-Dichloroethane | 62 | 6.780 | 6.783 | -0.003 | 97 | 697923 | 100.0 | 94.5 | |
| 55 Isobutyl alcohol | 41 | 6.695 | 6.698 | -0.003 | 42 | 262400 | 2500.0 | 2181.2 | |
| 59 n-Heptane | 43 | 7.072 | 7.075 | -0.003 | 90 | 434669 | 100.0 | 94.3 | |
| 61 Trichloroethene | 130 | 7.443 | 7.446 | -0.003 | 98 | 474544 | 100.0 | 97.7 | |
| 63 Methylcyclohexane | 83 | 7.681 | 7.677 | 0.004 | 86 | 819514 | 100.0 | 96.6 | |
| 64 1,2-Dichloropropane | 63 | 7.717 | 7.720 | -0.003 | 94 | 480287 | 100.0 | 96.1 | |
| 67 Dibromomethane | 93 | 7.808 | 7.811 | -0.003 | 96 | 306535 | 100.0 | 93.8 | |
| 65 1,4-Dioxane | 88 | 7.802 | 7.805 | -0.003 | 36 | 80878 | 2000.0 | 1774.8 | M |
| 68 Dichlorobromomethane | 83 | 8.009 | 8.006 | 0.003 | 100 | 544627 | 100.0 | 97.9 | |
| 70 2-Chloroethyl vinyl ether | 63 | 8.313 | 8.316 | -0.003 | 93 | 580737 | 200.0 | 175.4 | |
| 71 cis-1,3-Dichloropropene | 75 | 8.453 | 8.456 | -0.003 | 96 | 626305 | 100.0 | 99.2 | |
| 72 4-Methyl-2-pentanone (MIBK) | 43 | 8.611 | 8.614 | -0.003 | 93 | 970750 | 200.0 | 198.3 | |
| 73 Toluene | 91 | 8.782 | 8.784 | -0.002 | 98 | 1858285 | 100.0 | 98.4 | |
| 74 trans-1,3-Dichloropropene | 75 | 9.037 | 9.040 | -0.003 | 92 | 511188 | 100.0 | 101.0 | |
| 75 Ethyl methacrylate | 69 | 9.104 | 9.101 | 0.003 | 88 | 632371 | 100.0 | 99.2 | |
| 76 1,1,2-Trichloroethane | 97 | 9.232 | 9.229 | 0.003 | 91 | 416541 | 100.0 | 95.3 | |
| 77 Tetrachloroethene | 164 | 9.299 | 9.295 | 0.004 | 96 | 329342 | 100.0 | 97.4 | |
| 78 1,3-Dichloropropane | 76 | 9.384 | 9.387 | -0.003 | 89 | 757496 | 100.0 | 95.6 | |
| 79 2-Hexanone | 43 | 9.451 | 9.454 | -0.003 | 94 | 596567 | 200.0 | 189.4 | |
| 81 Chlorodibromomethane | 129 | 9.597 | 9.600 | -0.003 | 91 | 293309 | 100.0 | 99.0 | |
| 82 Ethylene Dibromide | 107 | 9.706 | 9.709 | -0.003 | 97 | 407201 | 100.0 | 96.8 | |
| 83 3-Chlorobenzotrifluoride | 180 | 10.181 | 10.184 | -0.003 | 91 | 536353 | 100.0 | 97.7 | |
| 84 Chlorobenzene | 112 | 10.199 | 10.202 | -0.003 | 92 | 1175123 | 100.0 | 95.7 | |
| 85 4-Chlorobenzotrifluoride | 180 | 10.266 | 10.269 | -0.003 | 97 | 489696 | 100.0 | 96.2 | |
| 86 1,1,1,2-Tetrachloroethane | 131 | 10.297 | 10.293 | 0.004 | 89 | 357566 | 100.0 | 102.4 | |
| 87 Ethylbenzene | 106 | 10.303 | 10.299 | 0.004 | 98 | 684943 | 100.0 | 97.5 | |
| 88 m-Xylene & p-Xylene | 106 | 10.430 | 10.433 | -0.003 | 99 | 838371 | 100.0 | 97.4 | |
| 89 o-Xylene | 106 | 10.814 | 10.816 | -0.002 | 95 | 844701 | 100.0 | 97.9 | |
| 90 Styrene | 104 | 10.838 | 10.835 | 0.003 | 94 | 1341052 | 100.0 | 96.1 | |
| 91 Bromoform | 173 | 11.014 | 11.017 | -0.003 | 95 | 160632 | 100.0 | 96.4 | |
| 92 2-Chlorobenzotrifluoride | 180 | 11.087 | 11.090 | -0.003 | 95 | 544128 | 100.0 | 97.8 | |
| 93 Isopropylbenzene | 105 | 11.185 | 11.181 | 0.004 | 97 | 1867348 | 100.0 | 96.7 | |
| 95 Bromobenzene | 156 | 11.489 | 11.492 | -0.003 | 96 | 474286 | 100.0 | 98.8 | |
| 96 1,1,2,2-Tetrachloroethane | 83 | 11.495 | 11.498 | -0.003 | 95 | 577708 | 100.0 | 93.9 | |
| 97 trans-1,4-Dichloro-2-butene | 53 | 11.531 | 11.534 | -0.003 | 78 | 133396 | 100.0 | 94.9 | |
| 98 1,2,3-Trichloropropane | 110 | 11.550 | 11.552 | -0.002 | 85 | 194531 | 100.0 | 97.6 | |
| 99 N-Propylbenzene | 120 | 11.598 | 11.601 | -0.003 | 97 | 529403 | 100.0 | 101.3 | |
| 100 2-Chlorotoluene | 126 | 11.684 | 11.680 | 0.004 | 96 | 453885 | 100.0 | 100.3 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 101 3-Chlorotoluene | 126 | 11.750 | 11.747 | 0.003 | 96 | 478005 | 100.0 | 100.6 | |
| 102 1,3,5-Trimethylbenzene | 105 | 11.787 | 11.784 | 0.003 | 95 | 1489442 | 100.0 | 100.4 | |
| 103 4-Chlorotoluene | 126 | 11.805 | 11.808 | -0.003 | 99 | 485508 | 100.0 | 98.3 | |
| 104 tert-Butylbenzene | 119 | 12.097 | 12.094 | 0.003 | 91 | 1203013 | 100.0 | 102.1 | |
| 106 1,2,4-Trimethylbenzene | 105 | 12.152 | 12.155 | -0.003 | 98 | 1554360 | 100.0 | 100.5 | |
| 107 1,2-dichloro-4-(trifluoromethyl) | 214 | 12.201 | 12.203 | -0.002 | 97 | 340323 | 100.0 | 99.1 | |
| 108 sec-Butylbenzene | 105 | 12.316 | 12.319 | -0.003 | 95 | 1659704 | 100.0 | 100.4 | |
| 109 1,3-Dichlorobenzene | 146 | 12.432 | 12.435 | -0.003 | 96 | 850676 | 100.0 | 98.7 | |
| 110 4-Isopropyltoluene | 119 | 12.474 | 12.477 | -0.003 | 95 | 1356230 | 100.0 | 99.4 | |
| 111 1,4-Dichlorobenzene | 146 | 12.535 | 12.538 | -0.003 | 93 | 880492 | 100.0 | 98.7 | |
| 113 2,4-Dichloro-1-(trifluoromethyl) | 214 | 12.572 | 12.568 | 0.004 | 96 | 314370 | 100.0 | 97.2 | |
| 114 2,5-Dichlorobenzotrifluoride | 214 | 12.614 | 12.611 | 0.003 | 97 | 352876 | 100.0 | 98.8 | |
| 116 n-Butylbenzene | 91 | 12.882 | 12.885 | -0.003 | 96 | 1244987 | 100.0 | 100.0 | |
| 117 1,2-Dichlorobenzene | 146 | 12.894 | 12.891 | 0.003 | 96 | 797769 | 100.0 | 97.3 | |
| 118 1,2-Dibromo-3-Chloropropan | 75 | 13.679 | 13.682 | -0.003 | 82 | 66004 | 100.0 | 90.3 | |
| 119 2,4- & 2,5- & 2,6- Dichlorobenzene | 125 | 13.819 | 13.822 | -0.003 | 99 | 1522613 | 300.0 | 291.5 | |
| 121 2,3- & 3,4- Dichlorotoluene | 125 | 14.239 | 14.235 | 0.004 | 99 | 1108272 | 200.0 | 194.3 | |
| 122 1,2,4-Trichlorobenzene | 180 | 14.500 | 14.503 | -0.003 | 93 | 406868 | 100.0 | 95.0 | |
| 123 Hexachlorobutadiene | 225 | 14.646 | 14.649 | -0.003 | 97 | 126465 | 100.0 | 93.2 | |
| 124 Naphthalene | 128 | 14.762 | 14.765 | -0.003 | 98 | 1215966 | 100.0 | 93.1 | |
| 125 1,2,3-Trichlorobenzene | 180 | 14.981 | 14.984 | -0.003 | 95 | 350907 | 100.0 | 91.2 | |
| 126 2,4,5-Trichlorotoluene | 159 | 15.778 | 15.780 | -0.002 | 0 | 205495 | 100.0 | 92.9 | |
| 127 2,3,6-Trichlorotoluene | 159 | 15.881 | 15.884 | -0.003 | 97 | 188457 | 100.0 | 95.0 | |
| 146 3,4-Dichlorotoluene | 1 | 0.000 | | | | | ND | ND | |
| S 130 1,2-Dichloroethene, Total | 96 | | | | 0 | | 200.0 | 195.2 | |
| S 131 Xylenes, Total | 106 | | | | 0 | | 200.0 | 195.2 | |
| S 132 1,3-Dichloropropene, Total | 1 | | | | 0 | | 200.0 | 200.3 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

Reagents:

| | | |
|---------------------|--------------------|-----------|
| voaWEEmix1stR_00009 | Amount Added: 4.00 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 4.00 | Units: uL |
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 4.00 | Units: uL |
| voaWVA1stRest_00016 | Amount Added: 4.00 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 8.00 | Units: uL |
| VOA8260VOAPRI_00263 | Amount Added: 4.00 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 4.00 | Units: uL |

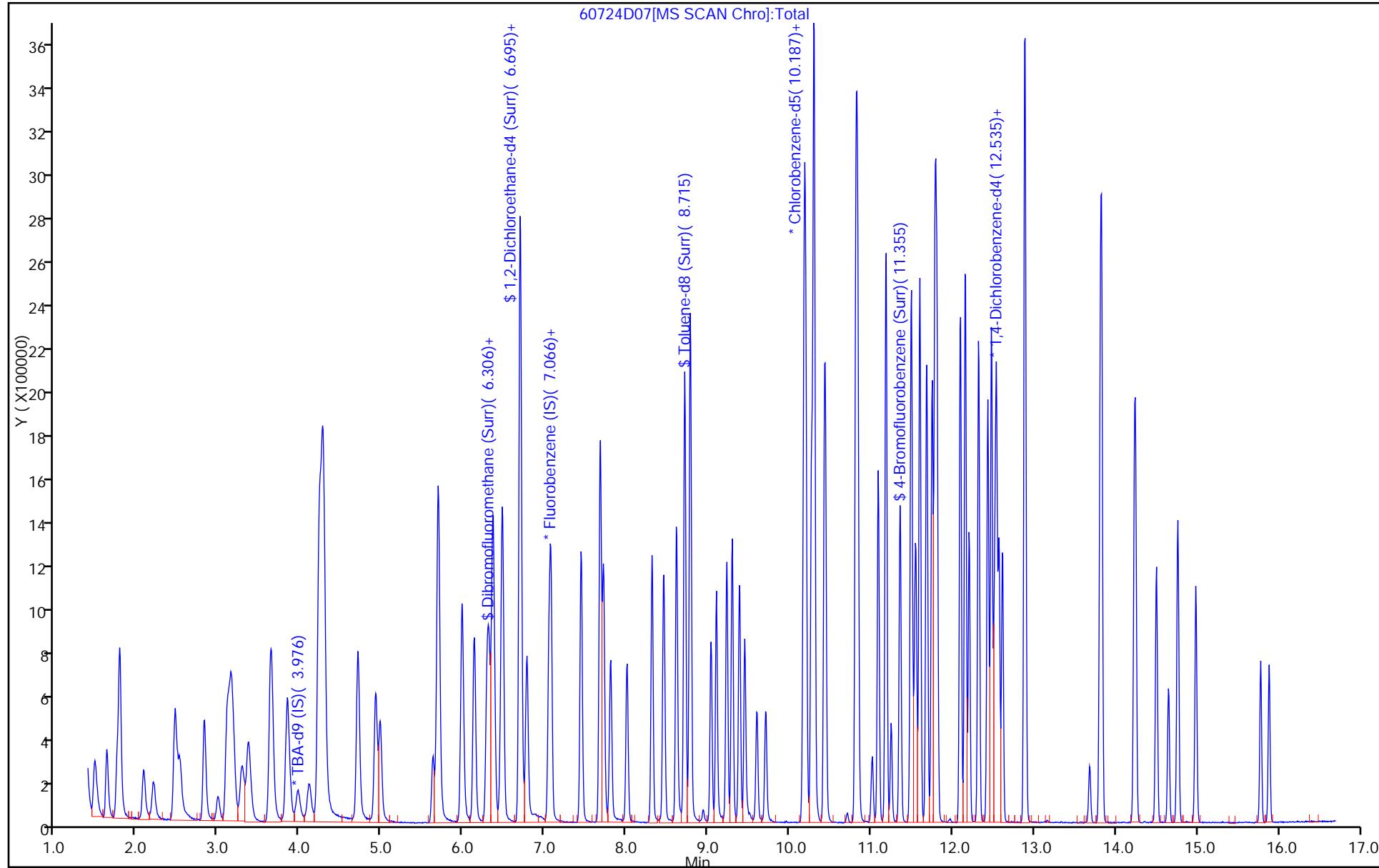
Report Date: 25-Jul-2017 01:44:38

Chrom Revision: 2.2 20-Jun-2017 07:42:38

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D07.D
Injection Date: 24-Jul-2017 08:16:30 Instrument ID: CHHP6
Lims ID: IC VSTD20 Operator ID: 034635
Client ID:
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 7
Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm)

Worklist Smp#: 7



TestAmerica Pittsburgh

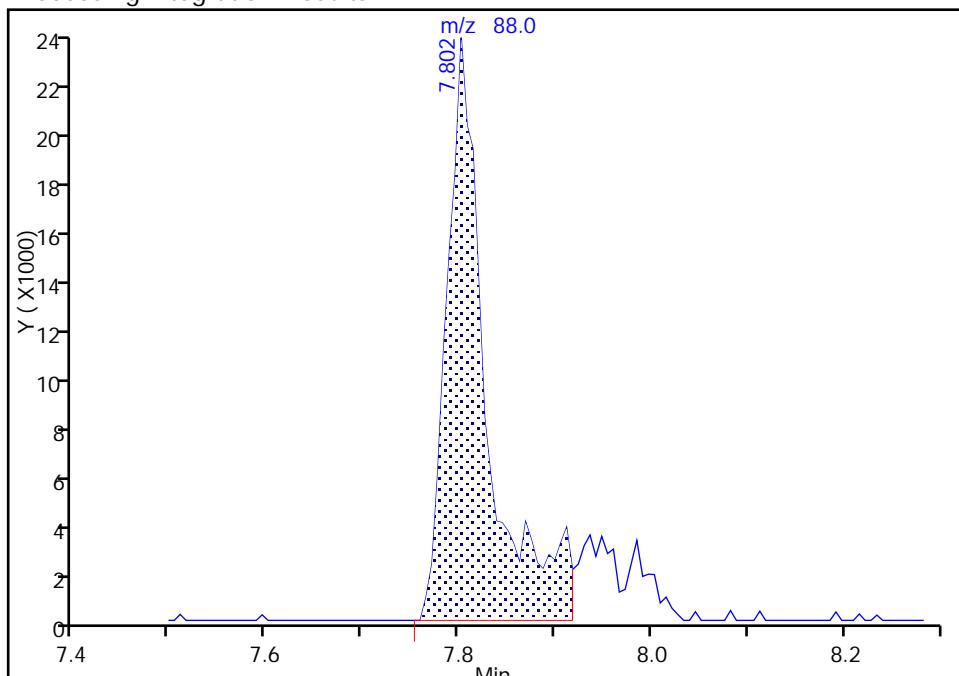
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 Injection Date: 24-Jul-2017 08:16:30 Instrument ID: CHHP6
 Lims ID: IC VSTD20
 Client ID:
 Operator ID: 034635 ALS Bottle#: 7 Worklist Smp#: 7
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
 Column: DB-624 (0.18 mm) Detector: MS SCAN

65 1,4-Dioxane, CAS: 123-91-1

Signal: 1

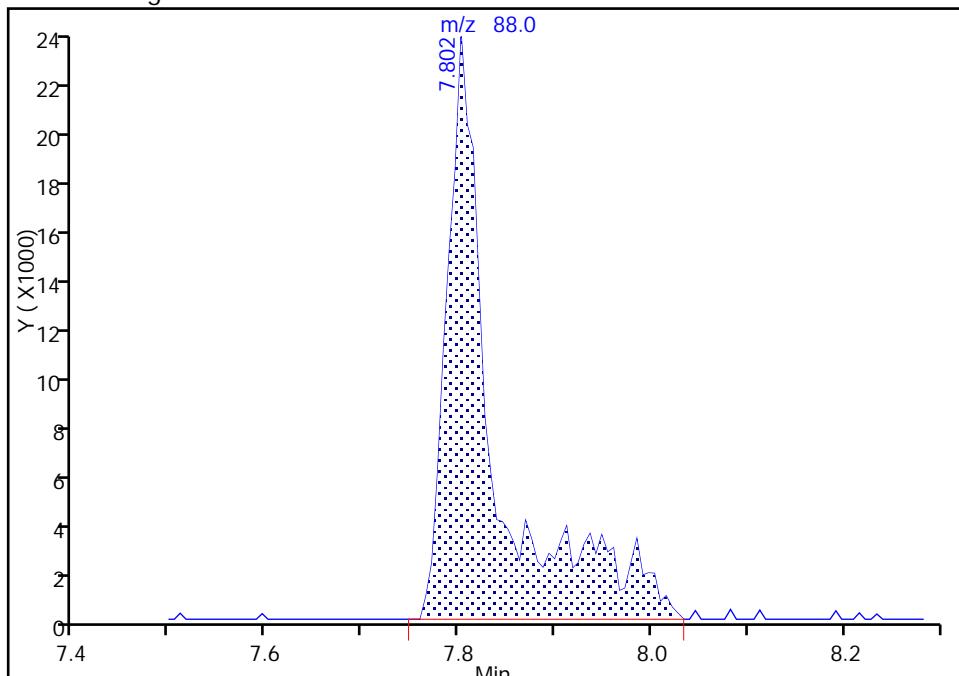
RT: 7.80
 Area: 67754
 Amount: 1445.4932
 Amount Units: ng

Processing Integration Results



RT: 7.80
 Area: 80878
 Amount: 1774.7944
 Amount Units: ng

Manual Integration Results



Reviewer: bungardf, 24-Jul-2017 08:44:40

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D08.D
 Lims ID: IC VSTD35
 Client ID:
 Sample Type: IC Calib Level: 6
 Inject. Date: 24-Jul-2017 08:40:30 ALS Bottle#: 8 Worklist Smp#: 8
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017705-008
 Misc. Info.: IC VSTD35
 Operator ID: 034635 Instrument ID: CHHP6
 Sublist: chrom-MSVOA_LL_CHHP6*sub10
 Method: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 25-Jul-2017 01:44:38 Calib Date: 24-Jul-2017 09:28:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D10.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK021

First Level Reviewer: bungardf Date: 24-Jul-2017 09:10:27

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 3.976 | 3.972 | 0.004 | 94 | 292725 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.048 | 7.051 | -0.003 | 99 | 784312 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.169 | 10.172 | -0.003 | 86 | 193416 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.511 | 12.514 | -0.003 | 96 | 277720 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr) | 113 | 6.318 | 6.315 | 0.003 | 94 | 729396 | 175.0 | 179.0 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur) | 65 | 6.689 | 6.692 | -0.003 | 70 | 1013596 | 175.0 | 174.0 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.715 | 8.718 | -0.003 | 93 | 2475360 | 175.0 | 166.2 | |
| \$ 8 4-Bromofluorobenzene (Surr) | 95 | 11.355 | 11.352 | 0.003 | 84 | 1116127 | 175.0 | 167.1 | |
| 11 Dichlorodifluoromethane | 85 | 1.488 | 1.484 | 0.004 | 99 | 839726 | 175.0 | 167.8 | |
| 12 Chloromethane | 50 | 1.634 | 1.630 | 0.004 | 98 | 792135 | 175.0 | 176.0 | |
| 13 Vinyl chloride | 62 | 1.762 | 1.758 | 0.004 | 97 | 788826 | 175.0 | 167.2 | |
| 14 Butadiene | 39 | 1.792 | 1.788 | 0.004 | 88 | 626705 | 175.0 | 160.2 | |
| 15 Bromomethane | 94 | 2.084 | 2.087 | -0.003 | 90 | 390167 | 175.0 | 177.4 | |
| 16 Chloroethane | 64 | 2.212 | 2.202 | 0.010 | 98 | 445022 | 175.0 | 174.6 | |
| 17 Dichlorofluoromethane | 67 | 2.473 | 2.470 | 0.003 | 96 | 986869 | 175.0 | 179.6 | |
| 18 Trichlorofluoromethane | 101 | 2.516 | 2.506 | 0.010 | 98 | 820933 | 175.0 | 176.7 | |
| 20 Ethyl ether | 59 | 2.826 | 2.823 | 0.003 | 87 | 720418 | 175.0 | 181.2 | |
| 21 Acrolein | 56 | 2.997 | 2.999 | -0.002 | 99 | 203936 | 225.0 | 239.5 | |
| 22 1,1-Dichloroethene | 96 | 3.118 | 3.115 | 0.003 | 98 | 719926 | 175.0 | 176.6 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.161 | 3.157 | 0.004 | 95 | 680574 | 175.0 | 174.2 | |
| 24 Acetone | 43 | 3.191 | 3.194 | -0.003 | 100 | 658887 | 350.0 | 391.6 | |
| 25 Iodomethane | 142 | 3.289 | 3.291 | -0.002 | 98 | 1045605 | 175.0 | 182.4 | |
| 26 Carbon disulfide | 76 | 3.368 | 3.364 | 0.004 | 99 | 1759930 | 175.0 | 193.9 | |
| 29 3-Chloro-1-propene | 76 | 3.635 | 3.638 | -0.003 | 91 | 457646 | 175.0 | 191.7 | |
| 30 Methyl acetate | 43 | 3.654 | 3.662 | -0.008 | 96 | 1378375 | 350.0 | 373.7 | |
| 31 Methylene Chloride | 84 | 3.842 | 3.839 | 0.003 | 88 | 940505 | 175.0 | 174.1 | |
| 32 2-Methyl-2-propanol | 59 | 4.116 | 4.112 | 0.004 | 94 | 572173 | 1750.0 | 1765.9 | |
| 33 Acrylonitrile | 53 | 4.238 | 4.240 | -0.002 | 97 | 3519987 | 1750.0 | 1786.6 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.274 | 4.277 | -0.003 | 99 | 817284 | 175.0 | 176.7 | |
| 35 Methyl tert-butyl ether | 73 | 4.286 | 4.295 | -0.009 | 96 | 2689634 | 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 | 4.712 | 4.709 | 0.003 | 91 | 940765 | 175.0 | 171.7 | |
| 37 1,1-Dichloroethane | 63 | 4.931 | 4.928 | 0.003 | 95 | 1429058 | 175.0 | 189.9 | |
| 38 Vinyl acetate | 43 | 4.980 | 4.982 | -0.002 | 97 | 1612184 | 175.0 | 184.0 | |
| 42 2,2-Dichloropropane | 97 | 5.692 | 5.682 | 0.010 | 92 | 147487 | 175.0 | 189.3 | |
| 43 cis-1,2-Dichloroethene | 96 | 5.692 | 5.694 | -0.002 | 80 | 997518 | 175.0 | 184.6 | |
| 44 2-Butanone (MEK) | 43 | 5.704 | 5.706 | -0.002 | 98 | 909301 | 350.0 | 377.0 | |
| 48 Chlorobromomethane | 128 | 5.984 | 5.986 | -0.002 | 97 | 441109 | 175.0 | 187.1 | |
| 49 Tetrahydrofuran | 42 | 5.990 | 5.992 | -0.002 | 87 | 565011 | 350.0 | 346.0 | |
| 50 Chloroform | 83 | 6.136 | 6.138 | -0.002 | 93 | 1486297 | 175.0 | 183.2 | |
| 51 1,1,1-Trichloroethane | 97 | 6.294 | 6.290 | 0.004 | 97 | 954279 | 175.0 | 185.1 | |
| 52 Cyclohexane | 56 | 6.361 | 6.363 | -0.002 | 89 | 1239342 | 175.0 | 169.1 | |
| 53 Carbon tetrachloride | 117 | 6.464 | 6.467 | -0.003 | 96 | 707925 | 175.0 | 190.9 | |
| 54 1,1-Dichloropropene | 75 | 6.482 | 6.485 | -0.003 | 97 | 1104618 | 175.0 | 177.8 | |
| 55 Isobutyl alcohol | 41 | 6.701 | 6.698 | 0.003 | 46 | 537905 | 4375.0 | 4944.3 | |
| 56 Benzene | 78 | 6.701 | 6.704 | -0.003 | 98 | 3132477 | 175.0 | 173.3 | |
| 57 1,2-Dichloroethane | 62 | 6.781 | 6.783 | -0.002 | 97 | 1211453 | 175.0 | 181.4 | |
| 59 n-Heptane | 43 | 7.073 | 7.075 | -0.002 | 89 | 728795 | 175.0 | 174.8 | |
| 61 Trichloroethene | 130 | 7.444 | 7.446 | -0.002 | 98 | 799778 | 175.0 | 182.1 | |
| 63 Methylcyclohexane | 83 | 7.681 | 7.677 | 0.004 | 87 | 1300055 | 175.0 | 169.5 | |
| 64 1,2-Dichloropropane | 63 | 7.717 | 7.720 | -0.003 | 94 | 853109 | 175.0 | 188.8 | |
| 65 1,4-Dioxane | 88 | 7.796 | 7.805 | -0.009 | 39 | 141286 | 3500.0 | 3428.3 | |
| 67 Dibromomethane | 93 | 7.803 | 7.811 | -0.008 | 97 | 555898 | 175.0 | 188.1 | |
| 68 Dichlorobromomethane | 83 | 8.003 | 8.006 | -0.003 | 99 | 1013623 | 175.0 | 201.4 | |
| 70 2-Chloroethyl vinyl ether | 63 | 8.314 | 8.316 | -0.002 | 93 | 1138561 | 350.0 | 380.2 | |
| 71 cis-1,3-Dichloropropene | 75 | 8.453 | 8.456 | -0.003 | 96 | 1210517 | 175.0 | 212.1 | |
| 72 4-Methyl-2-pentanone (MIBK) | 43 | 8.612 | 8.614 | -0.002 | 92 | 1737974 | 350.0 | 351.9 | |
| 73 Toluene | 91 | 8.782 | 8.784 | -0.002 | 96 | 3088570 | 175.0 | 162.0 | |
| 74 trans-1,3-Dichloropropene | 75 | 9.037 | 9.040 | -0.003 | 92 | 1035772 | 175.0 | 202.9 | |
| 75 Ethyl methacrylate | 69 | 9.104 | 9.101 | 0.003 | 88 | 1196636 | 175.0 | 186.0 | |
| 76 1,1,2-Trichloroethane | 97 | 9.232 | 9.229 | 0.003 | 91 | 786936 | 175.0 | 178.4 | |
| 77 Tetrachloroethene | 164 | 9.299 | 9.295 | 0.004 | 96 | 574638 | 175.0 | 168.4 | |
| 78 1,3-Dichloropropane | 76 | 9.384 | 9.387 | -0.003 | 89 | 1388403 | 175.0 | 173.6 | |
| 79 2-Hexanone | 43 | 9.451 | 9.454 | -0.003 | 91 | 1122326 | 350.0 | 353.1 | |
| 81 Chlorodibromomethane | 129 | 9.597 | 9.600 | -0.003 | 91 | 614068 | 175.0 | 205.5 | |
| 82 Ethylene Dibromide | 107 | 9.707 | 9.709 | -0.002 | 98 | 780325 | 175.0 | 183.9 | |
| 83 3-Chlorobenzotrifluoride | 180 | 10.181 | 10.184 | -0.003 | 92 | 925971 | 175.0 | 167.1 | |
| 84 Chlorobenzene | 112 | 10.199 | 10.202 | -0.003 | 92 | 2089428 | 175.0 | 168.6 | |
| 85 4-Chlorobenzotrifluoride | 180 | 10.272 | 10.269 | 0.003 | 96 | 878791 | 175.0 | 171.1 | |
| 86 1,1,1,2-Tetrachloroethane | 131 | 10.297 | 10.293 | 0.004 | 92 | 699176 | 175.0 | 198.4 | |
| 87 Ethylbenzene | 106 | 10.303 | 10.299 | 0.004 | 97 | 1218751 | 175.0 | 171.9 | |
| 88 m-Xylene & p-Xylene | 106 | 10.437 | 10.433 | 0.004 | 97 | 1499275 | 175.0 | 172.5 | |
| 89 o-Xylene | 106 | 10.814 | 10.816 | -0.002 | 94 | 1520782 | 175.0 | 174.6 | |
| 90 Styrene | 104 | 10.838 | 10.835 | 0.003 | 93 | 2430770 | 175.0 | 172.6 | |
| 91 Bromoform | 173 | 11.015 | 11.017 | -0.002 | 95 | 360803 | 175.0 | 214.6 | |
| 92 2-Chlorobenzotrifluoride | 180 | 11.088 | 11.090 | -0.002 | 94 | 976474 | 175.0 | 173.9 | |
| 93 Isopropylbenzene | 105 | 11.185 | 11.181 | 0.004 | 97 | 3139141 | 175.0 | 161.1 | |
| 95 Bromobenzene | 156 | 11.489 | 11.492 | -0.003 | 96 | 913147 | 175.0 | 176.9 | |
| 96 1,1,2,2-Tetrachloroethane | 83 | 11.501 | 11.498 | 0.003 | 95 | 1109880 | 175.0 | 178.8 | |
| 97 trans-1,4-Dichloro-2-butene | 53 | 11.538 | 11.534 | 0.004 | 80 | 280292 | 175.0 | 185.5 | |
| 98 1,2,3-Trichloropropane | 110 | 11.550 | 11.552 | -0.002 | 87 | 379404 | 175.0 | 177.0 | |
| 99 N-Propylbenzene | 120 | 11.599 | 11.601 | -0.002 | 96 | 962443 | 175.0 | 171.3 | |
| 100 2-Chlorotoluene | 126 | 11.684 | 11.680 | 0.004 | 96 | 862193 | 175.0 | 177.2 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 101 3-Chlorotoluene | 126 | 11.751 | 11.747 | 0.004 | 95 | 877299 | 175.0 | 171.8 | |
| 102 1,3,5-Trimethylbenzene | 105 | 11.787 | 11.784 | 0.003 | 96 | 2680034 | 175.0 | 168.0 | |
| 103 4-Chlorotoluene | 126 | 11.805 | 11.808 | -0.003 | 98 | 940075 | 175.0 | 177.0 | |
| 104 tert-Butylbenzene | 119 | 12.097 | 12.094 | 0.003 | 91 | 2145902 | 175.0 | 169.4 | |
| 106 1,2,4-Trimethylbenzene | 105 | 12.158 | 12.155 | 0.003 | 97 | 2764298 | 175.0 | 166.2 | |
| 107 1,2-dichloro-4-(trifluoromethyl) | 214 | 12.207 | 12.203 | 0.004 | 95 | 632603 | 175.0 | 171.4 | |
| 108 sec-Butylbenzene | 105 | 12.316 | 12.319 | -0.003 | 95 | 2889256 | 175.0 | 162.6 | |
| 109 1,3-Dichlorobenzene | 146 | 12.432 | 12.435 | -0.003 | 94 | 1574554 | 175.0 | 169.9 | |
| 110 4-Isopropyltoluene | 119 | 12.475 | 12.477 | -0.002 | 93 | 2441419 | 175.0 | 166.4 | |
| 111 1,4-Dichlorobenzene | 146 | 12.535 | 12.538 | -0.003 | 91 | 1638813 | 175.0 | 171.0 | |
| 113 2,4-Dichloro-1-(trifluoromethyl) | 214 | 12.572 | 12.568 | 0.004 | 95 | 599851 | 175.0 | 172.5 | |
| 114 2,5-Dichlorobenzotrifluoride | 214 | 12.615 | 12.611 | 0.004 | 97 | 682948 | 175.0 | 177.9 | |
| 116 n-Butylbenzene | 91 | 12.882 | 12.885 | -0.003 | 96 | 2222409 | 175.0 | 166.1 | |
| 117 1,2-Dichlorobenzene | 146 | 12.888 | 12.891 | -0.003 | 94 | 1520936 | 175.0 | 172.5 | |
| 118 1,2-Dibromo-3-Chloropropan | 75 | 13.679 | 13.682 | -0.003 | 81 | 156142 | 175.0 | 198.7 | |
| 119 2,4- & 2,5- & 2,6- Dichlorobenzene | 125 | 13.819 | 13.822 | -0.003 | 97 | 2790747 | 525.0 | 497.0 | |
| 121 2,3- & 3,4- Dichlorotoluene | 125 | 14.239 | 14.235 | 0.004 | 98 | 2076832 | 350.0 | 338.6 | |
| 122 1,2,4-Trichlorobenzene | 180 | 14.500 | 14.503 | -0.003 | 94 | 795349 | 175.0 | 172.6 | |
| 123 Hexachlorobutadiene | 225 | 14.646 | 14.649 | -0.003 | 97 | 239351 | 175.0 | 164.1 | |
| 124 Naphthalene | 128 | 14.762 | 14.765 | -0.003 | 99 | 2281539 | 175.0 | 162.5 | |
| 125 1,2,3-Trichlorobenzene | 180 | 14.981 | 14.984 | -0.003 | 95 | 706689 | 175.0 | 170.9 | |
| 126 2,4,5-Trichlorotoluene | 159 | 15.778 | 15.780 | -0.002 | 0 | 413111 | 175.0 | 173.7 | |
| 127 2,3,6-Trichlorotoluene | 159 | 15.881 | 15.884 | -0.003 | 96 | 356014 | 175.0 | 167.0 | |
| 146 3,4-Dichlorotoluene | 1 | 0.000 | | | | | ND | ND | |
| S 131 Xylenes, Total | 106 | | | | 0 | | 350.0 | 347.2 | |
| S 130 1,2-Dichloroethene, Total | 96 | | | | 0 | | 350.0 | 361.3 | |
| S 132 1,3-Dichloropropene, Total | 1 | | | | 0 | | 350.0 | 415.0 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Reagents:

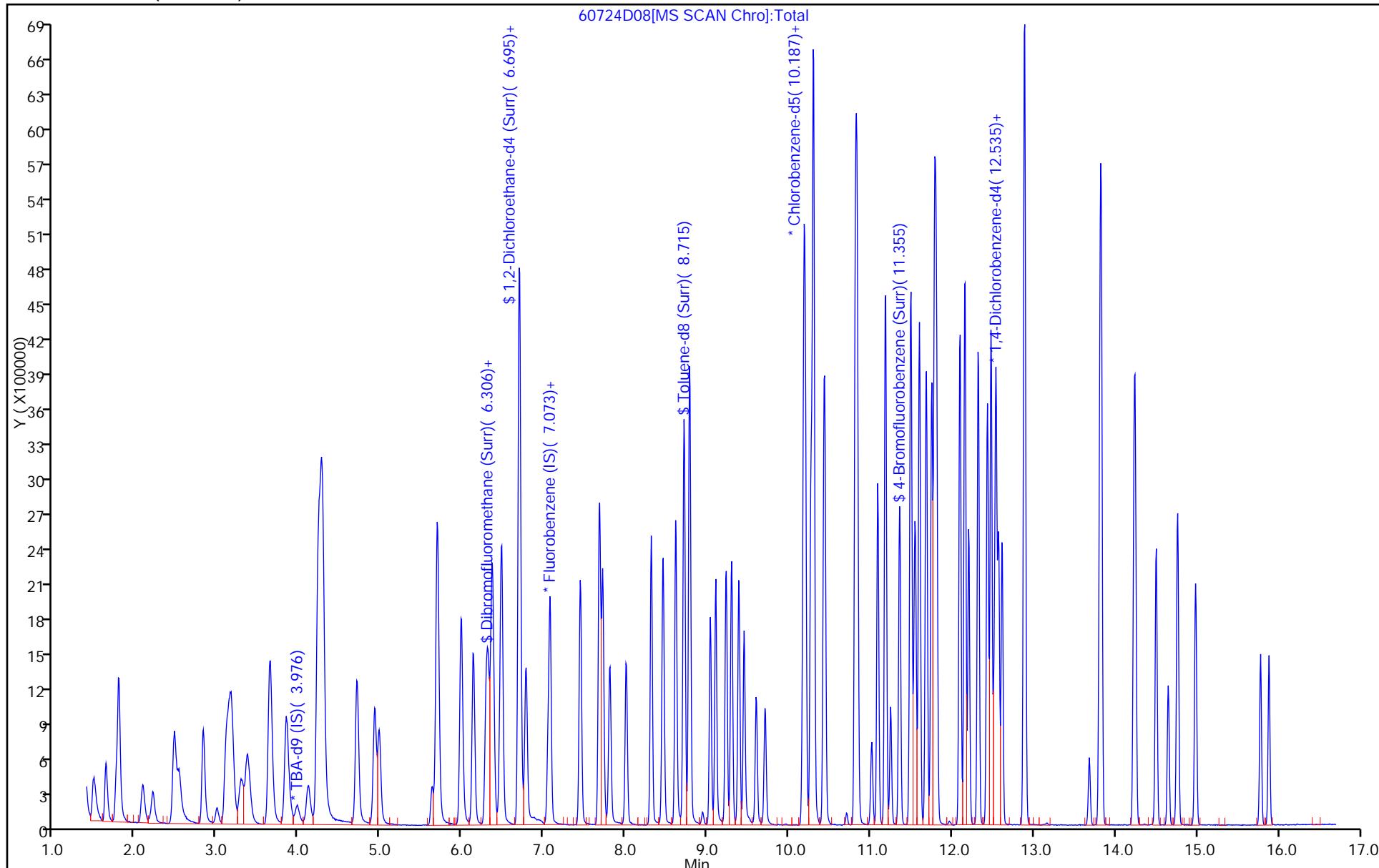
| | | |
|---------------------|--------------------|-----------|
| voaWEEmix1stR_00009 | Amount Added: 7.00 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 7.00 | Units: uL |
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 7.00 | Units: uL |
| voaWVA1stRest_00016 | Amount Added: 7.00 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 9.00 | Units: uL |
| VOA8260VOAPRI_00263 | Amount Added: 7.00 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 7.00 | Units: uL |

Report Date: 25-Jul-2017 01:44:40

Chrom Revision: 2.2 20-Jun-2017 07:42:38

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D08.D
Injection Date: 24-Jul-2017 08:40:30 Instrument ID: CHHP6
Lims ID: IC VSTD35 Operator ID: 034635
Client ID:
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 8
Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D09.D
 Lims ID: IC VSTD40
 Client ID:
 Sample Type: IC Calib Level: 7
 Inject. Date: 24-Jul-2017 09:04:30 ALS Bottle#: 9 Worklist Smp#: 9
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017705-009
 Misc. Info.: IC VSTD40
 Operator ID: 034635 Instrument ID: CHHP6
 Sublist: chrom-MSVOA_LL_CHHP6*sub10
 Method: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 25-Jul-2017 01:44:40 Calib Date: 24-Jul-2017 09:28:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D10.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK021

First Level Reviewer: bungardf Date: 24-Jul-2017 09:37:55

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 3.979 | 3.972 | 0.007 | 97 | 231119 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.051 | 7.051 | 0.000 | 98 | 870105 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.172 | 10.172 | 0.000 | 88 | 207878 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.514 | 12.514 | 0.000 | 97 | 276299 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr) | 113 | 6.321 | 6.315 | 0.006 | 94 | 842973 | 200.0 | 186.4 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur) | 65 | 6.692 | 6.692 | 0.000 | 70 | 1135703 | 200.0 | 175.7 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.718 | 8.718 | 0.000 | 93 | 2824683 | 200.0 | 176.6 | |
| \$ 8 4-Bromofluorobenzene (Surr) | 95 | 11.358 | 11.352 | 0.006 | 88 | 1222775 | 200.0 | 170.4 | |
| 11 Dichlorodifluoromethane | 85 | 1.484 | 1.484 | 0.000 | 99 | 998454 | 200.0 | 179.8 | |
| 12 Chloromethane | 50 | 1.630 | 1.630 | 0.000 | 98 | 904618 | 200.0 | 181.2 | |
| 13 Vinyl chloride | 62 | 1.764 | 1.758 | 0.006 | 98 | 942517 | 200.0 | 180.1 | |
| 14 Butadiene | 39 | 1.789 | 1.788 | 0.001 | 88 | 738623 | 200.0 | 170.2 | |
| 15 Bromomethane | 94 | 2.081 | 2.087 | -0.006 | 90 | 421777 | 200.0 | 172.9 | |
| 16 Chloroethane | 64 | 2.202 | 2.202 | 0.000 | 98 | 496292 | 200.0 | 175.5 | |
| 17 Dichlorofluoromethane | 67 | 2.464 | 2.470 | -0.006 | 96 | 1112294 | 200.0 | 182.5 | |
| 18 Trichlorofluoromethane | 101 | 2.506 | 2.506 | 0.000 | 97 | 958134 | 200.0 | 185.9 | |
| 20 Ethyl ether | 59 | 2.829 | 2.823 | 0.006 | 87 | 817030 | 200.0 | 185.2 | |
| 21 Acrolein | 56 | 2.993 | 2.999 | -0.006 | 98 | 222222 | 250.0 | 235.2 | |
| 22 1,1-Dichloroethene | 96 | 3.109 | 3.115 | -0.006 | 97 | 850942 | 200.0 | 188.2 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.157 | 3.157 | 0.000 | 95 | 799492 | 200.0 | 184.5 | |
| 24 Acetone | 43 | 3.194 | 3.194 | 0.000 | 98 | 631699 | 400.0 | 338.4 | |
| 25 Iodomethane | 142 | 3.285 | 3.291 | -0.006 | 98 | 1229211 | 200.0 | 193.3 | |
| 26 Carbon disulfide | 76 | 3.370 | 3.364 | 0.006 | 99 | 2106147 | 200.0 | 209.1 | |
| 29 3-Chloro-1-propene | 76 | 3.632 | 3.638 | -0.006 | 92 | 540192 | 200.0 | 204.0 | |
| 30 Methyl acetate | 43 | 3.650 | 3.662 | -0.012 | 96 | 1509491 | 400.0 | 368.9 | |
| 31 Methylene Chloride | 84 | 3.845 | 3.839 | 0.006 | 88 | 1081026 | 200.0 | 180.4 | |
| 32 2-Methyl-2-propanol | 59 | 4.119 | 4.112 | 0.006 | 94 | 544523 | 2000.0 | 2128.5 | |
| 33 Acrylonitrile | 53 | 4.234 | 4.240 | -0.006 | 97 | 3825550 | 2000.0 | 1750.2 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.277 | 4.277 | 0.000 | 98 | 968185 | 200.0 | 188.7 | |
| 35 Methyl tert-butyl ether | 73 | 4.295 | 4.295 | 0.000 | 96 | 3053893 | 200.0 | 186.1 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 36 Hexane | 57 | 4.709 | 4.709 | 0.000 | 90 | 1103506 | 200.0 | 181.6 | |
| 37 1,1-Dichloroethane | 63 | 4.928 | 4.928 | 0.000 | 95 | 1631648 | 200.0 | 195.5 | |
| 38 Vinyl acetate | 43 | 4.982 | 4.982 | 0.000 | 97 | 1962237 | 200.0 | 201.8 | |
| 42 2,2-Dichloropropane | 97 | 5.694 | 5.682 | 0.012 | 87 | 168555 | 200.0 | 195.0 | |
| 43 cis-1,2-Dichloroethene | 96 | 5.694 | 5.694 | 0.000 | 80 | 1143732 | 200.0 | 190.8 | |
| 44 2-Butanone (MEK) | 43 | 5.706 | 5.706 | 0.000 | 98 | 973759 | 400.0 | 363.9 | |
| 48 Chlorobromomethane | 128 | 5.980 | 5.986 | -0.006 | 98 | 507483 | 200.0 | 194.0 | |
| 49 Tetrahydrofuran | 42 | 5.992 | 5.992 | 0.000 | 84 | 629732 | 400.0 | 347.6 | |
| 50 Chloroform | 83 | 6.132 | 6.138 | -0.006 | 92 | 1701079 | 200.0 | 189.0 | |
| 51 1,1,1-Trichloroethane | 97 | 6.296 | 6.290 | 0.006 | 98 | 1112586 | 200.0 | 194.5 | |
| 52 Cyclohexane | 56 | 6.363 | 6.363 | 0.000 | 89 | 1473582 | 200.0 | 181.3 | |
| 53 Carbon tetrachloride | 117 | 6.467 | 6.467 | 0.000 | 96 | 848801 | 200.0 | 206.3 | |
| 54 1,1-Dichloropropene | 75 | 6.485 | 6.485 | 0.000 | 97 | 1298950 | 200.0 | 188.5 | |
| 55 Isobutyl alcohol | 41 | 6.698 | 6.698 | 0.000 | 88 | 631529 | 5000.0 | 5232.5 | |
| 56 Benzene | 78 | 6.698 | 6.704 | -0.006 | 98 | 3551507 | 200.0 | 177.1 | |
| 57 1,2-Dichloroethane | 62 | 6.783 | 6.783 | 0.000 | 97 | 1374193 | 200.0 | 185.5 | |
| 59 n-Heptane | 43 | 7.075 | 7.075 | 0.000 | 89 | 834225 | 200.0 | 180.4 | |
| 61 Trichloroethene | 130 | 7.446 | 7.446 | 0.000 | 98 | 926685 | 200.0 | 190.2 | |
| 63 Methylcyclohexane | 83 | 7.677 | 7.677 | 0.000 | 87 | 1530213 | 200.0 | 179.9 | |
| 64 1,2-Dichloropropane | 63 | 7.720 | 7.720 | 0.000 | 94 | 970394 | 200.0 | 193.5 | |
| 65 1,4-Dioxane | 88 | 7.805 | 7.805 | 0.000 | 56 | 152489 | 4000.0 | 3335.3 | M |
| 67 Dibromomethane | 93 | 7.805 | 7.811 | -0.006 | 97 | 636427 | 200.0 | 194.1 | |
| 68 Dichlorobromomethane | 83 | 8.006 | 8.006 | 0.000 | 99 | 1181877 | 200.0 | 211.7 | |
| 70 2-Chloroethyl vinyl ether | 63 | 8.316 | 8.316 | 0.000 | 93 | 1300177 | 400.0 | 391.4 | |
| 71 cis-1,3-Dichloropropene | 75 | 8.456 | 8.456 | 0.000 | 96 | 1391254 | 200.0 | 219.7 | |
| 72 4-Methyl-2-pentanone (MIBK) | 43 | 8.614 | 8.614 | 0.000 | 91 | 1932325 | 400.0 | 364.0 | |
| 73 Toluene | 91 | 8.785 | 8.784 | 0.001 | 95 | 3464609 | 200.0 | 169.1 | |
| 74 trans-1,3-Dichloropropene | 75 | 9.040 | 9.040 | 0.000 | 92 | 1215519 | 200.0 | 221.6 | |
| 75 Ethyl methacrylate | 69 | 9.101 | 9.101 | 0.000 | 88 | 1382390 | 200.0 | 200.0 | |
| 76 1,1,2-Trichloroethane | 97 | 9.229 | 9.229 | 0.000 | 91 | 896840 | 200.0 | 189.2 | |
| 77 Tetrachloroethene | 164 | 9.296 | 9.295 | 0.001 | 95 | 670325 | 200.0 | 182.8 | |
| 78 1,3-Dichloropropane | 76 | 9.387 | 9.387 | 0.000 | 89 | 1574600 | 200.0 | 183.2 | |
| 79 2-Hexanone | 43 | 9.448 | 9.454 | -0.006 | 92 | 1271094 | 400.0 | 372.0 | |
| 81 Chlorodibromomethane | 129 | 9.600 | 9.600 | 0.000 | 93 | 712324 | 200.0 | 221.8 | |
| 82 Ethylene Dibromide | 107 | 9.709 | 9.709 | 0.000 | 98 | 899338 | 200.0 | 197.2 | |
| 83 3-Chlorobenzotrifluoride | 180 | 10.184 | 10.184 | 0.000 | 91 | 1082251 | 200.0 | 181.8 | |
| 84 Chlorobenzene | 112 | 10.202 | 10.202 | 0.000 | 91 | 2335758 | 200.0 | 175.3 | |
| 85 4-Chlorobenzotrifluoride | 180 | 10.269 | 10.269 | 0.000 | 96 | 1026977 | 200.0 | 186.1 | |
| 86 1,1,1,2-Tetrachloroethane | 131 | 10.299 | 10.293 | 0.006 | 92 | 799872 | 200.0 | 211.2 | |
| 87 Ethylbenzene | 106 | 10.305 | 10.299 | 0.006 | 96 | 1390812 | 200.0 | 182.5 | |
| 88 m-Xylene & p-Xylene | 106 | 10.433 | 10.433 | 0.000 | 96 | 1695741 | 200.0 | 181.6 | |
| 89 o-Xylene | 106 | 10.816 | 10.816 | 0.000 | 94 | 1672026 | 200.0 | 178.7 | |
| 90 Styrene | 104 | 10.835 | 10.835 | 0.000 | 93 | 2669824 | 200.0 | 176.4 | |
| 91 Bromoform | 173 | 11.011 | 11.017 | -0.006 | 95 | 416604 | 200.0 | 230.5 | |
| 92 2-Chlorobenzotrifluoride | 180 | 11.090 | 11.090 | 0.000 | 94 | 1099473 | 200.0 | 182.2 | |
| 93 Isopropylbenzene | 105 | 11.181 | 11.181 | 0.000 | 97 | 3385367 | 200.0 | 161.6 | |
| 95 Bromobenzene | 156 | 11.492 | 11.492 | 0.000 | 97 | 1008153 | 200.0 | 196.3 | |
| 96 1,1,2,2-Tetrachloroethane | 83 | 11.498 | 11.498 | 0.000 | 95 | 1216769 | 200.0 | 182.4 | |
| 97 trans-1,4-Dichloro-2-butene | 53 | 11.534 | 11.534 | 0.000 | 80 | 313051 | 200.0 | 208.3 | |
| 98 1,2,3-Trichloropropane | 110 | 11.553 | 11.552 | 0.000 | 85 | 413676 | 200.0 | 194.0 | |
| 99 N-Propylbenzene | 120 | 11.601 | 11.601 | 0.000 | 95 | 1050858 | 200.0 | 188.0 | |
| 100 2-Chlorotoluene | 126 | 11.680 | 11.680 | 0.000 | 96 | 919859 | 200.0 | 190.0 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 101 3-Chlorotoluene | 126 | 11.747 | 11.747 | 0.000 | 94 | 965183 | 200.0 | 190.0 | |
| 102 1,3,5-Trimethylbenzene | 105 | 11.784 | 11.784 | 0.000 | 94 | 2803848 | 200.0 | 176.7 | |
| 103 4-Chlorotoluene | 126 | 11.808 | 11.808 | 0.000 | 98 | 999834 | 200.0 | 189.2 | |
| 104 tert-Butylbenzene | 119 | 12.094 | 12.094 | 0.000 | 91 | 2233746 | 200.0 | 177.3 | |
| 106 1,2,4-Trimethylbenzene | 105 | 12.155 | 12.155 | 0.000 | 97 | 2867121 | 200.0 | 173.3 | |
| 107 1,2-dichloro-4-(trifluoromethyl) | 214 | 12.203 | 12.203 | 0.000 | 96 | 684446 | 200.0 | 186.4 | |
| 108 sec-Butylbenzene | 105 | 12.319 | 12.319 | 0.000 | 95 | 2979536 | 200.0 | 168.6 | |
| 109 1,3-Dichlorobenzene | 146 | 12.435 | 12.435 | 0.000 | 94 | 1673821 | 200.0 | 181.5 | |
| 110 4-Isopropyltoluene | 119 | 12.477 | 12.477 | 0.000 | 94 | 2506437 | 200.0 | 171.8 | |
| 111 1,4-Dichlorobenzene | 146 | 12.538 | 12.538 | 0.000 | 91 | 1720245 | 200.0 | 180.4 | |
| 113 2,4-Dichloro-1-(trifluoromethyl) | 214 | 12.568 | 12.568 | 0.000 | 94 | 621645 | 200.0 | 179.7 | |
| 114 2,5-Dichlorobenzotrifluoride | 214 | 12.611 | 12.611 | 0.000 | 97 | 758780 | 200.0 | 198.7 | |
| 116 n-Butylbenzene | 91 | 12.885 | 12.885 | 0.000 | 94 | 2290785 | 200.0 | 172.1 | |
| 117 1,2-Dichlorobenzene | 146 | 12.891 | 12.891 | 0.000 | 95 | 1586590 | 200.0 | 180.9 | |
| 118 1,2-Dibromo-3-Chloropropan | 75 | 13.682 | 13.682 | 0.000 | 88 | 173322 | 200.0 | 221.7 | |
| 119 2,4- & 2,5- & 2,6- Dichlorobenzene | 125 | 13.822 | 13.822 | 0.000 | 96 | 3071993 | 600.0 | 549.9 | |
| 121 2,3- & 3,4- Dichlorotoluene | 125 | 14.235 | 14.235 | 0.000 | 97 | 2379503 | 400.0 | 389.9 | |
| 122 1,2,4-Trichlorobenzene | 180 | 14.503 | 14.503 | 0.000 | 94 | 937825 | 200.0 | 204.6 | |
| 123 Hexachlorobutadiene | 225 | 14.649 | 14.649 | 0.000 | 97 | 285364 | 200.0 | 196.7 | |
| 124 Naphthalene | 128 | 14.758 | 14.765 | -0.007 | 99 | 2669188 | 200.0 | 191.1 | |
| 125 1,2,3-Trichlorobenzene | 180 | 14.984 | 14.984 | 0.000 | 95 | 901210 | 200.0 | 219.0 | |
| 126 2,4,5-Trichlorotoluene | 159 | 15.781 | 15.780 | 0.000 | 0 | 580730 | 200.0 | 245.4 | |
| 127 2,3,6-Trichlorotoluene | 159 | 15.884 | 15.884 | 0.000 | 95 | 503740 | 200.0 | 237.4 | |
| 146 3,4-Dichlorotoluene | 1 | 0.000 | | | | | ND | ND | |
| S 130 1,2-Dichloroethene, Total | 96 | | | | 0 | | 400.0 | 379.5 | |
| S 131 Xylenes, Total | 106 | | | | 0 | | 400.0 | 360.2 | |
| S 132 1,3-Dichloropropene, Total | 1 | | | | 0 | | 400.0 | 441.2 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

Reagents:

| | | |
|---------------------|---------------------|-----------|
| voaWEEmix1stR_00009 | Amount Added: 8.00 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 8.00 | Units: uL |
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 8.00 | Units: uL |
| voaWVA1stRest_00016 | Amount Added: 8.00 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 10.00 | Units: uL |
| VOA8260VOAPRI_00263 | Amount Added: 8.00 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 8.00 | Units: uL |

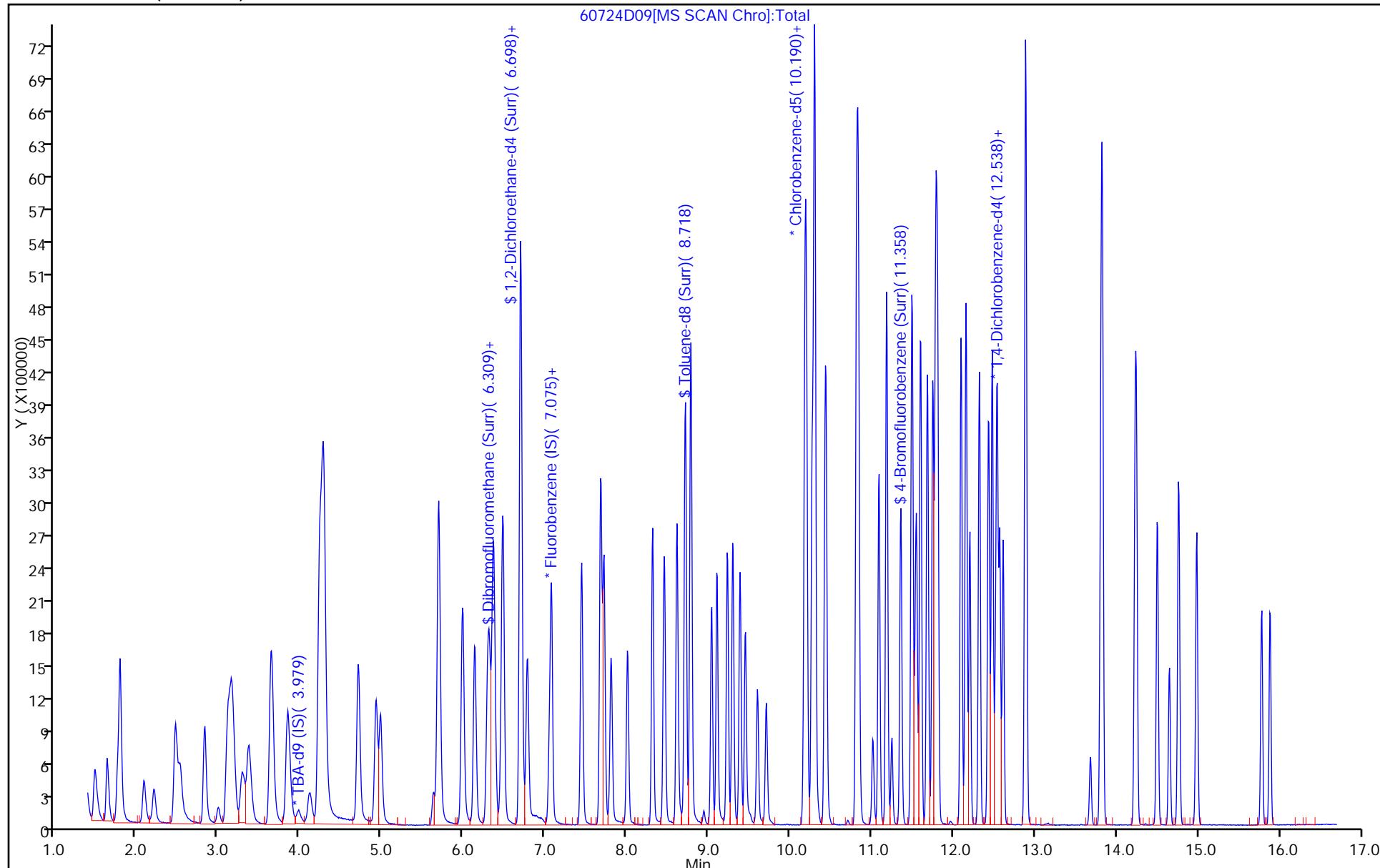
Report Date: 25-Jul-2017 01:44:42

Chrom Revision: 2.2 20-Jun-2017 07:42:38

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D09.D
Injection Date: 24-Jul-2017 09:04:30 Instrument ID: CHHP6
Lims ID: IC VSTD40 Operator ID: 034635
Client ID:
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 9
Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm)

Worklist Smp#: 9



TestAmerica Pittsburgh

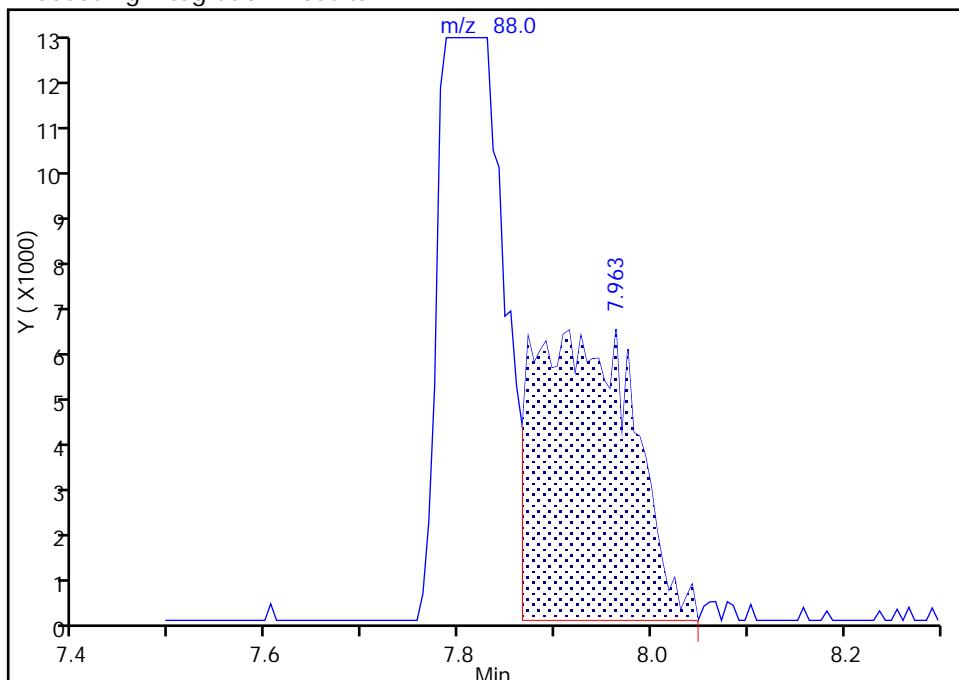
Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D09.D
 Injection Date: 24-Jul-2017 09:04:30 Instrument ID: CHHP6
 Lims ID: IC VSTD40
 Client ID:
 Operator ID: 034635 ALS Bottle#: 9 Worklist Smp#: 9
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
 Column: DB-624 (0.18 mm) Detector: MS SCAN

65 1,4-Dioxane, CAS: 123-91-1

Signal: 1

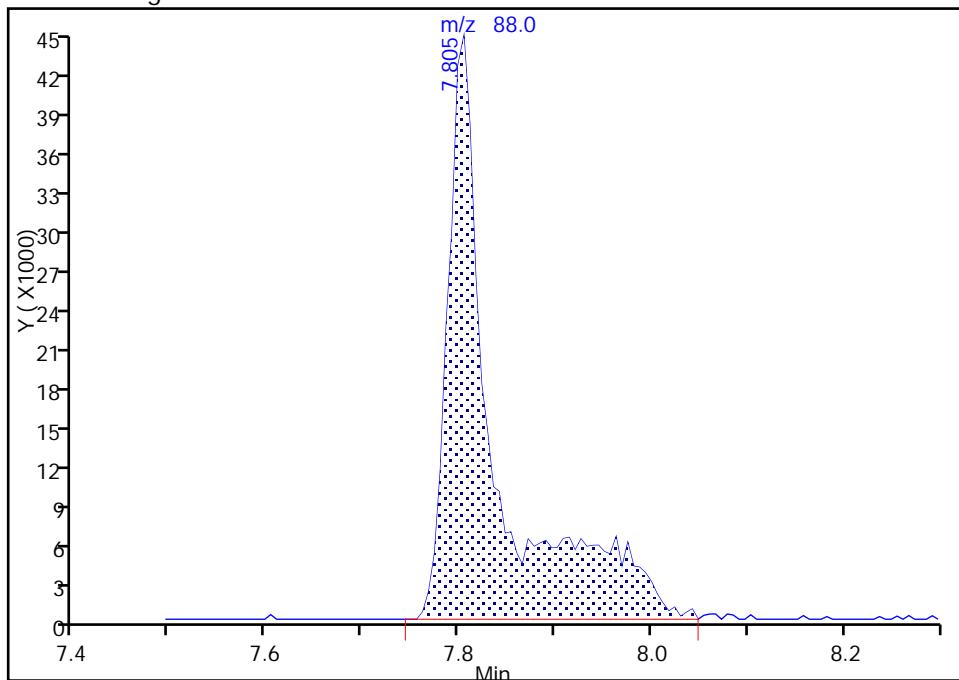
RT: 7.96
 Area: 45805
 Amount: 1076.1525
 Amount Units: ng

Processing Integration Results



RT: 7.81
 Area: 152489
 Amount: 3335.3454
 Amount Units: ng

Manual Integration Results



Reviewer: bungardf, 24-Jul-2017 09:26:43

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D10.D
 Lims ID: IC VSTD50
 Client ID:
 Sample Type: IC Calib Level: 8
 Inject. Date: 24-Jul-2017 09:28:30 ALS Bottle#: 10 Worklist Smp#: 10
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017705-010
 Misc. Info.: IC VSTD50
 Operator ID: 034635 Instrument ID: CHHP6
 Sublist: chrom-MSVOA_LL_CHHP6*sub10
 Method: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 25-Jul-2017 01:44:43 Calib Date: 24-Jul-2017 09:28:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D10.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK021

First Level Reviewer: bungardf Date: 24-Jul-2017 09:58:03

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 3.992 | 3.972 | 0.020 | 94 | 256331 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.052 | 7.051 | 0.001 | 99 | 859285 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.172 | 10.172 | 0.000 | 86 | 220905 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.515 | 12.514 | 0.001 | 96 | 278640 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr) | 113 | 6.316 | 6.315 | 0.001 | 94 | 1076618 | 250.0 | 241.1 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur) | 65 | 6.693 | 6.692 | 0.001 | 68 | 1435595 | 250.0 | 224.9 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.718 | 8.718 | 0.000 | 93 | 3487645 | 250.0 | 205.7 | |
| \$ 8 4-Bromofluorobenzene (Surr) | 95 | 11.353 | 11.352 | 0.001 | 84 | 1557524 | 250.0 | 204.2 | |
| 11 Dichlorodifluoromethane | 85 | 1.485 | 1.484 | 0.001 | 99 | 1341098 | 250.0 | 244.6 | |
| 12 Chloromethane | 50 | 1.637 | 1.630 | 0.007 | 98 | 1170186 | 250.0 | 237.3 | |
| 13 Vinyl chloride | 62 | 1.765 | 1.758 | 0.007 | 97 | 1244722 | 250.0 | 240.8 | |
| 14 Butadiene | 39 | 1.796 | 1.788 | 0.008 | 88 | 1011679 | 250.0 | 236.1 | |
| 15 Bromomethane | 94 | 2.081 | 2.087 | -0.006 | 90 | 482936 | 250.0 | 200.5 | |
| 16 Chloroethane | 64 | 2.209 | 2.202 | 0.007 | 99 | 595090 | 250.0 | 213.1 | |
| 17 Dichlorofluoromethane | 67 | 2.471 | 2.470 | 0.001 | 96 | 1369610 | 250.0 | 227.5 | |
| 18 Trichlorofluoromethane | 101 | 2.495 | 2.506 | -0.011 | 99 | 1263842 | 250.0 | 248.3 | |
| 20 Ethyl ether | 59 | 2.830 | 2.823 | 0.007 | 86 | 1004152 | 250.0 | 230.5 | |
| 21 Acrolein | 56 | 2.994 | 2.999 | -0.005 | 100 | 255428 | 275.0 | 273.8 | |
| 22 1,1-Dichloroethene | 96 | 3.116 | 3.115 | 0.001 | 97 | 1153420 | 250.0 | 258.3 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.158 | 3.157 | 0.001 | 95 | 1098540 | 250.0 | 256.7 | |
| 24 Acetone | 43 | 3.189 | 3.194 | -0.005 | 100 | 730103 | 500.0 | 396.1 | |
| 25 Iodomethane | 142 | 3.292 | 3.291 | 0.001 | 98 | 1578693 | 250.0 | 251.4 | |
| 26 Carbon disulfide | 76 | 3.371 | 3.364 | 0.007 | 99 | 2838844 | 250.0 | 285.5 | |
| 29 3-Chloro-1-propene | 76 | 3.633 | 3.638 | -0.005 | 91 | 717090 | 250.0 | 274.2 | |
| 30 Methyl acetate | 43 | 3.657 | 3.662 | -0.005 | 96 | 1923255 | 500.0 | 476.0 | |
| 31 Methylene Chloride | 84 | 3.846 | 3.839 | 0.007 | 87 | 1372983 | 250.0 | 232.0 | |
| 32 2-Methyl-2-propanol | 59 | 4.113 | 4.112 | 0.001 | 94 | 673044 | 2500.0 | 2372.2 | |
| 33 Acrylonitrile | 53 | 4.241 | 4.240 | 0.001 | 96 | 4723140 | 2500.0 | 2188.1 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.278 | 4.277 | 0.001 | 98 | 1277192 | 250.0 | 252.1 | |
| 35 Methyl tert-butyl ether | 73 | 4.290 | 4.295 | -0.005 | 97 | 3809189 | 250.0 | 235.1 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 36 Hexane | 57 | 4.703 | 4.709 | -0.006 | 90 | 1517069 | 250.0 | 252.8 | |
| 37 1,1-Dichloroethane | 63 | 4.928 | 4.928 | 0.000 | 95 | 2087088 | 250.0 | 253.2 | |
| 38 Vinyl acetate | 43 | 4.983 | 4.982 | 0.001 | 97 | 2492452 | 250.0 | 259.6 | |
| 42 2,2-Dichloropropane | 97 | 5.683 | 5.682 | 0.001 | 89 | 228817 | 250.0 | 268.0 | |
| 43 cis-1,2-Dichloroethene | 96 | 5.695 | 5.694 | 0.001 | 78 | 1462208 | 250.0 | 247.0 | |
| 44 2-Butanone (MEK) | 43 | 5.707 | 5.706 | 0.001 | 97 | 1217733 | 500.0 | 460.8 | |
| 48 Chlorobromomethane | 128 | 5.981 | 5.986 | -0.005 | 96 | 642539 | 250.0 | 248.8 | |
| 49 Tetrahydrofuran | 42 | 5.993 | 5.992 | 0.001 | 84 | 812537 | 500.0 | 454.2 | |
| 50 Chloroform | 83 | 6.139 | 6.138 | 0.001 | 92 | 2152497 | 250.0 | 242.1 | |
| 51 1,1,1-Trichloroethane | 97 | 6.291 | 6.290 | 0.001 | 97 | 1477890 | 250.0 | 261.6 | |
| 52 Cyclohexane | 56 | 6.364 | 6.363 | 0.001 | 87 | 1969875 | 250.0 | 245.4 | |
| 53 Carbon tetrachloride | 117 | 6.468 | 6.467 | 0.001 | 96 | 1172757 | 250.0 | 288.6 | |
| 54 1,1-Dichloropropene | 75 | 6.480 | 6.485 | -0.005 | 96 | 1715254 | 250.0 | 252.0 | |
| 55 Isobutyl alcohol | 41 | 6.705 | 6.698 | 0.007 | 81 | 789622 | 6250.0 | 6624.7 | |
| 56 Benzene | 78 | 6.699 | 6.704 | -0.005 | 97 | 4371437 | 250.0 | 220.7 | |
| 57 1,2-Dichloroethane | 62 | 6.778 | 6.783 | -0.005 | 96 | 1719102 | 250.0 | 234.9 | |
| 59 n-Heptane | 43 | 7.076 | 7.075 | 0.001 | 87 | 1149322 | 250.0 | 251.6 | |
| 61 Trichloroethene | 130 | 7.447 | 7.446 | 0.001 | 98 | 1194165 | 250.0 | 248.2 | |
| 63 Methylcyclohexane | 83 | 7.678 | 7.677 | 0.001 | 90 | 2038808 | 250.0 | 242.6 | |
| 64 1,2-Dichloropropane | 63 | 7.721 | 7.720 | 0.001 | 94 | 1251517 | 250.0 | 252.8 | |
| 65 1,4-Dioxane | 88 | 7.800 | 7.805 | -0.005 | 37 | 203123 | 5000.0 | 4498.8 | |
| 67 Dibromomethane | 93 | 7.806 | 7.811 | -0.005 | 97 | 802657 | 250.0 | 247.9 | |
| 68 Dichlorobromomethane | 83 | 8.007 | 8.006 | 0.001 | 99 | 1522204 | 250.0 | 276.1 | |
| 70 2-Chloroethyl vinyl ether | 63 | 8.317 | 8.316 | 0.001 | 93 | 1665721 | 500.0 | 507.7 | |
| 71 cis-1,3-Dichloropropene | 75 | 8.457 | 8.456 | 0.001 | 96 | 1779441 | 250.0 | 284.5 | |
| 72 4-Methyl-2-pentanone (MIBK) | 43 | 8.615 | 8.614 | 0.001 | 90 | 2362456 | 500.0 | 418.8 | |
| 73 Toluene | 91 | 8.785 | 8.784 | 0.001 | 94 | 4234419 | 250.0 | 194.5 | |
| 74 trans-1,3-Dichloropropene | 75 | 9.035 | 9.040 | -0.005 | 92 | 1562404 | 250.0 | 268.0 | |
| 75 Ethyl methacrylate | 69 | 9.102 | 9.101 | 0.001 | 88 | 1727540 | 250.0 | 235.2 | |
| 76 1,1,2-Trichloroethane | 97 | 9.230 | 9.229 | 0.001 | 92 | 1145832 | 250.0 | 227.5 | |
| 77 Tetrachloroethene | 164 | 9.296 | 9.295 | 0.001 | 95 | 881532 | 250.0 | 226.2 | |
| 78 1,3-Dichloropropane | 76 | 9.388 | 9.387 | 0.001 | 89 | 1979514 | 250.0 | 216.7 | |
| 79 2-Hexanone | 43 | 9.449 | 9.454 | -0.006 | 91 | 1593297 | 500.0 | 438.8 | |
| 81 Chlorodibromomethane | 129 | 9.601 | 9.600 | 0.001 | 91 | 931753 | 250.0 | 273.0 | |
| 82 Ethylene Dibromide | 107 | 9.710 | 9.709 | 0.001 | 98 | 1134020 | 250.0 | 234.0 | |
| 83 3-Chlorobenzotrifluoride | 180 | 10.185 | 10.184 | 0.001 | 91 | 1432766 | 250.0 | 226.4 | |
| 84 Chlorobenzene | 112 | 10.197 | 10.202 | -0.005 | 90 | 2898680 | 250.0 | 204.8 | |
| 85 4-Chlorobenzotrifluoride | 180 | 10.270 | 10.269 | 0.001 | 96 | 1366587 | 250.0 | 233.0 | |
| 86 1,1,1,2-Tetrachloroethane | 131 | 10.294 | 10.293 | 0.001 | 92 | 1029362 | 250.0 | 255.7 | |
| 87 Ethylbenzene | 106 | 10.300 | 10.299 | 0.001 | 95 | 1762716 | 250.0 | 217.7 | |
| 88 m-Xylene & p-Xylene | 106 | 10.434 | 10.433 | 0.001 | 95 | 2161557 | 250.0 | 217.8 | |
| 89 o-Xylene | 106 | 10.817 | 10.816 | 0.001 | 93 | 2111655 | 250.0 | 212.3 | |
| 90 Styrene | 104 | 10.836 | 10.835 | 0.001 | 90 | 3281557 | 250.0 | 204.0 | |
| 91 Bromoform | 173 | 11.012 | 11.017 | -0.005 | 94 | 548248 | 250.0 | 285.5 | |
| 92 2-Chlorobenzotrifluoride | 180 | 11.091 | 11.090 | 0.001 | 92 | 1439864 | 250.0 | 224.5 | |
| 93 Isopropylbenzene | 105 | 11.182 | 11.181 | 0.001 | 97 | 4143279 | 250.0 | 186.1 | |
| 95 Bromobenzene | 156 | 11.493 | 11.492 | 0.001 | 96 | 1259105 | 250.0 | 243.1 | |
| 96 1,1,2,2-Tetrachloroethane | 83 | 11.499 | 11.498 | 0.001 | 96 | 1511708 | 250.0 | 213.3 | |
| 97 trans-1,4-Dichloro-2-butene | 53 | 11.535 | 11.534 | 0.001 | 83 | 404523 | 250.0 | 266.9 | |
| 98 1,2,3-Trichloropropane | 110 | 11.553 | 11.552 | 0.001 | 86 | 529322 | 250.0 | 246.1 | |
| 99 N-Propylbenzene | 120 | 11.602 | 11.601 | 0.001 | 94 | 1357847 | 250.0 | 240.9 | |
| 100 2-Chlorotoluene | 126 | 11.681 | 11.680 | 0.001 | 96 | 1195161 | 250.0 | 244.8 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 101 3-Chlorotoluene | 126 | 11.748 | 11.747 | 0.001 | 93 | 1254125 | 250.0 | 244.8 | |
| 102 1,3,5-Trimethylbenzene | 105 | 11.785 | 11.784 | 0.001 | 95 | 3460126 | 250.0 | 216.2 | |
| 103 4-Chlorotoluene | 126 | 11.809 | 11.808 | 0.001 | 98 | 1276343 | 250.0 | 239.5 | |
| 104 tert-Butylbenzene | 119 | 12.095 | 12.094 | 0.001 | 90 | 2853353 | 250.0 | 224.6 | |
| 106 1,2,4-Trimethylbenzene | 105 | 12.156 | 12.155 | 0.001 | 96 | 3543615 | 250.0 | 212.4 | |
| 107 1,2-dichloro-4-(trifluoromethyl) | 214 | 12.204 | 12.203 | 0.001 | 95 | 945109 | 250.0 | 255.3 | |
| 108 sec-Butylbenzene | 105 | 12.320 | 12.319 | 0.001 | 95 | 3747062 | 250.0 | 210.2 | |
| 109 1,3-Dichlorobenzene | 146 | 12.435 | 12.435 | 0.000 | 93 | 2104721 | 250.0 | 226.3 | |
| 110 4-Isopropyltoluene | 119 | 12.478 | 12.477 | 0.001 | 92 | 3181497 | 250.0 | 216.2 | |
| 111 1,4-Dichlorobenzene | 146 | 12.539 | 12.538 | 0.001 | 90 | 2173124 | 250.0 | 225.9 | |
| 113 2,4-Dichloro-1-(trifluoromethyl) | 214 | 12.569 | 12.568 | 0.001 | 94 | 924522 | 250.0 | 265.0 | |
| 114 2,5-Dichlorobenzotrifluoride | 214 | 12.612 | 12.611 | 0.001 | 96 | 942783 | 250.0 | 244.8 | |
| 116 n-Butylbenzene | 91 | 12.886 | 12.885 | 0.001 | 93 | 2958420 | 250.0 | 220.4 | |
| 117 1,2-Dichlorobenzene | 146 | 12.892 | 12.891 | 0.001 | 93 | 2026312 | 250.0 | 229.1 | |
| 118 1,2-Dibromo-3-Chloropropan | 75 | 13.683 | 13.682 | 0.001 | 82 | 232535 | 250.0 | 294.9 | |
| 119 2,4- & 2,5- & 2,6- Dichlorobenzene | 125 | 13.823 | 13.822 | 0.000 | 94 | 4027755 | 750.0 | 714.9 | |
| 121 2,3- & 3,4- Dichlorotoluene | 125 | 14.236 | 14.235 | 0.001 | 96 | 3139946 | 500.0 | 510.2 | |
| 122 1,2,4-Trichlorobenzene | 180 | 14.498 | 14.503 | -0.005 | 95 | 1247374 | 250.0 | 269.9 | |
| 123 Hexachlorobutadiene | 225 | 14.650 | 14.649 | 0.001 | 95 | 411971 | 250.0 | 281.5 | |
| 124 Naphthalene | 128 | 14.759 | 14.765 | -0.006 | 99 | 3337709 | 250.0 | 236.9 | |
| 125 1,2,3-Trichlorobenzene | 180 | 14.984 | 14.984 | 0.000 | 94 | 1193234 | 250.0 | 287.5 | |
| 126 2,4,5-Trichlorotoluene | 159 | 15.781 | 15.780 | 0.001 | 0 | 796492 | 250.0 | 333.7 | |
| 127 2,3,6-Trichlorotoluene | 159 | 15.885 | 15.884 | 0.001 | 96 | 697018 | 250.0 | 325.8 | |
| 146 3,4-Dichlorotoluene | 1 | 0.000 | | | | | ND | ND | |
| S 131 Xylenes, Total | 106 | | | | 0 | | 500.0 | 430.1 | |
| S 130 1,2-Dichloroethene, Total | 96 | | | | 0 | | 500.0 | 499.1 | |
| S 132 1,3-Dichloropropene, Total | 1 | | | | 0 | | 500.0 | 552.5 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Reagents:

| | | |
|---------------------|---------------------|-----------|
| voaWEEmix1stR_00009 | Amount Added: 10.00 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 10.00 | Units: uL |
| voaWVA1stRest_00016 | Amount Added: 10.00 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 11.00 | Units: uL |
| VOA8260VOAPRI_00263 | Amount Added: 10.00 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 10.00 | Units: uL |
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 10.00 | Units: uL |

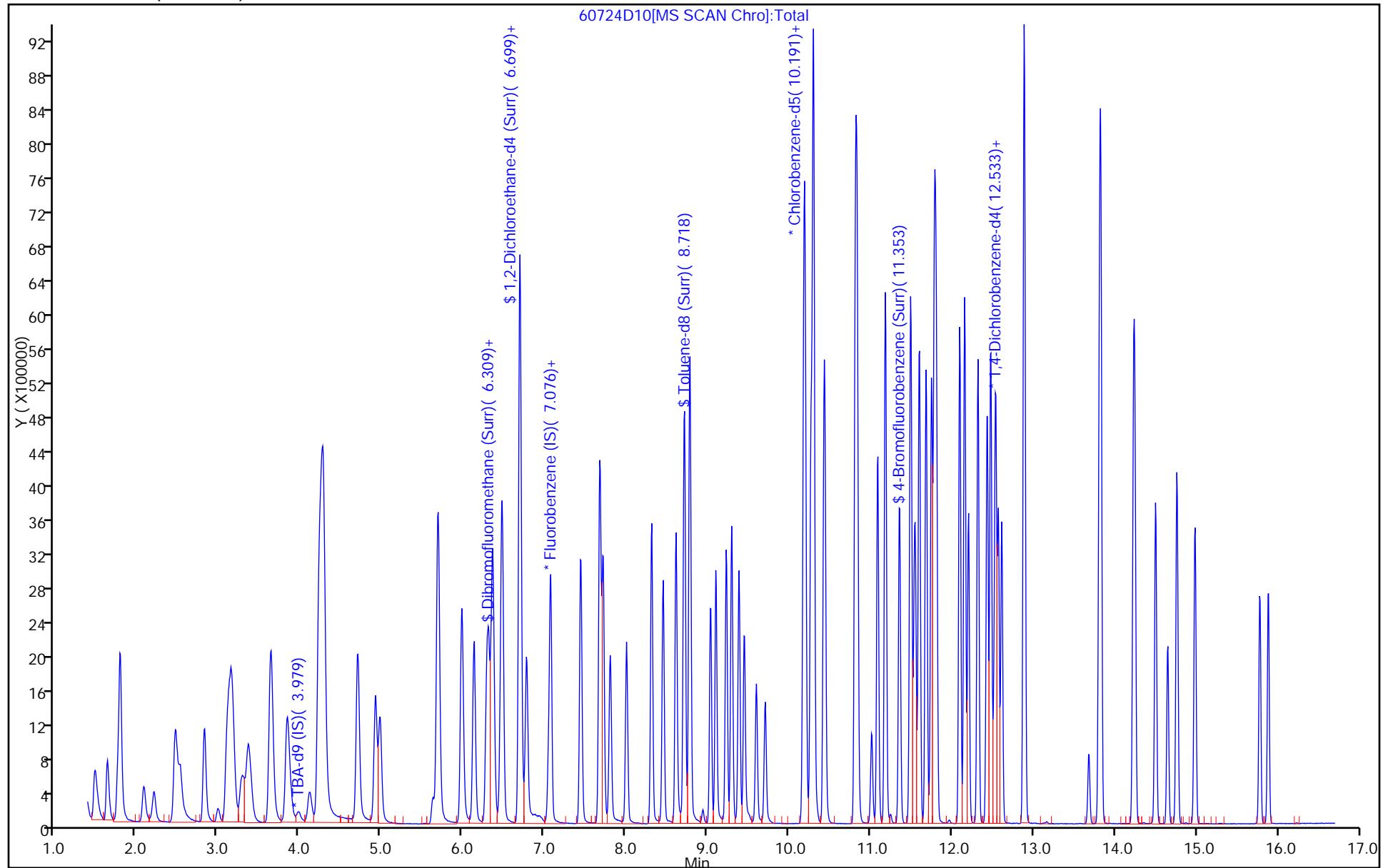
Report Date: 25-Jul-2017 01:44:44

Chrom Revision: 2.2 20-Jun-2017 07:42:38

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D10.D
Injection Date: 24-Jul-2017 09:28:30 Instrument ID: CHHP6
Lims ID: IC VSTD50 Operator ID: 034635
Client ID:
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 10
Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm)

Worklist Smp#: 10



FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

Lab Sample ID: ICV 180-217861/13

Calibration Date: 07/24/2017 10:40

Instrument ID: CHHP6

Calib Start Date: 07/24/2017 06:39

GC Column: DB-624 ID: 0.18 (mm)

Calib End Date: 07/24/2017 09:28

Lab File ID: 60724D13.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 | 0.3191 | 0.2879 | 0.1000 | 9.02 | 10.0 | -9.8 | 30.0 |
| Chloromethane | Ave | 0.2869 | 0.2661 | 0.1000 | 9.28 | 10.0 | -7.2 | 30.0 |
| Vinyl chloride | Ave | 0.3008 | 0.2776 | 0.1000 | 9.23 | 10.0 | -7.7 | 30.0 |
| 1,3-Butadiene | Ave | 0.2494 | 0.2286 | 0.0100 | 9.17 | 10.0 | -8.3 | 30.0 |
| Bromomethane | Ave | 0.1402 | 0.1617 | 0.0500 | 11.5 | 10.0 | 15.4 | 30.0 |
| Chloroethane | Ave | 0.1625 | 0.1750 | 0.0500 | 10.8 | 10.0 | 7.7 | 30.0 |
| Trichlorodifluoromethane | Ave | 0.2961 | 0.2915 | 0.1000 | 9.84 | 10.0 | -1.6 | 30.0 |
| Ethyl ether | Ave | 0.2534 | 0.2447 | 0.0100 | 9.65 | 10.0 | -3.5 | 30.0 |
| Acrolein | Ave | 0.0543 | 0.0504 | 0.0100 | 27.9 | 30.0 | -7.1 | 30.0 |
| 1,1-Dichloroethene | Ave | 0.2599 | 0.2344 | 0.1000 | 9.02 | 10.0 | -9.8 | 30.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.2490 | 0.2264 | 0.1000 | 9.09 | 10.0 | -9.1 | 30.0 |
| Acetone | Ave | 0.1073 | 0.0986 | 0.0500 | 18.4 | 20.0 | -8.1 | 30.0 |
| Iodomethane | Ave | 0.3654 | 0.3466 | 0.0100 | 9.49 | 10.0 | -5.1 | 30.0 |
| Carbon disulfide | Ave | 0.5787 | 0.5422 | 0.1000 | 9.37 | 10.0 | -6.3 | 30.0 |
| Allyl chloride | Ave | 0.1522 | 0.1372 | 0.0100 | 9.02 | 10.0 | -9.8 | 30.0 |
| Methyl acetate | Ave | 0.2351 | 0.2215 | 0.1000 | 18.8 | 20.0 | -5.8 | 30.0 |
| Methylene Chloride | Ave | 0.3444 | 0.3224 | 0.1000 | 9.36 | 10.0 | -6.4 | 30.0 |
| tert-Butyl alcohol | Ave | 1.107 | 1.111 | 0.0100 | 100 | 100 | 0.4 | 30.0 |
| Acrylonitrile | Ave | 0.1256 | 0.1222 | 0.0100 | 97.3 | 100 | -2.7 | 30.0 |
| trans-1,2-Dichloroethene | Ave | 0.2948 | 0.2794 | 0.1000 | 9.48 | 10.0 | -5.2 | 30.0 |
| Methyl tert-butyl ether | Ave | 0.9429 | 0.9239 | 0.1000 | 9.80 | 10.0 | -2.0 | 30.0 |
| Hexane | Ave | 0.3492 | 0.3102 | 0.0100 | 8.88 | 10.0 | -11.2 | 30.0 |
| 1,1-Dichloroethane | Ave | 0.4797 | 0.4818 | 0.2000 | 10.0 | 10.0 | 0.4 | 30.0 |
| Vinyl acetate | Ave | 0.5586 | 0.5200 | 0.0100 | 9.31 | 10.0 | -6.9 | 30.0 |
| 2,2-Dichloropropane | Ave | 0.0497 | 0.0477 | 0.0100 | 9.60 | 10.0 | -4.0 | 30.0 |
| cis-1,2-Dichloroethene | Ave | 0.3444 | 0.3334 | 0.1000 | 9.68 | 10.0 | -3.2 | 30.0 |
| 2-Butanone (MEK) | Ave | 0.1538 | 0.1428 | 0.0500 | 18.6 | 20.0 | -7.1 | 30.0 |
| Bromochloromethane | Ave | 0.1503 | 0.1414 | 0.0100 | 9.41 | 10.0 | -5.9 | 30.0 |
| Tetrahydrofuran | Ave | 0.1041 | 0.0920 | 0.0100 | 17.7 | 20.0 | -11.7 | 30.0 |
| Chloroform | Ave | 0.5173 | 0.5179 | 0.2000 | 10.0 | 10.0 | 0.1 | 30.0 |
| 1,1,1-Trichloroethane | Ave | 0.3287 | 0.3060 | 0.1000 | 9.31 | 10.0 | -6.9 | 30.0 |
| Cyclohexane | Ave | 0.4671 | 0.4280 | 0.1000 | 9.16 | 10.0 | -8.4 | 30.0 |
| Carbon tetrachloride | Ave | 0.2365 | 0.2139 | 0.1000 | 9.04 | 10.0 | -9.6 | 30.0 |
| 1,1-Dichloropropene | Ave | 0.3960 | 0.3728 | 0.0100 | 9.41 | 10.0 | -5.9 | 30.0 |
| Isobutyl alcohol | Ave | 0.0069 | 0.0066* | 0.0100 | 237 | 250 | -5.1 | 30.0 |
| Benzene | Ave | 1.152 | 1.142 | 0.5000 | 9.91 | 10.0 | -0.9 | 30.0 |
| 1,2-Dichloroethane | Ave | 0.4258 | 0.4091 | 0.1000 | 9.61 | 10.0 | -3.9 | 30.0 |
| n-Heptane | Ave | 0.2658 | 0.2342 | 0.0100 | 8.81 | 10.0 | -11.9 | 30.0 |
| Trichloroethene | Ave | 0.2800 | 0.2687 | 0.2000 | 9.60 | 10.0 | -4.0 | 30.0 |
| Methylcyclohexane | Ave | 0.4889 | 0.4513 | 0.1000 | 9.23 | 10.0 | -7.7 | 30.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

Lab Sample ID: ICV 180-217861/13

Calibration Date: 07/24/2017 10:40

Instrument ID: CHHP6

Calib Start Date: 07/24/2017 06:39

GC Column: DB-624 ID: 0.18 (mm)

Calib End Date: 07/24/2017 09:28

Lab File ID: 60724D13.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 | 0.2881 | 0.2658 | 0.1000 | 9.22 | 10.0 | -7.8 | 30.0 |
| 1,4-Dioxane | Ave | 0.0026 | 0.0023* | 0.0100 | 178 | 200 | -11.1 | 30.0 |
| Dibromomethane | Ave | 0.1884 | 0.1798 | 0.0100 | 9.54 | 10.0 | -4.6 | 30.0 |
| Bromodichloromethane | Ave | 0.3208 | 0.3081 | 0.2000 | 9.60 | 10.0 | -4.0 | 30.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.1909 | 0.1830 | 0.0100 | 19.2 | 20.0 | -4.1 | 30.0 |
| cis-1,3-Dichloropropene | Ave | 0.3639 | 0.3408 | 0.2000 | 9.36 | 10.0 | -6.4 | 30.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 1.277 | 1.191 | 0.1000 | 18.7 | 20.0 | -6.7 | 30.0 |
| Toluene | Ave | 4.927 | 4.579 | 0.4000 | 9.29 | 10.0 | -7.1 | 30.0 |
| trans-1,3-Dichloropropene | Ave | 1.320 | 1.145 | 0.1000 | 8.68 | 10.0 | -13.2 | 30.0 |
| Ethyl methacrylate | Ave | 1.663 | 1.503 | 0.0100 | 9.04 | 10.0 | -9.6 | 30.0 |
| 1,1,2-Trichloroethane | Ave | 1.140 | 1.029 | 0.1000 | 9.03 | 10.0 | -9.7 | 30.0 |
| Tetrachloroethene | Ave | 0.8822 | 0.7407 | 0.2000 | 8.40 | 10.0 | -16.0 | 30.0 |
| 1,3-Dichloropropane | Ave | 2.067 | 1.876 | 0.0100 | 9.07 | 10.0 | -9.3 | 30.0 |
| 2-Hexanone | Ave | 0.8218 | 0.7214 | 0.1000 | 17.6 | 20.0 | -12.2 | 30.0 |
| Dibromochloromethane | Ave | 0.7724 | 0.6533 | 0.1000 | 8.46 | 10.0 | -15.4 | 30.0 |
| 1,2-Dibromoethane (EDB) | Ave | 1.097 | 0.9910 | 0.1000 | 9.04 | 10.0 | -9.6 | 30.0 |
| 3-Chlorobenzotrifluoride | Ave | 1.432 | 1.244 | 0.0100 | 8.69 | 10.0 | -13.1 | 30.0 |
| Chlorobenzene | Ave | 3.204 | 2.875 | 0.5000 | 8.97 | 10.0 | -10.3 | 30.0 |
| 4-Chlorobenzotrifluoride | Ave | 1.328 | 1.161 | 0.0100 | 8.75 | 10.0 | -12.5 | 30.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.9111 | 0.7896 | 0.0100 | 8.67 | 10.0 | -13.3 | 30.0 |
| Ethylbenzene | Ave | 1.833 | 1.676 | 0.1000 | 9.14 | 10.0 | -8.6 | 30.0 |
| m-Xylene & p-Xylene | Ave | 2.246 | 2.011 | 0.1000 | 8.95 | 10.0 | -10.5 | 30.0 |
| o-Xylene | Ave | 2.251 | 2.090 | 0.3000 | 9.28 | 10.0 | -7.2 | 30.0 |
| Styrene | Ave | 3.641 | 3.382 | 0.3000 | 9.29 | 10.0 | -7.1 | 30.0 |
| Bromoform | Ave | 0.4347 | 0.3595 | 0.1000 | 8.27 | 10.0 | -17.3 | 30.0 |
| 2-Chlorobenzotrifluoride | Ave | 1.452 | 1.320 | 0.0100 | 9.09 | 10.0 | -9.1 | 30.0 |
| Isopropylbenzene | Ave | 5.039 | 4.803 | 0.1000 | 9.53 | 10.0 | -4.7 | 30.0 |
| Bromobenzene | Ave | 0.9293 | 0.7503 | 0.0100 | 8.07 | 10.0 | -19.3 | 30.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 1.604 | 1.508 | 0.3000 | 9.40 | 10.0 | -6.0 | 30.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2720 | 0.2040 | 0.0100 | 7.50 | 10.0 | -25.0 | 30.0 |
| 1,2,3-Trichloropropane | Ave | 0.3860 | 0.3026 | 0.0100 | 7.84 | 10.0 | -21.6 | 30.0 |
| N-Propylbenzene | Ave | 1.011 | 0.7830 | 0.0100 | 7.74 | 10.0 | -22.6 | 30.0 |
| 2-Chlorotoluene | Ave | 0.8762 | 0.6856 | 0.0100 | 7.82 | 10.0 | -21.8 | 30.0 |
| 3-Chlorotoluene | Ave | 0.9194 | 0.7153 | 0.0100 | 7.78 | 10.0 | -22.2 | 30.0 |
| 1,3,5-Trimethylbenzene | Ave | 2.872 | 2.404 | 0.0100 | 8.37 | 10.0 | -16.3 | 30.0 |
| 4-Chlorotoluene | Ave | 0.9565 | 0.7290 | 0.0100 | 7.62 | 10.0 | -23.8 | 30.0 |
| tert-Butylbenzene | Ave | 2.280 | 1.881 | 0.0100 | 8.25 | 10.0 | -17.5 | 30.0 |
| 1,2,4-Trimethylbenzene | Ave | 2.994 | 2.513 | 0.0100 | 8.39 | 10.0 | -16.1 | 30.0 |
| 3,4-Dichlorobenzotrifluoride | Ave | 0.6644 | 0.5039 | 0.0100 | 7.58 | 10.0 | -24.2 | 30.0 |
| sec-Butylbenzene | Ave | 3.198 | 2.650 | 0.0100 | 8.29 | 10.0 | -17.1 | 30.0 |
| 1,3-Dichlorobenzene | Ave | 1.669 | 1.347 | 0.6000 | 8.07 | 10.0 | -19.3 | 30.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.: _____

Lab Sample ID: ICV 180-217861/13 Calibration Date: 07/24/2017 10:40

Instrument ID: CHHP6 Calib Start Date: 07/24/2017 06:39

GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 07/24/2017 09:28

Lab File ID: 60724D13.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 |
|---------------------------------------|------------|---------|--------|---------|-------------|--------------|-------|--------|
| 4-Isopropyltoluene | Ave | 2.641 | 2.215 | 0.0100 | 8.39 | 10.0 | -16.1 | 30.0 |
| 1,4-Dichlorobenzene | Ave | 1.726 | 1.396 | 0.5000 | 8.09 | 10.0 | -19.1 | 30.0 |
| 2,4-Dichlorobenzotrifluoride | Ave | 0.6261 | 0.4701 | 0.0100 | 7.51 | 10.0 | -24.9 | 30.0 |
| 2,5-Dichlorobenzotrifluoride | Ave | 0.6911 | 0.5511 | 0.0100 | 7.97 | 10.0 | -20.3 | 30.0 |
| n-Butylbenzene | Ave | 2.409 | 1.985 | 0.0100 | 8.24 | 10.0 | -17.6 | 30.0 |
| 1,2-Dichlorobenzene | Ave | 1.587 | 1.343 | 0.4000 | 8.46 | 10.0 | -15.4 | 30.0 |
| 1,2-Dibromo-3-Chloropropane | Ave | 0.1415 | 0.1049 | 0.0500 | 7.42 | 10.0 | -25.8 | 30.0 |
| 2,4- & 2,5- & 2,6- Dichlorotoluene | Ave | 1.011 | 0.8617 | 0.0100 | 25.6 | 30.0 | -14.8 | 30.0 |
| 2,3- & 3,4- Dichlorotoluene | Ave | 1.104 | 0.9140 | 0.0100 | 16.6 | 20.0 | -17.2 | 30.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.8294 | 0.6700 | 0.2000 | 8.08 | 10.0 | -19.2 | 30.0 |
| Hexachlorobutadiene | Ave | 0.2626 | 0.2044 | 0.0100 | 7.78 | 10.0 | -22.2 | 30.0 |
| Naphthalene | Ave | 2.528 | 2.097 | 0.0100 | 8.30 | 10.0 | -17.0 | 30.0 |
| 1,2,3-Trichlorobenzene | Ave | 0.7447 | 0.5897 | 0.0100 | 7.92 | 10.0 | -20.8 | 30.0 |
| 2,4,5-Trichlorotoluene | Ave | 0.4283 | 0.3059 | 0.0100 | 7.14 | 10.0 | -28.6 | 30.0 |
| 2,3,6-Trichlorotoluene | Ave | 0.3839 | 0.2777 | 0.0100 | 7.23 | 10.0 | -27.7 | 30.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2598 | 0.2349 | | 9.04 | 10.0 | -9.6 | 30.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.3713 | 0.3133 | | 8.44 | 10.0 | -15.6 | 30.0 |
| Toluene-d8 (Surr) | Lin2 | | 3.852 | | 9.58 | 10.0 | -4.2 | 30.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 1.726 | 1.745 | | 10.1 | 10.0 | 1.1 | 30.0 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170724-17705.b\60724D13.D
 Lims ID: ICV
 Client ID:
 Sample Type: ICV
 Inject. Date: 24-Jul-2017 10:40:30 ALS Bottle#: 13 Worklist Smp#: 13
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017705-013
 Misc. Info.: ICV
 Operator ID: 034635 Instrument ID: CHHP6
 Sublist:
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170724-17705.b\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 25-Jul-2017 01:44:46 Calib Date: 24-Jul-2017 09:28:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170724-17705.b\60724D10.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK021

First Level Reviewer: bungardf Date: 25-Jul-2017 01:07:50

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 3.962 | 3.961 | 0.001 | 95 | 317387 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.053 | 7.051 | 0.002 | 98 | 870770 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.167 | 10.172 | -0.005 | 88 | 217522 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.510 | 12.514 | -0.004 | 94 | 360072 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr) | 113 | 6.317 | 6.315 | 0.002 | 92 | 204537 | 50.0 | 45.2 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur) | 65 | 6.694 | 6.692 | 0.002 | 88 | 272824 | 50.0 | 42.2 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.713 | 8.718 | -0.005 | 93 | 837799 | 50.0 | 47.9 | |
| \$ 8 4-Bromofluorobenzene (Surr) | 95 | 11.354 | 11.352 | 0.002 | 86 | 379643 | 50.0 | 50.5 | |
| 11 Dichlorodifluoromethane | 85 | 1.480 | 1.484 | -0.004 | 86 | 250657 | 50.0 | 45.1 | |
| 12 Chloromethane | 50 | 1.632 | 1.630 | 0.002 | 99 | 231750 | 50.0 | 46.4 | |
| 13 Vinyl chloride | 62 | 1.754 | 1.758 | -0.004 | 98 | 241724 | 50.0 | 46.1 | |
| 14 Butadiene | 39 | 1.790 | 1.788 | 0.002 | 88 | 199035 | 50.0 | 45.8 | |
| 15 Bromomethane | 94 | 2.095 | 2.087 | 0.008 | 89 | 140806 | 50.0 | 57.7 | |
| 16 Chloroethane | 64 | 2.210 | 2.202 | 0.008 | 99 | 152345 | 50.0 | 53.8 | |
| 17 Dichlorofluoromethane | 67 | 2.466 | 2.470 | -0.004 | 96 | 316935 | 50.0 | 52.0 | |
| 18 Trichlorofluoromethane | 101 | 2.502 | 2.506 | -0.004 | 98 | 253848 | 50.0 | 49.2 | |
| 20 Ethyl ether | 59 | 2.825 | 2.823 | 0.002 | 88 | 213070 | 50.0 | 48.3 | |
| 21 Acrolein | 56 | 2.995 | 2.999 | -0.004 | 98 | 131728 | 150.0 | 139.3 | |
| 22 1,1-Dichloroethene | 96 | 3.117 | 3.115 | 0.002 | 97 | 204120 | 50.0 | 45.1 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.159 | 3.157 | 0.002 | 94 | 197174 | 50.0 | 45.5 | |
| 24 Acetone | 43 | 3.184 | 3.194 | -0.010 | 96 | 171641 | 100.0 | 91.9 | |
| 25 Iodomethane | 142 | 3.293 | 3.291 | 0.002 | 98 | 301806 | 50.0 | 47.4 | |
| 26 Carbon disulfide | 76 | 3.372 | 3.364 | 0.008 | 99 | 472119 | 50.0 | 46.8 | |
| 29 3-Chloro-1-propene | 76 | 3.634 | 3.638 | -0.004 | 92 | 119495 | 50.0 | 45.1 | |
| 30 Methyl acetate | 43 | 3.658 | 3.662 | -0.004 | 97 | 385795 | 100.0 | 94.2 | |
| 31 Methylene Chloride | 84 | 3.841 | 3.839 | 0.002 | 89 | 280759 | 50.0 | 46.8 | |
| 32 2-Methyl-2-propanol | 59 | 4.096 | 4.112 | -0.016 | 94 | 176303 | 500.0 | 501.8 | |
| 33 Acrylonitrile | 53 | 4.236 | 4.240 | -0.004 | 99 | 1064053 | 500.0 | 486.4 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.279 | 4.277 | 0.002 | 99 | 243284 | 50.0 | 47.4 | |
| 35 Methyl tert-butyl ether | 73 | 4.285 | 4.295 | -0.010 | 95 | 804468 | 50.0 | 49.0 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 36 Hexane | 57 | 4.717 | 4.709 | 0.008 | 92 | 270066 | 50.0 | 44.4 | |
| 37 1,1-Dichloroethane | 63 | 4.930 | 4.928 | 0.002 | 96 | 419512 | 50.0 | 50.2 | |
| 38 Vinyl acetate | 43 | 4.984 | 4.982 | 0.002 | 97 | 452761 | 50.0 | 46.5 | |
| 42 2,2-Dichloropropane | 97 | 5.690 | 5.682 | 0.008 | 53 | 41509 | 50.0 | 48.0 | |
| 43 cis-1,2-Dichloroethene | 96 | 5.690 | 5.694 | -0.004 | 80 | 290340 | 50.0 | 48.4 | |
| 44 2-Butanone (MEK) | 43 | 5.702 | 5.706 | -0.004 | 97 | 248754 | 100.0 | 92.9 | |
| 48 Chlorobromomethane | 128 | 5.982 | 5.986 | -0.004 | 97 | 123123 | 50.0 | 47.0 | |
| 49 Tetrahydrofuran | 42 | 5.994 | 5.992 | 0.002 | 90 | 160129 | 100.0 | 88.3 | |
| 50 Chloroform | 83 | 6.134 | 6.138 | -0.004 | 93 | 450991 | 50.0 | 50.1 | |
| 51 1,1,1-Trichloroethane | 97 | 6.292 | 6.290 | 0.002 | 98 | 266443 | 50.0 | 46.5 | |
| 52 Cyclohexane | 56 | 6.365 | 6.363 | 0.002 | 90 | 372721 | 50.0 | 45.8 | |
| 53 Carbon tetrachloride | 117 | 6.463 | 6.467 | -0.004 | 96 | 186212 | 50.0 | 45.2 | |
| 54 1,1-Dichloropropene | 75 | 6.481 | 6.485 | -0.004 | 98 | 324612 | 50.0 | 47.1 | |
| 55 Isobutyl alcohol | 41 | 6.694 | 6.698 | -0.004 | 71 | 143231 | 1250.0 | 1185.8 | |
| 56 Benzene | 78 | 6.700 | 6.704 | -0.004 | 97 | 994365 | 50.0 | 49.5 | |
| 57 1,2-Dichloroethane | 62 | 6.779 | 6.783 | -0.004 | 98 | 356187 | 50.0 | 48.0 | |
| 59 n-Heptane | 43 | 7.077 | 7.075 | 0.002 | 91 | 203911 | 50.0 | 44.1 | |
| 61 Trichloroethene | 130 | 7.448 | 7.446 | 0.002 | 98 | 234008 | 50.0 | 48.0 | |
| 63 Methylcyclohexane | 83 | 7.679 | 7.677 | 0.002 | 86 | 392960 | 50.0 | 46.2 | |
| 64 1,2-Dichloropropane | 63 | 7.716 | 7.720 | -0.004 | 92 | 231431 | 50.0 | 46.1 | |
| 65 1,4-Dioxane | 88 | 7.801 | 7.805 | -0.004 | 64 | 40694 | 1000.0 | 889.4 | M |
| 67 Dibromomethane | 93 | 7.807 | 7.811 | -0.004 | 95 | 156565 | 50.0 | 47.7 | |
| 68 Dichlorobromomethane | 83 | 8.008 | 8.006 | 0.002 | 99 | 268266 | 50.0 | 48.0 | |
| 70 2-Chloroethyl vinyl ether | 63 | 8.312 | 8.316 | -0.004 | 93 | 318759 | 100.0 | 95.9 | |
| 71 cis-1,3-Dichloropropene | 75 | 8.452 | 8.456 | -0.004 | 96 | 296743 | 50.0 | 46.8 | |
| 72 4-Methyl-2-pentanone (MIBK) | 43 | 8.616 | 8.614 | 0.002 | 94 | 518033 | 100.0 | 93.3 | |
| 73 Toluene | 91 | 8.780 | 8.784 | -0.004 | 99 | 996103 | 50.0 | 46.5 | |
| 74 trans-1,3-Dichloropropene | 75 | 9.036 | 9.040 | -0.004 | 92 | 249017 | 50.0 | 43.4 | |
| 75 Ethyl methacrylate | 69 | 9.103 | 9.101 | 0.002 | 88 | 327011 | 50.0 | 45.2 | |
| 76 1,1,2-Trichloroethane | 97 | 9.231 | 9.229 | 0.002 | 90 | 223842 | 50.0 | 45.1 | |
| 77 Tetrachloroethene | 164 | 9.297 | 9.295 | 0.002 | 97 | 161123 | 50.0 | 42.0 | |
| 78 1,3-Dichloropropane | 76 | 9.389 | 9.387 | 0.002 | 89 | 408048 | 50.0 | 45.4 | |
| 79 2-Hexanone | 43 | 9.450 | 9.454 | -0.004 | 93 | 313824 | 100.0 | 87.8 | |
| 81 Chlorodibromomethane | 129 | 9.602 | 9.600 | 0.002 | 91 | 142099 | 50.0 | 42.3 | |
| 82 Ethylene Dibromide | 107 | 9.711 | 9.709 | 0.002 | 95 | 215564 | 50.0 | 45.2 | |
| 83 3-Chlorobenzotrifluoride | 180 | 10.186 | 10.184 | 0.002 | 91 | 270569 | 50.0 | 43.4 | |
| 84 Chlorobenzene | 112 | 10.198 | 10.202 | -0.004 | 92 | 625436 | 50.0 | 44.9 | |
| 85 4-Chlorobenzotrifluoride | 180 | 10.271 | 10.269 | 0.002 | 96 | 252548 | 50.0 | 43.7 | |
| 86 1,1,1,2-Tetrachloroethane | 131 | 10.295 | 10.293 | 0.002 | 89 | 171758 | 50.0 | 43.3 | |
| 87 Ethylbenzene | 106 | 10.301 | 10.299 | 0.002 | 98 | 364512 | 50.0 | 45.7 | |
| 88 m-Xylene & p-Xylene | 106 | 10.435 | 10.433 | 0.002 | 99 | 437513 | 50.0 | 44.8 | |
| 89 o-Xylene | 106 | 10.812 | 10.816 | -0.004 | 97 | 454567 | 50.0 | 46.4 | |
| 90 Styrene | 104 | 10.837 | 10.835 | 0.002 | 94 | 735565 | 50.0 | 46.4 | |
| 91 Bromoform | 173 | 11.013 | 11.017 | -0.004 | 95 | 78192 | 50.0 | 41.3 | |
| 92 2-Chlorobenzotrifluoride | 180 | 11.092 | 11.090 | 0.002 | 97 | 287032 | 50.0 | 45.4 | |
| 93 Isopropylbenzene | 105 | 11.183 | 11.181 | 0.002 | 97 | 1044738 | 50.0 | 47.7 | |
| 95 Bromobenzene | 156 | 11.494 | 11.492 | 0.002 | 95 | 270177 | 50.0 | 40.4 | |
| 96 1,1,2,2-Tetrachloroethane | 83 | 11.500 | 11.498 | 0.002 | 95 | 328079 | 50.0 | 47.0 | |
| 97 trans-1,4-Dichloro-2-butene | 53 | 11.530 | 11.534 | -0.004 | 69 | 73439 | 50.0 | 37.5 | |
| 98 1,2,3-Trichloropropane | 110 | 11.548 | 11.552 | -0.004 | 85 | 108942 | 50.0 | 39.2 | |
| 99 N-Propylbenzene | 120 | 11.597 | 11.601 | -0.004 | 98 | 281925 | 50.0 | 38.7 | |
| 100 2-Chlorotoluene | 126 | 11.682 | 11.680 | 0.002 | 95 | 246870 | 50.0 | 39.1 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 101 3-Chlorotoluene | 126 | 11.749 | 11.747 | 0.002 | 96 | 257552 | 50.0 | 38.9 | |
| 102 1,3,5-Trimethylbenzene | 105 | 11.786 | 11.784 | 0.002 | 93 | 865530 | 50.0 | 41.8 | |
| 103 4-Chlorotoluene | 126 | 11.804 | 11.808 | -0.004 | 99 | 262485 | 50.0 | 38.1 | |
| 104 tert-Butylbenzene | 119 | 12.096 | 12.094 | 0.002 | 91 | 677153 | 50.0 | 41.2 | |
| 106 1,2,4-Trimethylbenzene | 105 | 12.157 | 12.155 | 0.002 | 98 | 905005 | 50.0 | 42.0 | |
| 107 1,2-dichloro-4-(trifluoromethyl) | 214 | 12.205 | 12.203 | 0.002 | 96 | 181425 | 50.0 | 37.9 | |
| 108 sec-Butylbenzene | 105 | 12.321 | 12.319 | 0.002 | 95 | 954305 | 50.0 | 41.4 | |
| 109 1,3-Dichlorobenzene | 146 | 12.430 | 12.435 | -0.005 | 96 | 484901 | 50.0 | 40.3 | |
| 110 4-Isopropyltoluene | 119 | 12.479 | 12.477 | 0.002 | 96 | 797606 | 50.0 | 41.9 | |
| 111 1,4-Dichlorobenzene | 146 | 12.534 | 12.538 | -0.004 | 94 | 502627 | 50.0 | 40.4 | |
| 113 2,4-Dichloro-1-(trifluoromethyl) | 214 | 12.570 | 12.568 | 0.002 | 96 | 169284 | 50.0 | 37.5 | |
| 114 2,5-Dichlorobenzotrifluoride | 214 | 12.613 | 12.611 | 0.002 | 98 | 198432 | 50.0 | 39.9 | |
| 116 n-Butylbenzene | 91 | 12.881 | 12.885 | -0.004 | 96 | 714807 | 50.0 | 41.2 | |
| 117 1,2-Dichlorobenzene | 146 | 12.893 | 12.891 | 0.002 | 95 | 483712 | 50.0 | 42.3 | |
| 118 1,2-Dibromo-3-Chloropropan | 75 | 13.678 | 13.682 | -0.004 | 74 | 37785 | 50.0 | 37.1 | |
| 119 2,4- & 2,5- & 2,6- Dichlorobenzene | 125 | 13.817 | 13.822 | -0.005 | 99 | 930835 | 150.0 | 127.9 | |
| 121 2,3- & 3,4- Dichlorotoluene | 125 | 14.237 | 14.235 | 0.002 | 99 | 658230 | 100.0 | 82.8 | |
| 122 1,2,4-Trichlorobenzene | 180 | 14.499 | 14.503 | -0.004 | 94 | 241256 | 50.0 | 40.4 | |
| 123 Hexachlorobutadiene | 225 | 14.651 | 14.649 | 0.002 | 97 | 73584 | 50.0 | 38.9 | |
| 124 Naphthalene | 128 | 14.760 | 14.765 | -0.005 | 98 | 755128 | 50.0 | 41.5 | |
| 125 1,2,3-Trichlorobenzene | 180 | 14.986 | 14.984 | 0.002 | 94 | 212348 | 50.0 | 39.6 | |
| 126 2,4,5-Trichlorotoluene | 159 | 15.776 | 15.780 | -0.004 | 0 | 110148 | 50.0 | 35.7 | |
| 127 2,3,6-Trichlorotoluene | 159 | 15.880 | 15.884 | -0.004 | 97 | 99997 | 50.0 | 36.2 | |
| 146 3,4-Dichlorotoluene | 1 | 0.000 | | | | | ND | ND | |
| S 130 1,2-Dichloroethene, Total | 96 | | | | | 0 | 100.0 | 95.8 | |
| S 131 Xylenes, Total | 106 | | | | | 0 | 100.0 | 91.2 | |
| S 132 1,3-Dichloropropene, Total | 1 | | | | | 0 | 100.0 | 90.2 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

Reagents:

| | | | |
|---------------------|--------------------|-----------|-------------|
| VOA8260VOA2ND_00253 | Amount Added: 2.00 | Units: uL | |
| voaW2cleve2nd_00012 | Amount Added: 2.00 | Units: uL | |
| voaWva2ndRete_00003 | Amount Added: 2.00 | Units: uL | |
| voaWAcro2ndRe_00012 | Amount Added: 6.00 | Units: uL | |
| voaWKet2ndRes_00021 | Amount Added: 2.00 | Units: uL | |
| voaWee2ndRest_00014 | Amount Added: 2.00 | Units: uL | |
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL | Run Reagent |
| VOA8260SURR_00071 | Amount Added: 2.00 | Units: uL | Run Reagent |

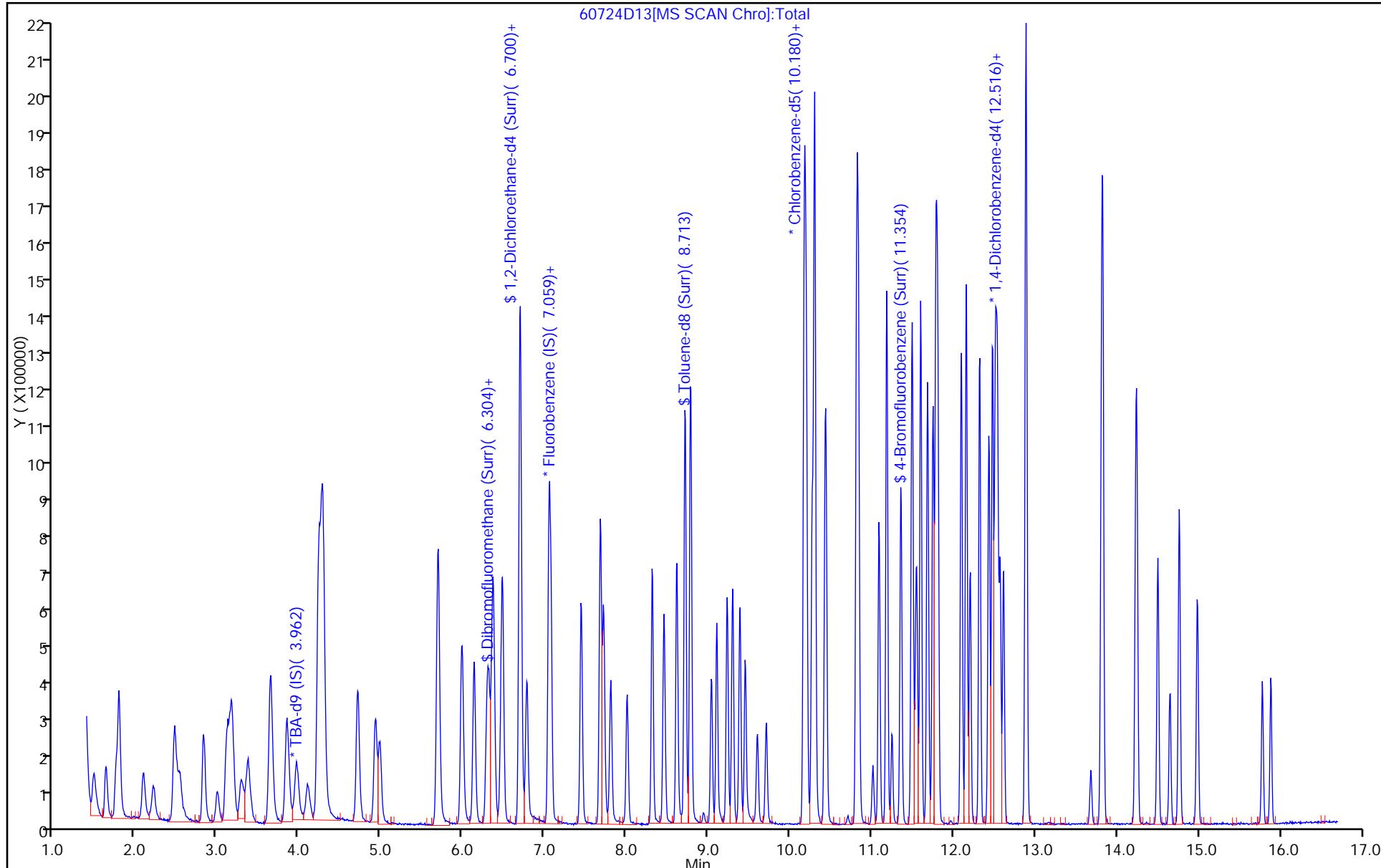
Report Date: 25-Jul-2017 01:44:50

Chrom Revision: 2.2 20-Jun-2017 07:42:38

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D13.D
Injection Date: 24-Jul-2017 10:40:30 Instrument ID: CHHP6
Lims ID: ICV Operator ID: 034635
Client ID:
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 13
Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm)

Worklist Smp#: 13



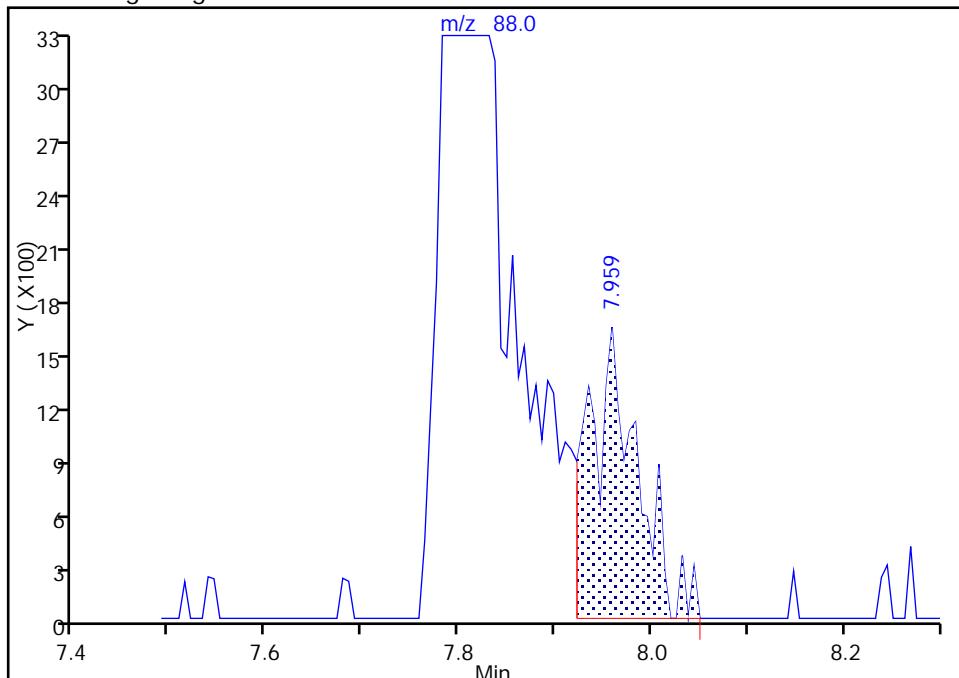
TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D13.D
 Injection Date: 24-Jul-2017 10:40:30 Instrument ID: CHHP6
 Lims ID: ICV
 Client ID:
 Operator ID: 034635 ALS Bottle#: 13 Worklist Smp#: 13
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
 Column: DB-624 (0.18 mm) Detector: MS SCAN

65 1,4-Dioxane, CAS: 123-91-1
Signal: 1

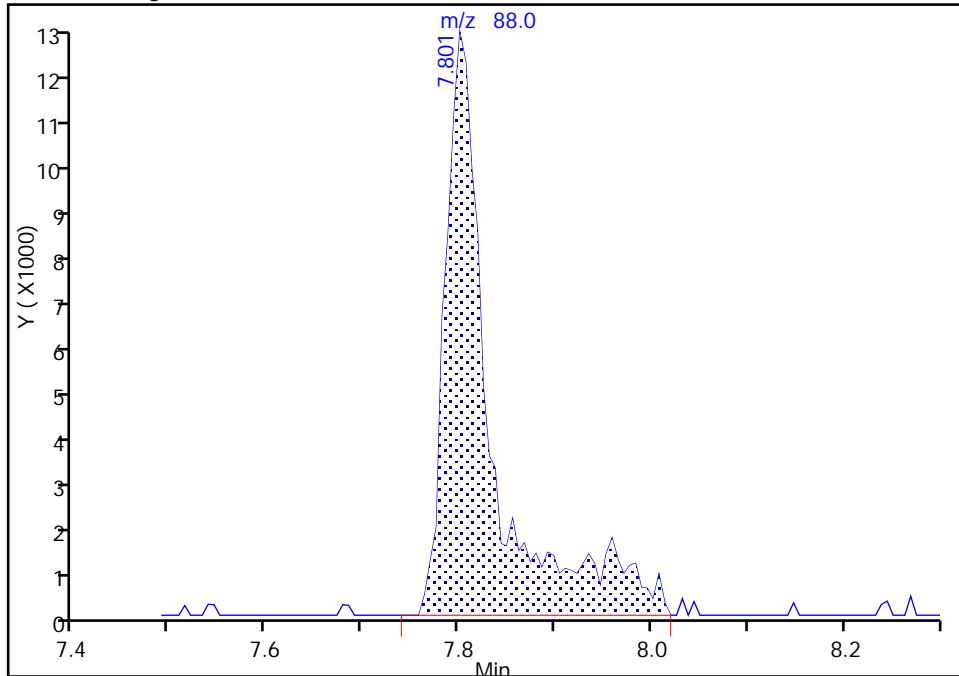
RT: 7.96
 Area: 5522
 Amount: 120.6888
 Amount Units: ng

Processing Integration Results



RT: 7.80
 Area: 40694
 Amount: 889.4077
 Amount Units: ng

Manual Integration Results



Reviewer: bungardf, 25-Jul-2017 01:35:35

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

Lab Sample ID: CCVIS 180-226148/2 Calibration Date: 10/17/2017 23:55

Instrument ID: CHHP6 Calib Start Date: 07/24/2017 06:39

GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 07/24/2017 09:28

Lab File ID: 6101802D.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 | 0.3191 | 0.3616 | 0.1000 | 11.3 | 10.0 | 13.3 | 20.0 |
| Chloromethane | Ave | 0.2869 | 0.2441 | 0.1000 | 8.51 | 10.0 | -14.9 | 20.0 |
| Vinyl chloride | Ave | 0.3008 | 0.3093 | 0.1000 | 10.3 | 10.0 | 2.8 | 20.0 |
| 1,3-Butadiene | Ave | 0.2494 | 0.2428 | 0.0100 | 9.74 | 10.0 | -2.6 | 20.0 |
| Bromomethane | Ave | 0.1402 | 0.1524 | 0.0500 | 10.9 | 10.0 | 8.7 | 20.0 |
| Chloroethane | Ave | 0.1625 | 0.1765 | 0.0500 | 10.9 | 10.0 | 8.6 | 20.0 |
| Trichlorofluoromethane | Ave | 0.2961 | 0.4104 | 0.1000 | 13.9 | 10.0 | 38.6* | 20.0 |
| Ethyl ether | Ave | 0.2534 | 0.2535 | 0.0100 | 10.0 | 10.0 | 0.0 | 20.0 |
| Acrolein | Ave | 0.0543 | 0.0515 | 0.0100 | 28.5 | 30.0 | -5.1 | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2599 | 0.2769 | 0.1000 | 10.7 | 10.0 | 6.6 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.2490 | 0.3032 | 0.1000 | 12.2 | 10.0 | 21.8* | 20.0 |
| Acetone | Ave | 0.1073 | 0.1452 | 0.0500 | 27.1 | 20.0 | 35.4* | 20.0 |
| Iodomethane | Ave | 0.3654 | 0.3809 | 0.0100 | 10.4 | 10.0 | 4.3 | 20.0 |
| Carbon disulfide | Ave | 0.5787 | 0.5757 | 0.1000 | 9.95 | 10.0 | -0.5 | 20.0 |
| Allyl chloride | Ave | 0.1522 | 0.1467 | 0.0100 | 9.64 | 10.0 | -3.6 | 20.0 |
| Methyl acetate | Ave | 0.2351 | 0.1853 | 0.1000 | 15.8 | 20.0 | -21.2* | 20.0 |
| Methylene Chloride | Ave | 0.3444 | 0.3389 | 0.1000 | 9.84 | 10.0 | -1.6 | 20.0 |
| tert-Butyl alcohol | Ave | 1.107 | 1.047 | 0.0100 | 94.6 | 100 | -5.4 | 20.0 |
| Acrylonitrile | Ave | 0.1256 | 0.0947 | 0.0100 | 75.4 | 100 | -24.6* | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2948 | 0.3058 | 0.1000 | 10.4 | 10.0 | 3.7 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.9429 | 0.8659 | 0.1000 | 9.18 | 10.0 | -8.2 | 20.0 |
| Hexane | Ave | 0.3492 | 0.3345 | 0.0100 | 9.58 | 10.0 | -4.2 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.4797 | 0.4881 | 0.2000 | 10.2 | 10.0 | 1.7 | 20.0 |
| Vinyl acetate | Ave | 0.5586 | 0.4171 | 0.0100 | 7.47 | 10.0 | -25.3* | 20.0 |
| 2,2-Dichloropropane | Ave | 0.0497 | 0.0585 | 0.0100 | 11.8 | 10.0 | 17.8 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.3444 | 0.3462 | 0.1000 | 10.1 | 10.0 | 0.5 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.1538 | 0.1621 | 0.0500 | 21.1 | 20.0 | 5.4 | 20.0 |
| Bromochloromethane | Ave | 0.1503 | 0.1547 | 0.0100 | 10.3 | 10.0 | 2.9 | 20.0 |
| Tetrahydrofuran | Ave | 0.1041 | 0.0671 | 0.0100 | 12.9 | 20.0 | -35.6* | 20.0 |
| Chloroform | Ave | 0.5173 | 0.5451 | 0.2000 | 10.5 | 10.0 | 5.4 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.3287 | 0.3880 | 0.1000 | 11.8 | 10.0 | 18.0 | 20.0 |
| Cyclohexane | Ave | 0.4671 | 0.4191 | 0.1000 | 8.97 | 10.0 | -10.3 | 20.0 |
| Carbon tetrachloride | Ave | 0.2365 | 0.3178 | 0.1000 | 13.4 | 10.0 | 34.4* | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3960 | 0.4302 | 0.0100 | 10.9 | 10.0 | 8.6 | 20.0 |
| Benzene | Ave | 1.152 | 1.225 | 0.5000 | 10.6 | 10.0 | 6.3 | 20.0 |
| Isobutyl alcohol | Ave | 0.0069 | 0.0060* | 0.0100 | 216 | 250 | -13.6 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.4258 | 0.4382 | 0.1000 | 10.3 | 10.0 | 2.9 | 20.0 |
| n-Heptane | Ave | 0.2658 | 0.2318 | 0.0100 | 8.72 | 10.0 | -12.8 | 20.0 |
| Trichloroethene | Ave | 0.2800 | 0.2896 | 0.2000 | 10.3 | 10.0 | 3.4 | 20.0 |
| Methylcyclohexane | Ave | 0.4889 | 0.5045 | 0.1000 | 10.3 | 10.0 | 3.2 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

Lab Sample ID: CCVIS 180-226148/2

Calibration Date: 10/17/2017 23:55

Instrument ID: CHHP6

Calib Start Date: 07/24/2017 06:39

GC Column: DB-624 ID: 0.18 (mm)

Calib End Date: 07/24/2017 09:28

Lab File ID: 6101802D.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 | 0.2881 | 0.2638 | 0.1000 | 9.15 | 10.0 | -8.5 | 20.0 |
| 1,4-Dioxane | Ave | 0.0026 | 0.0022* | 0.0100 | 168 | 200 | -16.1 | 20.0 |
| Dibromomethane | Ave | 0.1884 | 0.1906 | 0.0100 | 10.1 | 10.0 | 1.1 | 20.0 |
| Bromodichloromethane | Ave | 0.3208 | 0.3221 | 0.2000 | 10.0 | 10.0 | 0.4 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.1909 | 0.1969 | 0.0100 | 20.6 | 20.0 | 3.1 | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3639 | 0.3672 | 0.2000 | 10.1 | 10.0 | 0.9 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 1.277 | 0.9695 | 0.1000 | 15.2 | 20.0 | -24.1* | 20.0 |
| Toluene | Ave | 4.927 | 4.828 | 0.4000 | 9.80 | 10.0 | -2.0 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 1.320 | 1.292 | 0.1000 | 9.79 | 10.0 | -2.1 | 20.0 |
| Ethyl methacrylate | Ave | 1.663 | 1.498 | 0.0100 | 9.01 | 10.0 | -9.9 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 1.140 | 1.091 | 0.1000 | 9.57 | 10.0 | -4.3 | 20.0 |
| Tetrachloroethene | Ave | 0.8822 | 0.8422 | 0.2000 | 9.55 | 10.0 | -4.5 | 20.0 |
| 1,3-Dichloropropane | Ave | 2.067 | 2.022 | 0.0100 | 9.78 | 10.0 | -2.2 | 20.0 |
| 2-Hexanone | Ave | 0.8218 | 0.8839 | 0.1000 | 21.5 | 20.0 | 7.6 | 20.0 |
| Dibromochloromethane | Ave | 0.7724 | 0.7623 | 0.1000 | 9.87 | 10.0 | -1.3 | 20.0 |
| 1,2-Dibromoethane (EDB) | Ave | 1.097 | 1.041 | 0.1000 | 9.49 | 10.0 | -5.1 | 20.0 |
| 3-Chlorobenzotrifluoride | Ave | 1.432 | 1.728 | 0.0100 | 12.1 | 10.0 | 20.7* | 20.0 |
| Chlorobenzene | Ave | 3.204 | 3.213 | 0.5000 | 10.0 | 10.0 | 0.3 | 20.0 |
| 4-Chlorobenzotrifluoride | Ave | 1.328 | 1.597 | 0.0100 | 12.0 | 10.0 | 20.3* | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.9111 | 0.9800 | 0.0100 | 10.8 | 10.0 | 7.6 | 20.0 |
| Ethylbenzene | Ave | 1.833 | 1.767 | 0.1000 | 9.64 | 10.0 | -3.6 | 20.0 |
| m-Xylene & p-Xylene | Ave | 2.246 | 2.173 | 0.1000 | 9.67 | 10.0 | -3.3 | 20.0 |
| o-Xylene | Ave | 2.251 | 2.118 | 0.3000 | 9.41 | 10.0 | -5.9 | 20.0 |
| Styrene | Ave | 3.641 | 3.710 | 0.3000 | 10.2 | 10.0 | 1.9 | 20.0 |
| Bromoform | Ave | 0.4347 | 0.4102 | 0.1000 | 9.44 | 10.0 | -5.6 | 20.0 |
| 2-Chlorobenzotrifluoride | Ave | 1.452 | 1.755 | 0.0100 | 12.1 | 10.0 | 20.9* | 20.0 |
| Isopropylbenzene | Ave | 5.039 | 5.164 | 0.1000 | 10.2 | 10.0 | 2.5 | 20.0 |
| Bromobenzene | Ave | 0.9293 | 0.8562 | 0.0100 | 9.21 | 10.0 | -7.9 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 1.604 | 1.556 | 0.3000 | 9.70 | 10.0 | -3.0 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2720 | 0.2124 | 0.0100 | 7.81 | 10.0 | -21.9* | 20.0 |
| 1,2,3-Trichloropropene | Ave | 0.3860 | 0.3476 | 0.0100 | 9.01 | 10.0 | -9.9 | 20.0 |
| N-Propylbenzene | Ave | 1.011 | 0.9478 | 0.0100 | 9.37 | 10.0 | -6.3 | 20.0 |
| 2-Chlorotoluene | Ave | 0.8762 | 0.8263 | 0.0100 | 9.43 | 10.0 | -5.7 | 20.0 |
| 3-Chlorotoluene | Ave | 0.9194 | 1.039 | 0.0100 | 11.3 | 10.0 | 13.0 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 2.872 | 2.902 | 0.0100 | 10.1 | 10.0 | 1.0 | 20.0 |
| 4-Chlorotoluene | Ave | 0.9565 | 0.9026 | 0.0100 | 9.44 | 10.0 | -5.6 | 20.0 |
| tert-Butylbenzene | Ave | 2.280 | 2.319 | 0.0100 | 10.2 | 10.0 | 1.7 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 2.994 | 3.011 | 0.0100 | 10.1 | 10.0 | 0.6 | 20.0 |
| 3,4-Dichlorobenzotrifluoride | Ave | 0.6644 | 0.7918 | 0.0100 | 11.9 | 10.0 | 19.2 | 20.0 |
| sec-Butylbenzene | Ave | 3.198 | 3.312 | 0.0100 | 10.4 | 10.0 | 3.6 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.669 | 1.619 | 0.6000 | 9.70 | 10.0 | -3.0 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.: _____

Lab Sample ID: CCVIS 180-226148/2 Calibration Date: 10/17/2017 23:55

Instrument ID: CHHP6 Calib Start Date: 07/24/2017 06:39

GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 07/24/2017 09:28

Lab File ID: 6101802D.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 |
|---------------------------------------|------------|---------|--------|---------|-------------|--------------|-------|--------|
| 4-Isopropyltoluene | Ave | 2.641 | 2.790 | 0.0100 | 10.6 | 10.0 | 5.6 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.726 | 1.694 | 0.5000 | 9.81 | 10.0 | -1.9 | 20.0 |
| 2,4-Dichlorobenzotrifluoride | Ave | 0.6261 | 0.7390 | 0.0100 | 11.8 | 10.0 | 18.0 | 20.0 |
| 2,5-Dichlorobenzotrifluoride | Ave | 0.6911 | 0.8749 | 0.0100 | 12.7 | 10.0 | 26.6* | 20.0 |
| n-Butylbenzene | Ave | 2.409 | 2.570 | 0.0100 | 10.7 | 10.0 | 6.7 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.587 | 1.593 | 0.4000 | 10.0 | 10.0 | 0.3 | 20.0 |
| 1,2-Dibromo-3-Chloropropane | Ave | 0.1415 | 0.1284 | 0.0500 | 9.07 | 10.0 | -9.3 | 20.0 |
| 2,4- & 2,5- & 2,6- Dichlorotoluene | Ave | 1.011 | 1.374 | 0.0100 | 40.8 | 30.0 | 35.9* | 20.0 |
| 2,3- & 3,4- Dichlorotoluene | Ave | 1.104 | 1.514 | 0.0100 | 27.4 | 20.0 | 37.1* | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.8294 | 0.9420 | 0.2000 | 11.4 | 10.0 | 13.6 | 20.0 |
| Hexachlorobutadiene | Ave | 0.2626 | 0.3243 | 0.0100 | 12.3 | 10.0 | 23.5* | 20.0 |
| Naphthalene | Ave | 2.528 | 2.478 | 0.0100 | 9.80 | 10.0 | -2.0 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 0.7447 | 0.8532 | 0.0100 | 11.5 | 10.0 | 14.6 | 20.0 |
| 2,4,5-Trichlorotoluene | Ave | 0.4283 | 0.5282 | 0.0100 | 12.3 | 10.0 | 23.3* | 20.0 |
| 2,3,6-Trichlorotoluene | Ave | 0.3839 | 0.5346 | 0.0100 | 13.9 | 10.0 | 39.2* | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2598 | 0.2629 | | 10.1 | 10.0 | 1.2 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.3713 | 0.3771 | | 10.2 | 10.0 | 1.6 | 20.0 |
| Toluene-d8 (Surr) | Lin2 | | 3.866 | | 9.62 | 10.0 | -3.8 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 1.726 | 1.701 | | 9.85 | 10.0 | -1.5 | 20.0 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101802D.D
 Lims ID: CCVIS
 Client ID:
 Sample Type: CCVIS
 Inject. Date: 17-Oct-2017 23:55:30 ALS Bottle#: 2 Worklist Smp#: 2
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018914-002
 Misc. Info.: CCVIS
 Operator ID: 034635 Instrument ID: CHHP6
 Sublist: chrom-MSVOA_LL_CHHP6*sub10
 Method: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 18-Oct-2017 20:30:36 Calib Date: 28-Sep-2017 15:13:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170928-18631.b\\60928P06.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK031

First Level Reviewer: bungardf Date: 18-Oct-2017 00:15:54

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 3.954 | 3.954 | 0.000 | 92 | 360770 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.044 | 7.044 | 0.000 | 98 | 1145374 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.171 | 10.171 | 0.000 | 88 | 290716 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.507 | 12.507 | 0.000 | 96 | 448308 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr) | 113 | 6.314 | 6.314 | 0.000 | 92 | 301086 | 50.0 | 50.6 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur) | 65 | 6.685 | 6.685 | 0.000 | 83 | 431946 | 50.0 | 50.8 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.711 | 8.711 | 0.000 | 93 | 1124001 | 50.0 | 48.1 | |
| \$ 8 4-Bromofluorobenzene (Surr) | 95 | 11.351 | 11.351 | 0.000 | 82 | 494388 | 50.0 | 49.3 | |
| 11 Dichlorodifluoromethane | 85 | 1.472 | 1.472 | 0.000 | 99 | 414133 | 50.0 | 56.7 | |
| 12 Chloromethane | 50 | 1.624 | 1.624 | 0.000 | 97 | 279627 | 50.0 | 42.5 | |
| 13 Vinyl chloride | 62 | 1.740 | 1.740 | 0.000 | 97 | 354290 | 50.0 | 51.4 | |
| 14 Butadiene | 39 | 1.788 | 1.788 | 0.000 | 92 | 278145 | 50.0 | 48.7 | |
| 15 Bromomethane | 94 | 2.068 | 2.068 | 0.000 | 92 | 174522 | 50.0 | 54.3 | |
| 16 Chloroethane | 64 | 2.202 | 2.202 | 0.000 | 98 | 202206 | 50.0 | 54.3 | |
| 17 Dichlorofluoromethane | 67 | 2.457 | 2.457 | 0.000 | 96 | 450565 | 50.0 | 56.2 | |
| 18 Trichlorofluoromethane | 101 | 2.470 | 2.470 | 0.000 | 93 | 470086 | 50.0 | 69.3 | |
| 20 Ethyl ether | 59 | 2.816 | 2.816 | 0.000 | 83 | 290326 | 50.0 | 50.0 | |
| 21 Acrolein | 56 | 2.981 | 2.981 | 0.000 | 99 | 176944 | 150.0 | 142.3 | |
| 22 1,1-Dichloroethene | 96 | 3.090 | 3.090 | 0.000 | 98 | 317197 | 50.0 | 53.3 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.151 | 3.151 | 0.000 | 95 | 347303 | 50.0 | 60.9 | |
| 24 Acetone | 43 | 3.181 | 3.181 | 0.000 | 99 | 332642 | 100.0 | 135.4 | |
| 25 Iodomethane | 142 | 3.273 | 3.273 | 0.000 | 99 | 436319 | 50.0 | 52.1 | |
| 26 Carbon disulfide | 76 | 3.364 | 3.364 | 0.000 | 99 | 659410 | 50.0 | 49.7 | |
| 29 3-Chloro-1-propene | 76 | 3.631 | 3.631 | 0.000 | 85 | 168033 | 50.0 | 48.2 | |
| 30 Methyl acetate | 43 | 3.644 | 3.644 | 0.000 | 96 | 424505 | 100.0 | 78.8 | |
| 31 Methylene Chloride | 84 | 3.838 | 3.838 | 0.000 | 82 | 388136 | 50.0 | 49.2 | |
| 32 2-Methyl-2-propanol | 59 | 4.094 | 4.094 | 0.000 | 91 | 188867 | 500.0 | 473.0 | |
| 33 Acrylonitrile | 53 | 4.222 | 4.222 | 0.000 | 98 | 1084711 | 500.0 | 377.0 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.270 | 4.270 | 0.000 | 97 | 350195 | 50.0 | 51.8 | |
| 35 Methyl tert-butyl ether | 73 | 4.276 | 4.276 | 0.000 | 94 | 991746 | 50.0 | 45.9 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 36 Hexane | 57 | 4.696 | 4.696 | 0.000 | 89 | 383088 | 50.0 | 47.9 | |
| 37 1,1-Dichloroethane | 63 | 4.921 | 4.921 | 0.000 | 96 | 559012 | 50.0 | 50.9 | |
| 38 Vinyl acetate | 43 | 4.976 | 4.976 | 0.000 | 97 | 477750 | 50.0 | 37.3 | |
| 42 2,2-Dichloropropane | 97 | 5.676 | 5.676 | 0.000 | 62 | 67006 | 50.0 | 58.9 | |
| 43 cis-1,2-Dichloroethene | 96 | 5.688 | 5.688 | 0.000 | 78 | 396522 | 50.0 | 50.3 | |
| 44 2-Butanone (MEK) | 43 | 5.694 | 5.694 | 0.000 | 88 | 371245 | 100.0 | 105.4 | |
| 48 Chlorobromomethane | 128 | 5.974 | 5.974 | 0.000 | 87 | 177171 | 50.0 | 51.5 | |
| 49 Tetrahydrofuran | 42 | 5.992 | 5.992 | 0.000 | 80 | 153635 | 100.0 | 64.4 | |
| 50 Chloroform | 83 | 6.126 | 6.126 | 0.000 | 92 | 624295 | 50.0 | 52.7 | |
| 51 1,1,1-Trichloroethane | 97 | 6.290 | 6.290 | 0.000 | 98 | 444403 | 50.0 | 59.0 | |
| 52 Cyclohexane | 56 | 6.357 | 6.357 | 0.000 | 81 | 480001 | 50.0 | 44.9 | |
| 53 Carbon tetrachloride | 117 | 6.454 | 6.454 | 0.000 | 95 | 364006 | 50.0 | 67.2 | |
| 54 1,1-Dichloropropene | 75 | 6.479 | 6.479 | 0.000 | 98 | 492726 | 50.0 | 54.3 | |
| 56 Benzene | 78 | 6.691 | 6.691 | 0.000 | 97 | 1403557 | 50.0 | 53.2 | |
| 55 Isobutyl alcohol | 41 | 6.691 | 6.691 | 0.000 | 62 | 171526 | 1250.0 | 1079.6 | |
| 57 1,2-Dichloroethane | 62 | 6.777 | 6.777 | 0.000 | 99 | 501890 | 50.0 | 51.5 | |
| 59 n-Heptane | 43 | 7.069 | 7.069 | 0.000 | 81 | 265500 | 50.0 | 43.6 | |
| 61 Trichloroethene | 130 | 7.440 | 7.440 | 0.000 | 97 | 331659 | 50.0 | 51.7 | |
| 63 Methylcyclohexane | 83 | 7.671 | 7.671 | 0.000 | 84 | 577855 | 50.0 | 51.6 | |
| 64 1,2-Dichloropropane | 63 | 7.713 | 7.713 | 0.000 | 91 | 302111 | 50.0 | 45.8 | |
| 65 1,4-Dioxane | 88 | 7.793 | 7.793 | 0.000 | 38 | 50488 | 1000.0 | 838.9 | M |
| 67 Dibromomethane | 93 | 7.799 | 7.799 | 0.000 | 93 | 218269 | 50.0 | 50.6 | |
| 68 Dichlorobromomethane | 83 | 7.999 | 7.999 | 0.000 | 99 | 368884 | 50.0 | 50.2 | |
| 70 2-Chloroethyl vinyl ether | 63 | 8.310 | 8.310 | 0.000 | 93 | 450947 | 100.0 | 103.1 | |
| 71 cis-1,3-Dichloropropene | 75 | 8.450 | 8.450 | 0.000 | 96 | 420558 | 50.0 | 50.5 | |
| 72 4-Methyl-2-pentanone (MIBK) | 43 | 8.608 | 8.608 | 0.000 | 92 | 563691 | 100.0 | 75.9 | |
| 73 Toluene | 91 | 8.778 | 8.778 | 0.000 | 99 | 1403520 | 50.0 | 49.0 | |
| 74 trans-1,3-Dichloropropene | 75 | 9.034 | 9.034 | 0.000 | 91 | 375688 | 50.0 | 49.0 | |
| 75 Ethyl methacrylate | 69 | 9.100 | 9.100 | 0.000 | 85 | 435415 | 50.0 | 45.0 | |
| 76 1,1,2-Trichloroethane | 97 | 9.222 | 9.222 | 0.000 | 92 | 317088 | 50.0 | 47.8 | |
| 77 Tetrachloroethene | 164 | 9.295 | 9.295 | 0.000 | 93 | 244830 | 50.0 | 47.7 | |
| 78 1,3-Dichloropropane | 76 | 9.380 | 9.380 | 0.000 | 87 | 587805 | 50.0 | 48.9 | |
| 79 2-Hexanone | 43 | 9.447 | 9.447 | 0.000 | 92 | 513905 | 100.0 | 107.6 | |
| 81 Chlorodibromomethane | 129 | 9.593 | 9.593 | 0.000 | 89 | 221611 | 50.0 | 49.3 | |
| 82 Ethylene Dibromide | 107 | 9.709 | 9.709 | 0.000 | 99 | 302607 | 50.0 | 47.5 | |
| 83 3-Chlorobenzotrifluoride | 180 | 10.183 | 10.183 | 0.000 | 89 | 502373 | 50.0 | 60.3 | |
| 84 Chlorobenzene | 112 | 10.196 | 10.196 | 0.000 | 93 | 934005 | 50.0 | 50.1 | |
| 85 4-Chlorobenzotrifluoride | 180 | 10.269 | 10.269 | 0.000 | 96 | 464194 | 50.0 | 60.1 | |
| 86 1,1,1,2-Tetrachloroethane | 131 | 10.293 | 10.293 | 0.000 | 89 | 284886 | 50.0 | 53.8 | |
| 87 Ethylbenzene | 106 | 10.299 | 10.299 | 0.000 | 98 | 513832 | 50.0 | 48.2 | |
| 88 m-Xylene & p-Xylene | 106 | 10.433 | 10.433 | 0.000 | 99 | 631786 | 50.0 | 48.4 | |
| 89 o-Xylene | 106 | 10.816 | 10.816 | 0.000 | 96 | 615703 | 50.0 | 47.0 | |
| 90 Styrene | 104 | 10.834 | 10.834 | 0.000 | 93 | 1078665 | 50.0 | 51.0 | |
| 91 Bromoform | 173 | 11.011 | 11.011 | 0.000 | 92 | 119244 | 50.0 | 47.2 | |
| 92 2-Chlorobenzotrifluoride | 180 | 11.084 | 11.084 | 0.000 | 93 | 510216 | 50.0 | 60.4 | |
| 93 Isopropylbenzene | 105 | 11.181 | 11.181 | 0.000 | 97 | 1501372 | 50.0 | 51.2 | |
| 95 Bromobenzene | 156 | 11.485 | 11.485 | 0.000 | 95 | 383822 | 50.0 | 46.1 | |
| 96 1,1,2,2-Tetrachloroethane | 83 | 11.497 | 11.497 | 0.000 | 96 | 452431 | 50.0 | 48.5 | |
| 97 trans-1,4-Dichloro-2-butene | 53 | 11.540 | 11.540 | 0.000 | 72 | 95229 | 50.0 | 39.1 | |
| 98 1,2,3-Trichloropropane | 110 | 11.546 | 11.546 | 0.000 | 86 | 155841 | 50.0 | 45.0 | |
| 99 N-Propylbenzene | 120 | 11.595 | 11.595 | 0.000 | 98 | 424892 | 50.0 | 46.9 | |
| 100 2-Chlorotoluene | 126 | 11.680 | 11.680 | 0.000 | 95 | 370449 | 50.0 | 47.2 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 101 3-Chlorotoluene | 126 | 11.747 | 11.747 | 0.000 | 96 | 465929 | 50.0 | 56.5 | |
| 102 1,3,5-Trimethylbenzene | 105 | 11.783 | 11.783 | 0.000 | 93 | 1300833 | 50.0 | 50.5 | |
| 103 4-Chlorotoluene | 126 | 11.802 | 11.802 | 0.000 | 99 | 404640 | 50.0 | 47.2 | |
| 104 tert-Butylbenzene | 119 | 12.094 | 12.094 | 0.000 | 91 | 1039835 | 50.0 | 50.9 | |
| 106 1,2,4-Trimethylbenzene | 105 | 12.154 | 12.154 | 0.000 | 98 | 1349978 | 50.0 | 50.3 | |
| 107 1,2-dichloro-4-(trifluoromethyl) | 214 | 12.203 | 12.203 | 0.000 | 93 | 354962 | 50.0 | 59.6 | |
| 108 sec-Butylbenzene | 105 | 12.319 | 12.319 | 0.000 | 96 | 1484959 | 50.0 | 51.8 | |
| 109 1,3-Dichlorobenzene | 146 | 12.428 | 12.428 | 0.000 | 95 | 725865 | 50.0 | 48.5 | |
| 110 4-Isopropyltoluene | 119 | 12.471 | 12.471 | 0.000 | 95 | 1250691 | 50.0 | 52.8 | |
| 111 1,4-Dichlorobenzene | 146 | 12.532 | 12.532 | 0.000 | 93 | 759417 | 50.0 | 49.1 | |
| 113 2,4-Dichloro-1-(trifluoromethyl) | 214 | 12.568 | 12.568 | 0.000 | 94 | 331314 | 50.0 | 59.0 | |
| 114 2,5-Dichlorobenzotrifluoride | 214 | 12.611 | 12.611 | 0.000 | 97 | 392222 | 50.0 | 63.3 | |
| 116 n-Butylbenzene | 91 | 12.878 | 12.878 | 0.000 | 96 | 1152283 | 50.0 | 53.4 | |
| 117 1,2-Dichlorobenzene | 146 | 12.890 | 12.890 | 0.000 | 95 | 713947 | 50.0 | 50.2 | |
| 118 1,2-Dibromo-3-Chloropropan | 75 | 13.681 | 13.675 | 0.006 | 79 | 57542 | 50.0 | 45.4 | |
| 119 2,4- & 2,5- & 2,6- Dichlorobenzene | 125 | 13.815 | 13.815 | 0.000 | 98 | 1848132 | 150.0 | 203.9 | |
| 121 2,3- & 3,4- Dichlorotoluene | 125 | 14.235 | 14.235 | 0.000 | 99 | 1357764 | 100.0 | 137.1 | |
| 122 1,2,4-Trichlorobenzene | 180 | 14.497 | 14.497 | 0.000 | 93 | 422291 | 50.0 | 56.8 | |
| 123 Hexachlorobutadiene | 225 | 14.643 | 14.643 | 0.000 | 94 | 145382 | 50.0 | 61.7 | |
| 124 Naphthalene | 128 | 14.758 | 14.758 | 0.000 | 98 | 1110813 | 50.0 | 49.0 | |
| 125 1,2,3-Trichlorobenzene | 180 | 14.977 | 14.977 | 0.000 | 97 | 382474 | 50.0 | 57.3 | |
| 126 2,4,5-Trichlorotoluene | 159 | 15.774 | 15.774 | 0.000 | 0 | 236806 | 50.0 | 61.7 | |
| 127 2,3,6-Trichlorotoluene | 159 | 15.877 | 15.877 | 0.000 | 95 | 239668 | 50.0 | 69.6 | |
| 146 3,4-Dichlorotoluene | 1 | 0.000 | | | | | ND | ND | |
| S 131 Xylenes, Total | 106 | | | | 0 | | 100.0 | 95.4 | |
| S 130 1,2-Dichloroethene, Total | 96 | | | | 0 | | 100.0 | 102.1 | |
| S 132 1,3-Dichloropropene, Total | 1 | | | | 0 | | 100.0 | 99.4 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

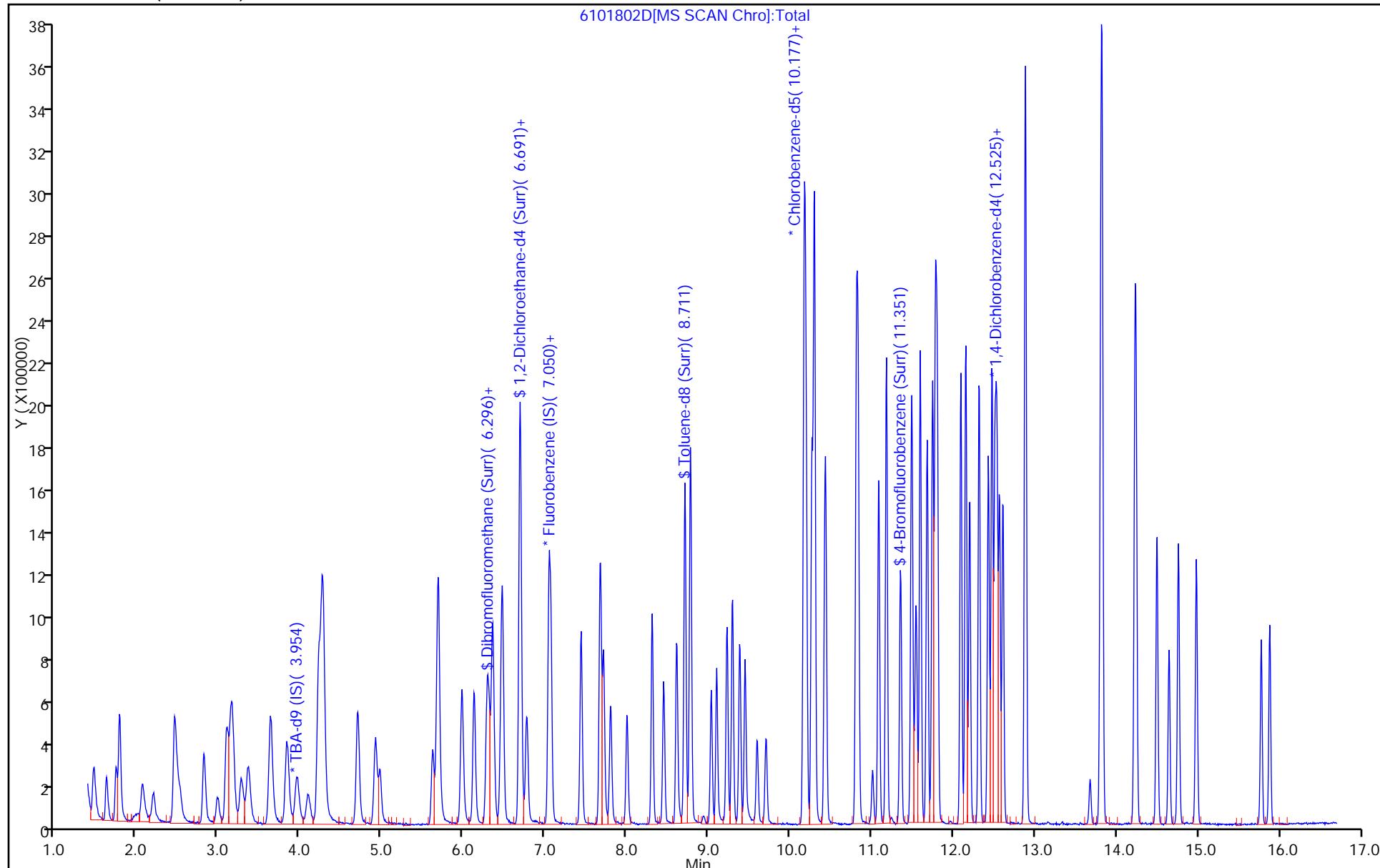
Reagents:

| | | | |
|---------------------|--------------------|-----------|-------------|
| voaWKetmix1st_00006 | Amount Added: 2.00 | Units: uL | |
| voaWVA1stRest_00021 | Amount Added: 2.00 | Units: uL | |
| voaWAcro1stRe_00021 | Amount Added: 6.00 | Units: uL | |
| voaWEEmix1stR_00014 | Amount Added: 2.00 | Units: uL | |
| VOA2CEVE2ND_00010 | Amount Added: 2.00 | Units: uL | |
| VOA8260VOAPRI_00266 | Amount Added: 2.00 | Units: uL | |
| VOA8260INT_00074 | Amount Added: 2.00 | Units: uL | Run Reagent |
| VOA8260SURR_00073 | Amount Added: 2.00 | Units: uL | Run Reagent |

Report Date: 18-Oct-2017 20:30:37

Chrom Revision: 2.2 16-Aug-2017 16:24:46

TestAmerica Pittsburgh
Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101802D.D
Injection Date: 17-Oct-2017 23:55:30 Instrument ID: CHHP6
Lims ID: CCVIS Operator ID: 034635
Client ID:
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 2
Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh

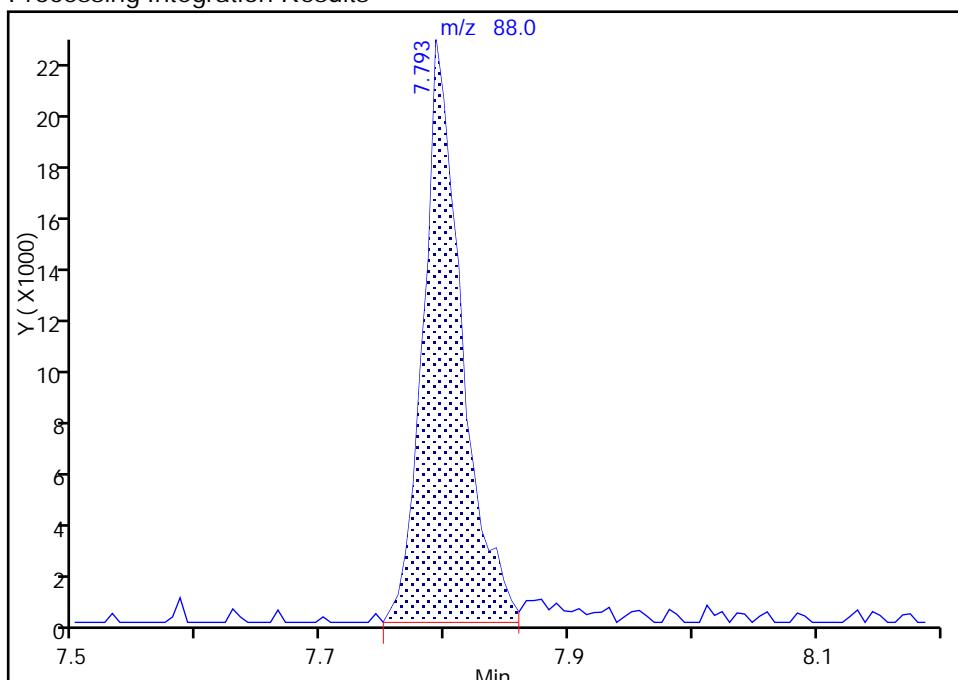
Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101802D.D
 Injection Date: 17-Oct-2017 23:55:30 Instrument ID: CHHP6
 Lims ID: CCVIS
 Client ID:
 Operator ID: 034635 ALS Bottle#: 2 Worklist Smp#: 2
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
 Column: DB-624 (0.18 mm) Detector: MS SCAN

65 1,4-Dioxane, CAS: 123-91-1

Signal: 1

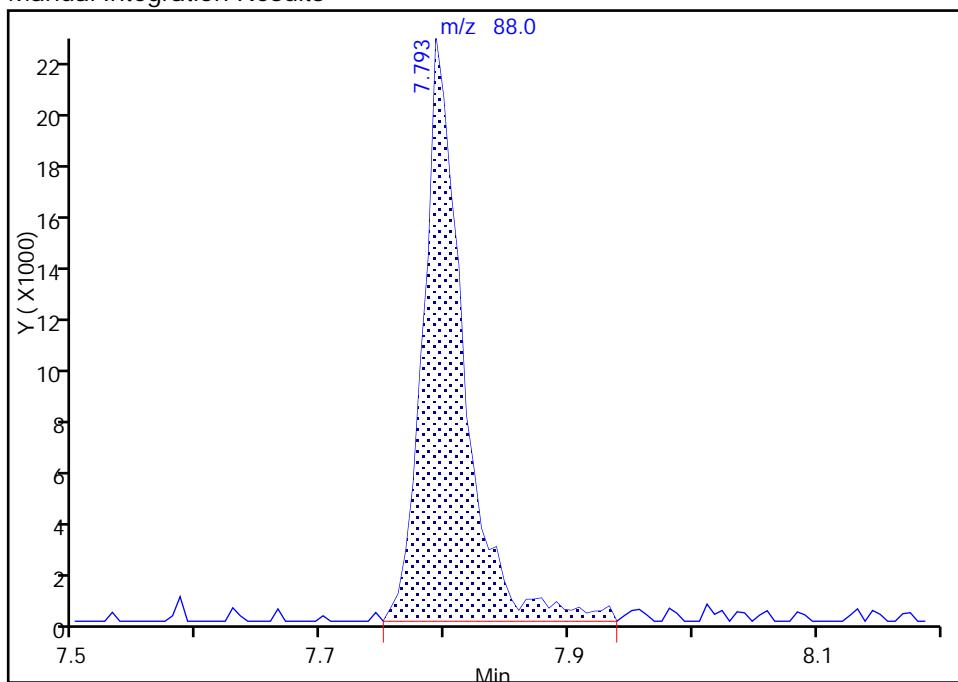
RT: 7.79
 Area: 48010
 Amount: 797.7345
 Amount Units: ng

Processing Integration Results



RT: 7.79
 Area: 50488
 Amount: 838.9089
 Amount Units: ng

Manual Integration Results



Reviewer: bungardf, 18-Oct-2017 00:15:52

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170602-17015.b\60602003.D
 Lims ID: BFB
 Client ID:
 Sample Type: BFB
 Inject. Date: 02-Jun-2017 06:02:30 ALS Bottle#: 1 Worklist Smp#: 3
 Injection Vol: 5.0 mL Dil. Factor: 1.0000
 Sample Info: 180-0017015-003
 Misc. Info.: BFB
 Operator ID: 034635 Instrument ID: CHHP6
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170602-17015.b\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 05-Jun-2017 02:18:35 Calib Date: 02-Jun-2017 14:26:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170602-17015.b\60602021.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK004

First Level Reviewer: bungardf Date: 02-Jun-2017 06:25:15

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------|-----|-----------|---------------|---------------|---|----------|------------|--------------|-------|
|----------|-----|-----------|---------------|---------------|---|----------|------------|--------------|-------|

\$ 10 BFB

95 8.213 8.213 0.000 0 92923

NR NR

QC Flag Legend

Processing Flags

NR - Missing Quant Standard

Reagents:

VOABFB25_00088

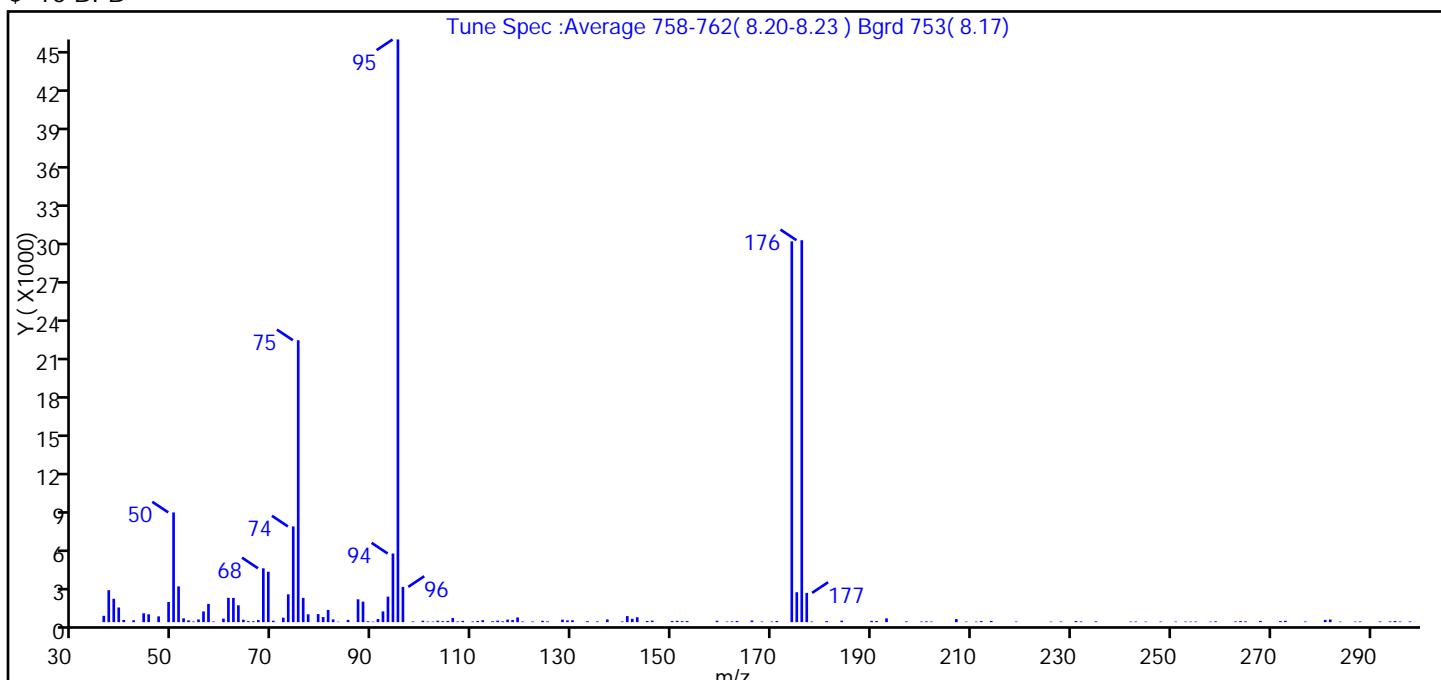
Amount Added: 1.00

Units: uL

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170602-17015.b\\60602003.D
 Injection Date: 02-Jun-2017 06:02:30 Instrument ID: CHHP6
 Lims ID: BFB
 Client ID:
 Operator ID: 034635 ALS Bottle#: 1 Worklist Smp#: 3
 Injection Vol: 5.0 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 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 | 18.8 |
| 75 | 30 to 60% of m/z 95 | 48.4 |
| 96 | 5 to 9% of m/z 95 | 6.0 |
| 173 | Less than 2% of m/z 174 | 0.0 (0.0) |
| 174 | 50 to 120% of m/z 95 | 65.4 |
| 175 | 5 to 9% of m/z 174 | 5.1 (7.9) |
| 176 | Greater than 95% but less than 101% of m/z 174 | 65.5 (100.3) |
| 177 | 5 to 9% of m/z 176 | 5.0 (7.7) |

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170602-17015.b\\60602003.D\MSVOA_LL_CHHP6.rslt\spec
 Injection Date: 02-Jun-2017 06:02:30
 Spectrum: Tune Spec :Average 758-762(8.20-8.23) Bgrd 753(8.17)
 Base Peak: 95.00
 Minimum % Base Peak: 0
 Number of Points: 151

| m/z | Y | m/z | Y | m/z | Y | m/z | Y |
|-------|------|--------|-------|--------|-------|--------|-----|
| 36.00 | 491 | 81.00 | 948 | 129.00 | 138 | 209.00 | 41 |
| 37.00 | 2479 | 82.00 | 215 | 130.00 | 148 | 211.00 | 44 |
| 38.00 | 1815 | 83.00 | 32 | 133.00 | 75 | 212.00 | 81 |
| 39.00 | 1138 | 85.00 | 175 | 135.00 | 55 | 214.00 | 91 |
| 40.00 | 168 | 87.00 | 1772 | 137.00 | 211 | 219.00 | 45 |
| 42.00 | 145 | 88.00 | 1592 | 140.00 | 55 | 226.00 | 40 |
| 44.00 | 690 | 89.00 | 69 | 141.00 | 469 | 228.00 | 46 |
| 45.00 | 609 | 90.00 | 40 | 142.00 | 261 | 231.00 | 88 |
| 47.00 | 446 | 91.00 | 242 | 143.00 | 380 | 232.00 | 56 |
| 49.00 | 1560 | 92.00 | 834 | 145.00 | 88 | 235.00 | 63 |
| 50.00 | 8513 | 93.00 | 1980 | 146.00 | 118 | 242.00 | 45 |
| 51.00 | 2775 | 94.00 | 5325 | 150.00 | 86 | 243.00 | 48 |
| 52.00 | 295 | 95.00 | 45208 | 151.00 | 105 | 245.00 | 40 |
| 53.00 | 139 | 96.00 | 2735 | 152.00 | 77 | 248.00 | 43 |
| 54.00 | 49 | 98.00 | 55 | 153.00 | 84 | 251.00 | 43 |
| 55.00 | 209 | 100.00 | 111 | 159.00 | 109 | 253.00 | 44 |
| 56.00 | 833 | 101.00 | 45 | 161.00 | 43 | 254.00 | 41 |
| 57.00 | 1414 | 102.00 | 40 | 162.00 | 57 | 255.00 | 47 |
| 58.00 | 48 | 103.00 | 112 | 163.00 | 88 | 258.00 | 42 |
| 60.00 | 273 | 104.00 | 73 | 166.00 | 126 | 259.00 | 63 |
| 61.00 | 1885 | 105.00 | 94 | 168.00 | 46 | 263.00 | 55 |
| 62.00 | 1879 | 106.00 | 320 | 170.00 | 47 | 264.00 | 81 |
| 63.00 | 1308 | 107.00 | 54 | 171.00 | 85 | 265.00 | 62 |
| 64.00 | 188 | 108.00 | 101 | 174.00 | 29544 | 268.00 | 97 |
| 65.00 | 77 | 110.00 | 52 | 175.00 | 2320 | 272.00 | 83 |
| 66.00 | 68 | 111.00 | 89 | 176.00 | 29632 | 273.00 | 106 |
| 67.00 | 162 | 112.00 | 154 | 177.00 | 2269 | 277.00 | 48 |
| 68.00 | 4177 | 114.00 | 68 | 178.00 | 48 | 281.00 | 177 |
| 69.00 | 3914 | 115.00 | 118 | 181.00 | 82 | 282.00 | 204 |
| 70.00 | 109 | 116.00 | 61 | 184.00 | 116 | 284.00 | 40 |
| 72.00 | 348 | 117.00 | 198 | 190.00 | 102 | 287.00 | 40 |
| 73.00 | 2157 | 118.00 | 161 | 191.00 | 84 | 288.00 | 52 |
| 74.00 | 7426 | 119.00 | 366 | 193.00 | 294 | 292.00 | 56 |

Report Date: 05-Jun-2017 02:18:37

Chrom Revision: 2.2 25-Apr-2017 13:27:22

Data File:

\\\ChromNA\Pittsburgh\ChromData\CHHP6\20170602-17015.b\60602003.D\MSVOA_LL_CHHP6.rslt\spec

Injection Date:

02-Jun-2017 06:02:30

Spectrum:

Tune Spec :Average 758-762(8.20-8.23) Bgrd 753(8.17)

Base Peak:

95.00

Minimum % Base Peak: 0

Number of Points: 151

| m/z | Y | m/z | Y | m/z | Y | m/z | Y |
|-------|-------|--------|-----|--------|-----|--------|----|
| 75.00 | 21872 | 120.00 | 51 | 197.00 | 60 | 294.00 | 57 |
| 76.00 | 1884 | 122.00 | 51 | 200.00 | 48 | 295.00 | 90 |
| 77.00 | 608 | 124.00 | 107 | 201.00 | 63 | 296.00 | 47 |
| 79.00 | 627 | 125.00 | 60 | 202.00 | 44 | 298.00 | 54 |
| 80.00 | 394 | 128.00 | 194 | 207.00 | 235 | | |

Report Date: 05-Jun-2017 02:18:37

Chrom Revision: 2.2 25-Apr-2017 13:27:22

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170602-17015.b\\60602003.D

Injection Date: 02-Jun-2017 06:02:30

Instrument ID: CHHP6

Operator ID: 034635

Lims ID: BFB

Worklist Smp#: 3

Client ID:

Injection Vol: 5.0 mL

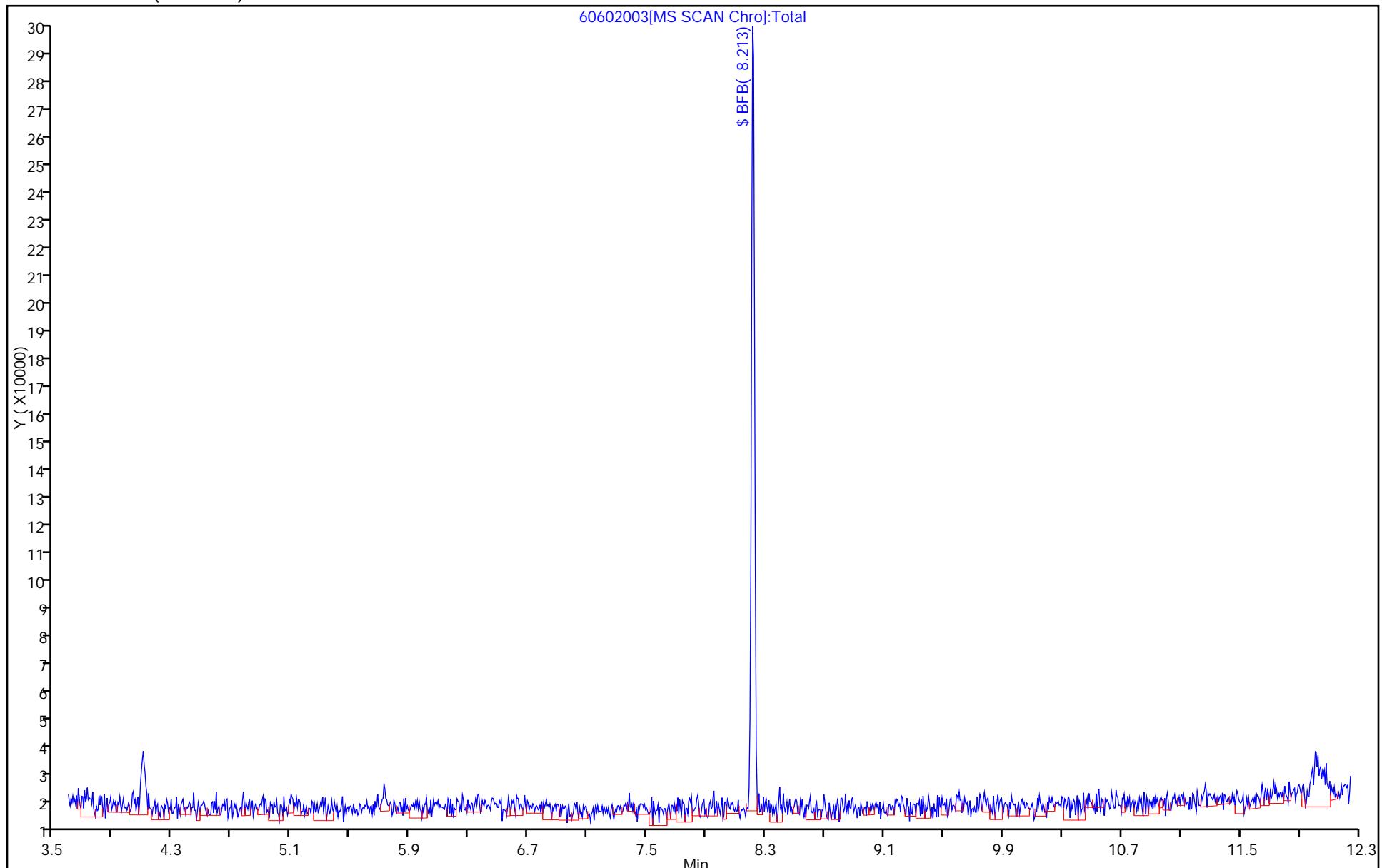
Dil. Factor: 1.0000

ALS Bottle#: 1

Method: MSVOA_LL_CHHP6

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170724-17705.b\60724D01.D
 Lims ID: BFB
 Client ID:
 Sample Type: BFB
 Inject. Date: 24-Jul-2017 04:58:30 ALS Bottle#: 1 Worklist Smp#: 1
 Injection Vol: 5.0 mL Dil. Factor: 1.0000
 Sample Info: 180-0017705-001
 Misc. Info.: BFB
 Operator ID: 034635 Instrument ID: CHHP6
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170724-17705.b\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 25-Jul-2017 01:44:23 Calib Date: 24-Jul-2017 09:28:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170724-17705.b\60724D10.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK021

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------|-----|-----------|---------------|---------------|---|----------|------------|--------------|-------|
|----------|-----|-----------|---------------|---------------|---|----------|------------|--------------|-------|

| | | | | | | | | |
|-----------|----|-------|-------|-------|---|-------|----|----|
| \$ 10 BFB | 95 | 8.182 | 8.182 | 0.000 | 0 | 80019 | NR | NR |
|-----------|----|-------|-------|-------|---|-------|----|----|

QC Flag Legend

Processing Flags

NR - Missing Quant Standard

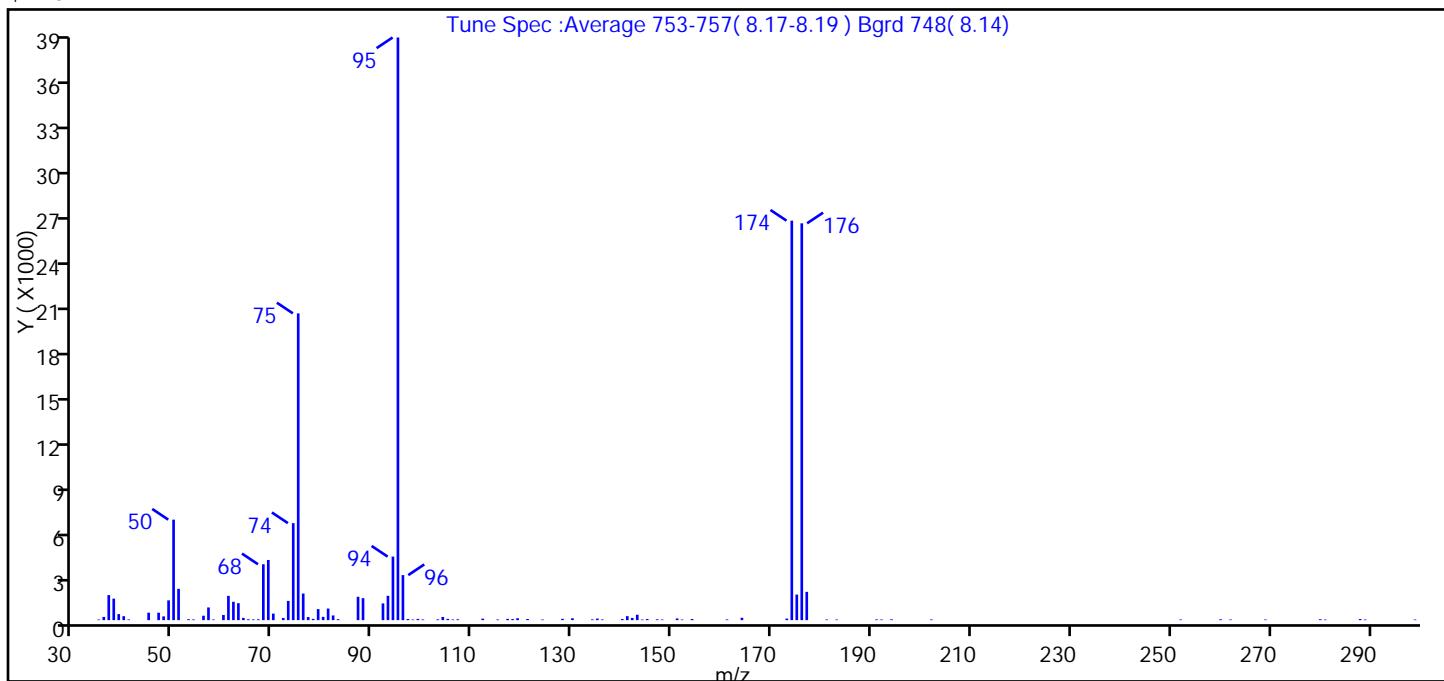
Reagents:

VOABFB25_00090 Amount Added: 1.00 Units: uL

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D01.D
 Injection Date: 24-Jul-2017 04:58:30 Instrument ID: CHHP6
 Lims ID: BFB
 Client ID:
 Operator ID: 034635 ALS Bottle#: 1 Worklist Smp#: 1
 Injection Vol: 5.0 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 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 | 17.3 |
| 75 | 30 to 60% of m/z 95 | 52.6 |
| 96 | 5 to 9% of m/z 95 | 7.7 |
| 173 | Less than 2% of m/z 174 | 0.3 (0.4) |
| 174 | 50 to 120% of m/z 95 | 68.6 |
| 175 | 5 to 9% of m/z 174 | 4.4 (6.4) |
| 176 | Greater than 95% but less than 101% of m/z 174 | 68.1 (99.3) |
| 177 | 5 to 9% of m/z 176 | 4.9 (7.1) |

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170724-17705.b\\60724D01.D\\MSVOA_LL_CHHP6.rslt\\spec
 Injection Date: 24-Jul-2017 04:58:30
 Spectrum: Tune Spec :Average 753-757(8.17-8.19) Bgrd 748(8.14)
 Base Peak: 95.00
 Minimum % Base Peak: 0
 Number of Points: 102

| m/z | Y | m/z | Y | m/z | Y | m/z | Y |
|-------|------|--------|-------|--------|-----|--------|-------|
| 35.00 | 41 | 68.00 | 3693 | 103.00 | 52 | 152.00 | 40 |
| 36.00 | 211 | 69.00 | 3974 | 104.00 | 211 | 154.00 | 76 |
| 37.00 | 1655 | 70.00 | 433 | 105.00 | 85 | 161.00 | 49 |
| 38.00 | 1420 | 72.00 | 145 | 106.00 | 41 | 164.00 | 157 |
| 39.00 | 400 | 73.00 | 1271 | 107.00 | 54 | 173.00 | 103 |
| 40.00 | 262 | 74.00 | 6411 | 112.00 | 99 | 174.00 | 26400 |
| 41.00 | 45 | 75.00 | 20272 | 115.00 | 45 | 175.00 | 1689 |
| 45.00 | 492 | 76.00 | 1760 | 117.00 | 85 | 176.00 | 26224 |
| 47.00 | 490 | 77.00 | 205 | 118.00 | 72 | 177.00 | 1870 |
| 48.00 | 248 | 78.00 | 83 | 119.00 | 127 | 181.00 | 43 |
| 49.00 | 1307 | 79.00 | 730 | 121.00 | 74 | 183.00 | 41 |
| 50.00 | 6642 | 80.00 | 217 | 124.00 | 40 | 191.00 | 49 |
| 51.00 | 2068 | 81.00 | 766 | 128.00 | 87 | 192.00 | 42 |
| 53.00 | 63 | 82.00 | 309 | 130.00 | 124 | 194.00 | 54 |
| 54.00 | 47 | 83.00 | 65 | 134.00 | 50 | 202.00 | 45 |
| 56.00 | 292 | 87.00 | 1545 | 135.00 | 104 | 252.00 | 43 |
| 57.00 | 839 | 88.00 | 1453 | 136.00 | 44 | 260.00 | 58 |
| 58.00 | 42 | 92.00 | 1107 | 140.00 | 73 | 262.00 | 42 |
| 60.00 | 343 | 93.00 | 1608 | 141.00 | 265 | 269.00 | 42 |
| 61.00 | 1604 | 94.00 | 4199 | 142.00 | 143 | 280.00 | 58 |
| 62.00 | 1212 | 95.00 | 38504 | 143.00 | 358 | 281.00 | 40 |
| 63.00 | 1119 | 96.00 | 2976 | 144.00 | 40 | 288.00 | 73 |
| 64.00 | 149 | 97.00 | 76 | 145.00 | 73 | 289.00 | 43 |
| 65.00 | 56 | 98.00 | 40 | 147.00 | 56 | 299.00 | 42 |
| 66.00 | 47 | 99.00 | 82 | 148.00 | 40 | | |
| 67.00 | 56 | 100.00 | 44 | 151.00 | 113 | | |

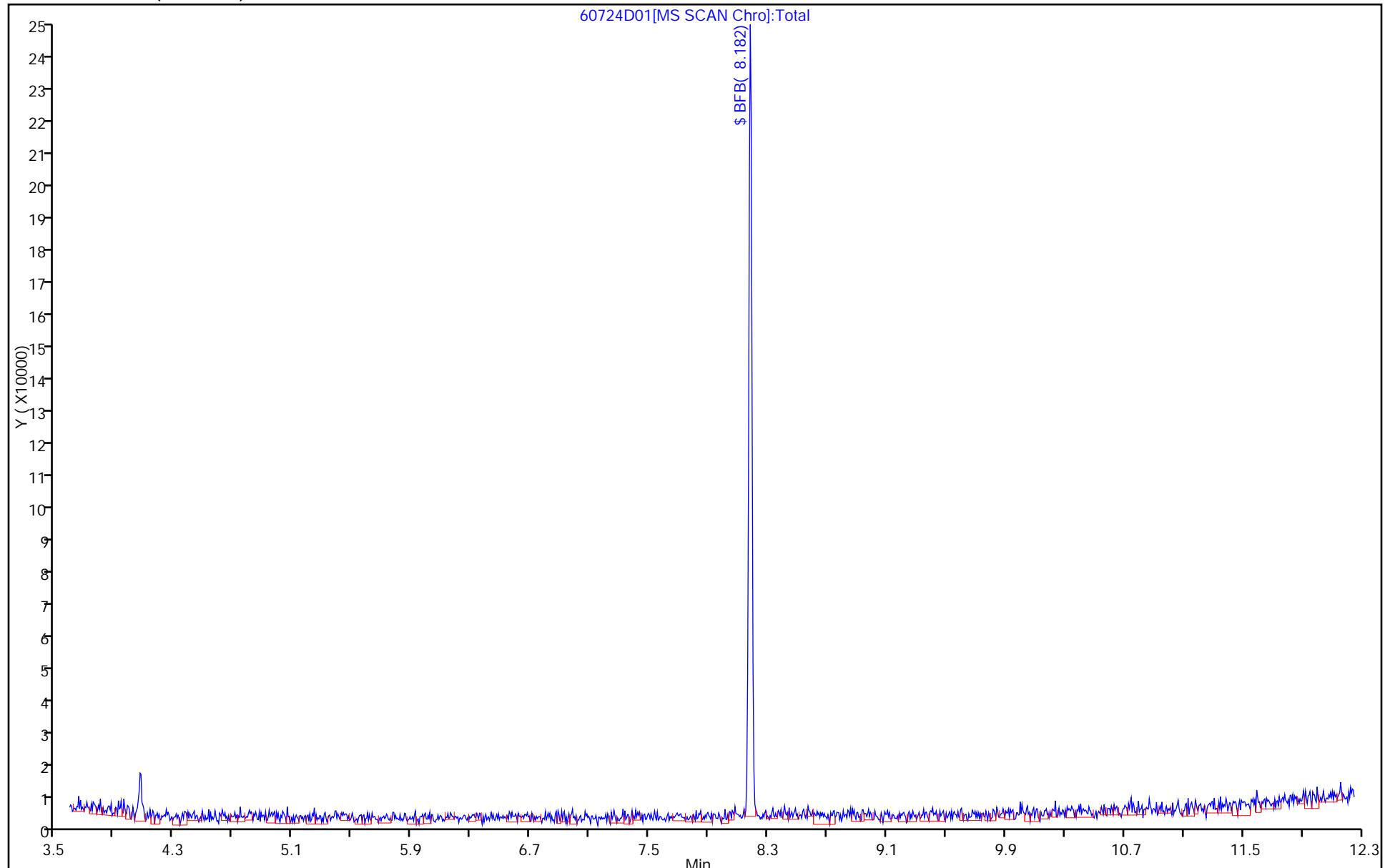
Report Date: 25-Jul-2017 01:44:25

Chrom Revision: 2.2 20-Jun-2017 07:42:38

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20170724-17705.b\\60724D01.D
Injection Date: 24-Jul-2017 04:58:30 Instrument ID: CHHP6
Lims ID: BFB Operator ID: 034635
Client ID:
Injection Vol: 5.0 mL Dil. Factor: 1.0000 ALS Bottle#: 1
Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm)

Worklist Smp#: 1



TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\6101801D.D
 Lims ID: BFB
 Client ID:
 Sample Type: BFB
 Inject. Date: 17-Oct-2017 21:58:30 ALS Bottle#: 1 Worklist Smp#: 1
 Injection Vol: 5.0 mL Dil. Factor: 1.0000
 Sample Info: 180-0018914-001
 Misc. Info.:
 Operator ID: 034635 Instrument ID: CHHP6
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 18-Oct-2017 20:30:34 Calib Date: 28-Sep-2017 15:13:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICAL File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170928-18631.b\60928P06.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK031

First Level Reviewer: bungardf Date: 17-Oct-2017 22:50:15

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------|-----|-----------|---------------|---------------|---|----------|------------|--------------|-------|
|----------|-----|-----------|---------------|---------------|---|----------|------------|--------------|-------|

\$ 10 BFB

95 8.183 8.183 0.000 0 63554

NR NR

QC Flag Legend

Processing Flags

NR - Missing Quant Standard

Reagents:

VOABFB25_00094

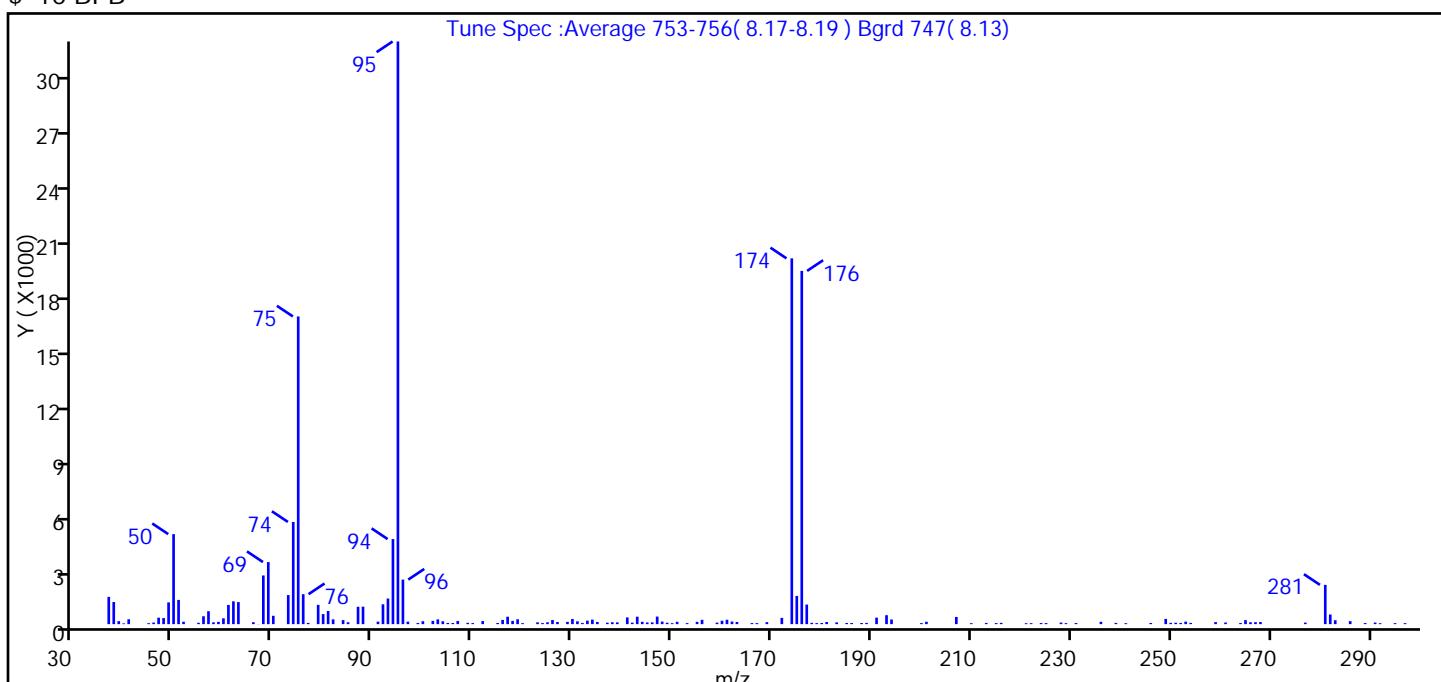
Amount Added: 1.00

Units: uL

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101801D.D
 Injection Date: 17-Oct-2017 21:58:30 Instrument ID: CHHP6
 Lims ID: BFB
 Client ID:
 Operator ID: 034635 ALS Bottle#: 1 Worklist Smp#: 1
 Injection Vol: 5.0 mL Dil. Factor: 1.0000
 Method: MSVOA_LL_CHHP6 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 | 15.4 |
| 75 | 30 to 60% of m/z 95 | 52.8 |
| 96 | 5 to 9% of m/z 95 | 7.6 |
| 173 | Less than 2% of m/z 174 | 0.0 (0.0) |
| 174 | 50 to 120% of m/z 95 | 62.8 |
| 175 | 5 to 9% of m/z 174 | 4.8 (7.7) |
| 176 | Greater than 95% but less than 101% of m/z 174 | 60.6 (96.6) |
| 177 | 5 to 9% of m/z 176 | 3.4 (5.5) |

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\6101801D.D\MSVOA_LL_CHHP6.rslt\spec
 Injection Date: 17-Oct-2017 21:58:30
 Spectrum: Tune Spec :Average 753-756(8.17-8.19) Bgrd 747(8.13)
 Base Peak: 95.00
 Minimum % Base Peak: 0
 Number of Points: 159

| m/z | Y | m/z | Y | m/z | Y | m/z | Y |
|-------|-------|--------|-------|--------|-------|--------|------|
| 37.00 | 1460 | 91.00 | 135 | 142.00 | 81 | 207.00 | 386 |
| 38.00 | 1185 | 92.00 | 1063 | 143.00 | 399 | 210.00 | 52 |
| 39.00 | 160 | 93.00 | 1372 | 144.00 | 127 | 213.00 | 65 |
| 40.00 | 41 | 94.00 | 4557 | 145.00 | 90 | 215.00 | 58 |
| 41.00 | 260 | 95.00 | 31208 | 146.00 | 95 | 216.00 | 70 |
| 45.00 | 52 | 96.00 | 2385 | 147.00 | 405 | 221.00 | 50 |
| 46.00 | 84 | 97.00 | 133 | 148.00 | 139 | 222.00 | 53 |
| 47.00 | 345 | 99.00 | 63 | 149.00 | 76 | 224.00 | 60 |
| 48.00 | 326 | 100.00 | 159 | 150.00 | 52 | 225.00 | 58 |
| 49.00 | 1164 | 102.00 | 173 | 151.00 | 125 | 228.00 | 85 |
| 50.00 | 4821 | 103.00 | 254 | 153.00 | 56 | 229.00 | 55 |
| 51.00 | 1297 | 104.00 | 154 | 155.00 | 123 | 231.00 | 61 |
| 52.00 | 129 | 105.00 | 65 | 156.00 | 232 | 236.00 | 125 |
| 55.00 | 73 | 106.00 | 62 | 159.00 | 82 | 239.00 | 67 |
| 56.00 | 422 | 107.00 | 167 | 160.00 | 183 | 241.00 | 52 |
| 57.00 | 692 | 109.00 | 69 | 161.00 | 234 | 246.00 | 61 |
| 58.00 | 98 | 110.00 | 57 | 162.00 | 140 | 249.00 | 280 |
| 59.00 | 123 | 112.00 | 165 | 163.00 | 115 | 250.00 | 68 |
| 60.00 | 313 | 115.00 | 61 | 166.00 | 57 | 251.00 | 76 |
| 61.00 | 1029 | 116.00 | 226 | 167.00 | 55 | 252.00 | 64 |
| 62.00 | 1223 | 117.00 | 397 | 169.00 | 106 | 253.00 | 135 |
| 63.00 | 1181 | 118.00 | 175 | 172.00 | 336 | 254.00 | 66 |
| 66.00 | 103 | 119.00 | 259 | 174.00 | 19592 | 259.00 | 110 |
| 67.00 | 7 | 120.00 | 55 | 175.00 | 1512 | 261.00 | 88 |
| 68.00 | 2608 | 123.00 | 98 | 176.00 | 18920 | 264.00 | 58 |
| 69.00 | 3325 | 124.00 | 58 | 177.00 | 1048 | 265.00 | 215 |
| 70.00 | 447 | 125.00 | 99 | 178.00 | 70 | 266.00 | 101 |
| 73.00 | 1556 | 126.00 | 222 | 179.00 | 56 | 267.00 | 106 |
| 74.00 | 5474 | 127.00 | 115 | 180.00 | 64 | 268.00 | 112 |
| 75.00 | 16480 | 129.00 | 121 | 181.00 | 114 | 277.00 | 82 |
| 76.00 | 1605 | 130.00 | 277 | 183.00 | 94 | 281.00 | 2104 |
| 77.00 | 66 | 131.00 | 151 | 185.00 | 61 | 282.00 | 517 |
| 79.00 | 1033 | 132.00 | 64 | 186.00 | 62 | 283.00 | 210 |

Report Date: 18-Oct-2017 20:30:35

Chrom Revision: 2.2 16-Aug-2017 16:24:46

Data File:

\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\6101801D.D\MSVOA_LL_CHHP6.rslt\spec

Injection Date:

17-Oct-2017 21:58:30

Spectrum:

Tune Spec :Average 753-756(8.17-8.19) Bgrd 747(8.13)

Base Peak:

95.00

Minimum % Base Peak: 0

Number of Points: 159

| m/z | Y | m/z | Y | m/z | Y | m/z | Y |
|-------|-----|--------|-----|--------|-----|--------|-----|
| 80.00 | 538 | 133.00 | 187 | 188.00 | 60 | 286.00 | 161 |
| 81.00 | 706 | 134.00 | 243 | 189.00 | 65 | 289.00 | 58 |
| 82.00 | 257 | 135.00 | 117 | 191.00 | 349 | 291.00 | 85 |
| 84.00 | 217 | 137.00 | 79 | 193.00 | 480 | 292.00 | 50 |
| 85.00 | 100 | 138.00 | 101 | 194.00 | 250 | 295.00 | 55 |
| 87.00 | 931 | 139.00 | 95 | 200.00 | 57 | 297.00 | 54 |
| 88.00 | 936 | 141.00 | 361 | 201.00 | 129 | | |

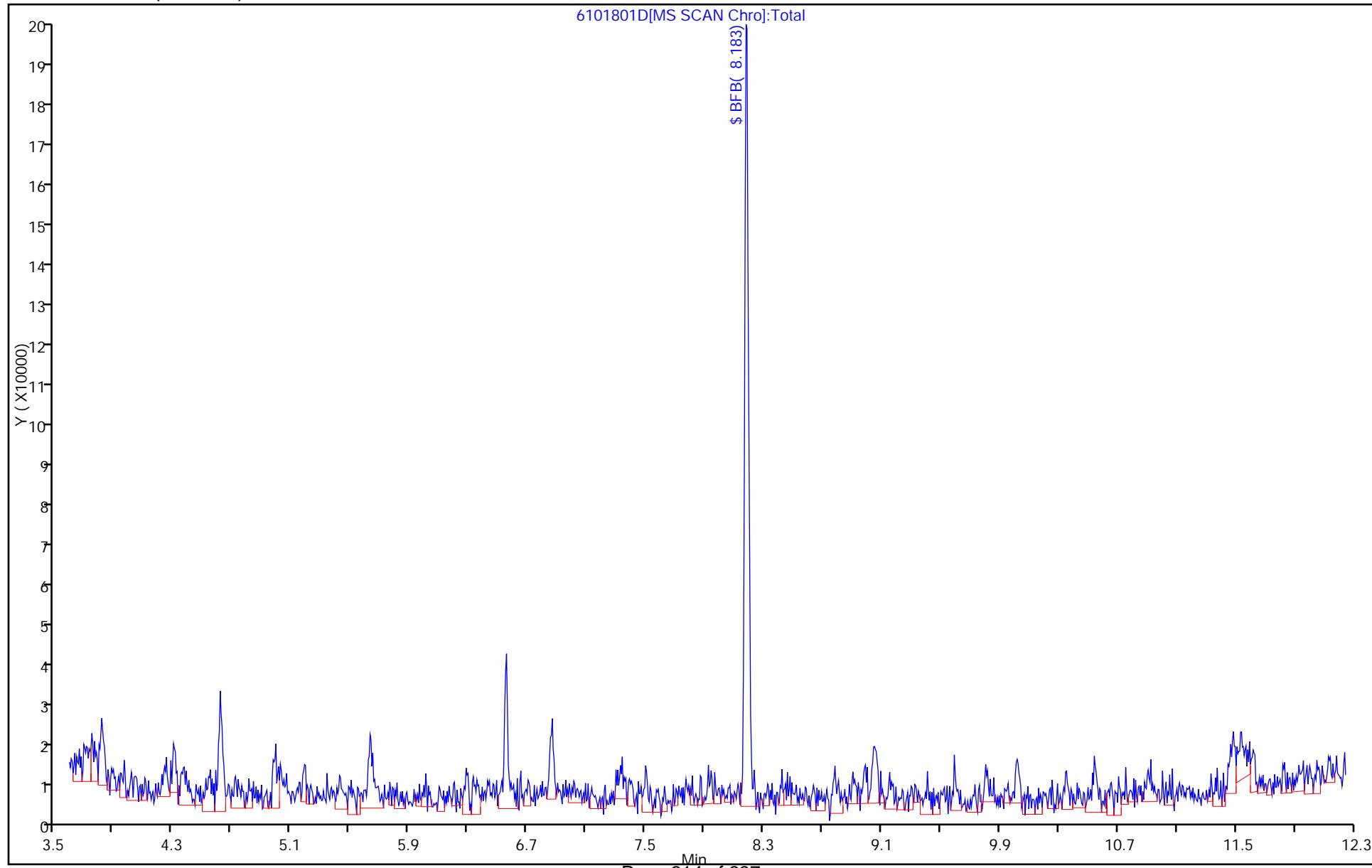
Report Date: 18-Oct-2017 20:30:35

Chrom Revision: 2.2 16-Aug-2017 16:24:46

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101801D.D
Injection Date: 17-Oct-2017 21:58:30 Instrument ID: CHHP6
Lims ID: BFB Operator ID: 034635
Client ID:
Injection Vol: 5.0 mL Dil. Factor: 1.0000 ALS Bottle#: 1
Method: MSVOA_LL_CHHP6 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.: 180-71131-1

SDG No.: _____

Client Sample ID: _____

Lab Sample ID: MB 180-226148/6

Matrix: Water

Lab File ID: 6101806D.D

Analysis Method: 8260C

Date Collected: _____

Sample wt/vol: 5 (mL)

Date Analyzed: 10/18/2017 02:09

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.: 226148

Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|------------|-----------------------------|--------|---|-----|------|
| 74-87-3 | Chloromethane | 1.0 | U | 1.0 | 0.90 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 1.0 | 0.88 |
| 74-83-9 | Bromomethane | 1.0 | U | 1.0 | 0.89 |
| 75-00-3 | Chloroethane | 1.0 | U | 1.0 | 0.90 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.55 |
| 67-64-1 | Acetone | 5.0 | U | 5.0 | 3.4 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 1.0 | 0.88 |
| 75-09-2 | Methylene Chloride | 1.0 | U | 1.0 | 0.36 |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.67 |
| 1634-04-4 | Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.59 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.63 |
| 156-59-2 | cis-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.71 |
| 74-97-5 | Bromochloromethane | 1.0 | U | 1.0 | 0.63 |
| 78-93-3 | 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 |
| 67-66-3 | Chloroform | 1.0 | U | 1.0 | 0.60 |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.60 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 1.0 | 0.88 |
| 71-43-2 | Benzene | 1.0 | U | 1.0 | 0.60 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.57 |
| 79-01-6 | Trichloroethene | 1.0 | U | 1.0 | 0.69 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.66 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 1.0 | 0.64 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.59 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 5.0 | U | 5.0 | 3.1 |
| 108-88-3 | Toluene | 1.0 | U | 1.0 | 0.46 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.58 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.45 |
| 127-18-4 | Tetrachloroethene | 1.0 | U | 1.0 | 0.47 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 5.0 | 3.3 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 1.0 | 0.84 |
| 106-93-4 | 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.50 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 1.0 | 0.50 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.57 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 1.0 | 0.51 |
| 1330-20-7 | Xylenes, Total | 2.0 | U | 2.0 | 0.89 |
| 100-42-5 | Styrene | 1.0 | U | 1.0 | 0.47 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.: _____

Client Sample ID: _____

Lab Sample ID: MB 180-226148/6

Matrix: Water

Lab File ID: 6101806.D.D

Analysis Method: 8260C

Date Collected: _____

Sample wt/vol: 5 (mL)

Date Analyzed: 10/18/2017 02:09

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.: 226148

Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|----------|---------------------------|--------|---|-----|------|
| 75-25-2 | Bromoform | 1.0 | U | 1.0 | 0.98 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.60 |
| 107-13-1 | Acrylonitrile | 20 | U | 20 | 7.8 |
| 123-91-1 | 1,4-Dioxane | 200 | U | 200 | 14 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|------------|------------------------------|------|---|--------|
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 93 | | 65-121 |
| 2037-26-5 | Toluene-d8 (Surr) | 89 | | 73-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 90 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 91 | | 73-120 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\6101806.D.D
 Lims ID: MB
 Client ID:
 Sample Type: MB
 Inject. Date: 18-Oct-2017 02:09:30 ALS Bottle#: 6 Worklist Smp#: 6
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018914-006
 Misc. Info.: MB
 Operator ID: 034635 Instrument ID: CHHP6
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 18-Oct-2017 20:30:42 Calib Date: 28-Sep-2017 15:13:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICAL File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170928-18631.b\60928P06.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK031

First Level Reviewer: bungardf Date: 18-Oct-2017 02:37:50

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 3.955 | 3.949 | 0.006 | 92 | 312973 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.045 | 7.045 | 0.000 | 98 | 1062992 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.172 | 10.166 | 0.006 | 88 | 286612 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.508 | 12.508 | 0.000 | 97 | 425493 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr) | 113 | 6.315 | 6.314 | 0.001 | 91 | 251413 | 50.0 | 45.5 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur) | 65 | 6.686 | 6.685 | 0.001 | 70 | 368853 | 50.0 | 46.7 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.712 | 8.711 | 0.001 | 93 | 1035065 | 50.0 | 44.7 | |
| \$ 8 4-Bromofluorobenzene (Surr) | 95 | 11.352 | 11.351 | 0.001 | 83 | 443998 | 50.0 | 44.9 | |
| 11 Dichlorodifluoromethane | 85 | | 1.472 | | | | | ND | |
| 12 Chloromethane | 50 | | 1.624 | | | | | ND | |
| 13 Vinyl chloride | 62 | | 1.740 | | | | | ND | |
| 14 Butadiene | 39 | | 1.788 | | | | | ND | |
| 15 Bromomethane | 94 | | 2.068 | | | | | ND | |
| 16 Chloroethane | 64 | | 2.202 | | | | | ND | |
| 17 Dichlorofluoromethane | 67 | | 2.457 | | | | | ND | |
| 18 Trichlorofluoromethane | 101 | | 2.470 | | | | | ND | |
| 20 Ethyl ether | 59 | | 2.816 | | | | | ND | |
| 19 Ethanol | 45 | | 2.823 | | | | | ND | |
| 21 Acrolein | 56 | | 2.981 | | | | | ND | |
| 22 1,1-Dichloroethene | 96 | | 3.090 | | | | | ND | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | | 3.151 | | | | | ND | |
| 24 Acetone | 43 | | 3.181 | | | | | ND | |
| 25 Iodomethane | 142 | | 3.273 | | | | | ND | |
| 26 Carbon disulfide | 76 | | 3.364 | | | | | ND | |
| 27 Isopropyl alcohol | 45 | | 3.432 | | | | | ND | |
| 28 Acetonitrile | 41 | | 3.572 | | | | | ND | |
| 29 3-Chloro-1-propene | 76 | | 3.631 | | | | | ND | |
| 30 Methyl acetate | 43 | | 3.644 | | | | | ND | |
| 31 Methylene Chloride | 84 | 3.833 | 3.838 | -0.005 | 83 | 8391 | | 1.15 | |
| 32 2-Methyl-2-propanol | 59 | | 4.094 | | | | | ND | |
| 33 Acrylonitrile | 53 | | 4.222 | | | | | 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 | | 4.270 | | | | | ND | |
| 35 Methyl tert-butyl ether | 73 | | 4.276 | | | | | ND | |
| 36 Hexane | 57 | | 4.696 | | | | | ND | |
| 37 1,1-Dichloroethane | 63 | | 4.921 | | | | | ND | |
| 38 Vinyl acetate | 43 | | 4.976 | | | | | ND | |
| 40 Isopropyl ether | 45 | | 5.020 | | | | | ND | |
| 39 2-Chloro-1,3-butadiene | 53 | | 5.020 | | | | | ND | |
| 41 Tert-butyl ethyl ether | 59 | | 5.506 | | | | | ND | |
| 42 2,2-Dichloropropane | 97 | | 5.676 | | | | | ND | |
| 43 cis-1,2-Dichloroethene | 96 | | 5.688 | | | | | ND | |
| 44 2-Butanone (MEK) | 43 | | 5.694 | | | | | ND | |
| 45 Propionitrile | 54 | | 5.768 | | | | | ND | |
| 46 Ethyl acetate | 43 | | 5.774 | | | | | ND | |
| 47 Methacrylonitrile | 41 | | 5.950 | | | | | ND | |
| 48 Chlorobromomethane | 128 | | 5.974 | | | | | ND | |
| 49 Tetrahydrofuran | 42 | | 5.992 | | | | | ND | |
| 50 Chloroform | 83 | 6.120 | 6.126 | -0.006 | 53 | 5787 | | 0.5262 | |
| 51 1,1,1-Trichloroethane | 97 | | 6.290 | | | | | ND | |
| 52 Cyclohexane | 56 | | 6.357 | | | | | ND | |
| 53 Carbon tetrachloride | 117 | | 6.454 | | | | | ND | |
| 54 1,1-Dichloropropene | 75 | | 6.479 | | | | | ND | |
| 56 Benzene | 78 | | 6.691 | | | | | ND | |
| 55 Isobutyl alcohol | 41 | | 6.691 | | | | | ND | |
| 57 1,2-Dichloroethane | 62 | | 6.777 | | | | | ND | |
| 148 Isooctane | 57 | | 6.857 | | | | | ND | |
| 58 Tert-amyl methyl ether | 73 | | 6.881 | | | | | ND | |
| 59 n-Heptane | 43 | | 7.069 | | | | | ND | |
| 60 n-Butanol | 56 | | 7.410 | | | | | ND | |
| 61 Trichloroethene | 130 | | 7.440 | | | | | ND | |
| 62 Ethyl acrylate | 55 | | 7.569 | | | | | ND | |
| 63 Methylcyclohexane | 83 | | 7.671 | | | | | ND | |
| 64 1,2-Dichloropropane | 63 | | 7.713 | | | | | ND | |
| 65 1,4-Dioxane | 88 | | 7.793 | | | | | ND | |
| 67 Dibromomethane | 93 | | 7.799 | | | | | ND | |
| 66 Methyl methacrylate | 69 | | 7.806 | | | | | ND | |
| 68 Dichlorobromomethane | 83 | | 7.999 | | | | | ND | |
| 70 2-Chloroethyl vinyl ether | 63 | | 8.310 | | | | | ND | |
| 71 cis-1,3-Dichloropropene | 75 | | 8.450 | | | | | ND | |
| 72 4-Methyl-2-pentanone (MIBK) | 43 | | 8.608 | | | | | ND | |
| 73 Toluene | 91 | | 8.778 | | | | | ND | |
| 74 trans-1,3-Dichloropropene | 75 | | 9.034 | | | | | ND | |
| 75 Ethyl methacrylate | 69 | | 9.100 | | | | | ND | |
| 76 1,1,2-Trichloroethane | 97 | | 9.222 | | | | | ND | |
| 77 Tetrachloroethene | 164 | | 9.295 | | | | | ND | |
| 78 1,3-Dichloropropane | 76 | | 9.380 | | | | | ND | |
| 79 2-Hexanone | 43 | | 9.447 | | | | | ND | |
| 80 n-Butyl acetate | 43 | | 9.576 | | | | | ND | |
| 81 Chlorodibromomethane | 129 | | 9.593 | | | | | ND | |
| 82 Ethylene Dibromide | 107 | | 9.709 | | | | | ND | |
| 83 3-Chlorobenzotrifluoride | 180 | | 10.183 | | | | | ND | |
| 84 Chlorobenzene | 112 | | 10.196 | | | | | ND | |
| 85 4-Chlorobenzotrifluoride | 180 | | 10.269 | | | | | ND | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|--------------|------------------|------------------|---|----------|---------------|-----------------|-------|
| 86 1,1,1,2-Tetrachloroethane | 131 | | 10.293 | | | | | ND | |
| 87 Ethylbenzene | 106 | | 10.299 | | | | | ND | |
| 88 m-Xylene & p-Xylene | 106 | | 10.433 | | | | | ND | |
| 89 o-Xylene | 106 | | 10.816 | | | | | ND | |
| 90 Styrene | 104 | | 10.834 | | | | | ND | |
| 91 Bromoform | 173 | | 11.011 | | | | | ND | |
| 129 Cyclohexanol | 57 | | 11.037 | | | | | ND | |
| 92 2-Chlorobenzotrifluoride | 180 | | 11.084 | | | | | ND | |
| 93 Isopropylbenzene | 105 | | 11.181 | | | | | ND | |
| 94 Cyclohexanone | 55 | | 11.267 | | | | | ND | |
| 95 Bromobenzene | 156 | | 11.485 | | | | | ND | |
| 96 1,1,2,2-Tetrachloroethane | 83 | | 11.497 | | | | | ND | |
| 97 trans-1,4-Dichloro-2-butene | 53 | | 11.540 | | | | | ND | |
| 98 1,2,3-Trichloropropane | 110 | | 11.546 | | | | | ND | |
| 99 N-Propylbenzene | 120 | | 11.595 | | | | | ND | |
| 100 2-Chlorotoluene | 126 | | 11.680 | | | | | ND | |
| 101 3-Chlorotoluene | 126 | | 11.747 | | | | | ND | |
| 102 1,3,5-Trimethylbenzene | 105 | | 11.783 | | | | | ND | |
| 103 4-Chlorotoluene | 126 | | 11.802 | | | | | ND | |
| 104 tert-Butylbenzene | 119 | | 12.094 | | | | | ND | |
| 106 1,2,4-Trimethylbenzene | 105 | | 12.154 | | | | | ND | |
| 107 1,2-dichloro-4-(trifluorom | 214 | | 12.203 | | | | | ND | |
| 108 sec-Butylbenzene | 105 | | 12.319 | | | | | ND | |
| 109 1,3-Dichlorobenzene | 146 | | 12.428 | | | | | ND | |
| 110 4-Isopropyltoluene | 119 | | 12.471 | | | | | ND | |
| 111 1,4-Dichlorobenzene | 146 | | 12.532 | | | | | ND | |
| 112 1,2,3-Trimethylbenzene | 105 | | 12.563 | | | | | ND | |
| 113 2,4-Dichloro-1-(trifluorom | 214 | | 12.568 | | | | | ND | |
| 114 2,5-Dichlorobenzotrifluori | 214 | | 12.611 | | | | | ND | |
| 115 Benzyl chloride | 91 | | 12.648 | | | | | ND | |
| 116 n-Butylbenzene | 91 | | 12.878 | | | | | ND | |
| 117 1,2-Dichlorobenzene | 146 | | 12.890 | | | | | ND | |
| 118 1,2-Dibromo-3-Chloropropan | 75 | | 13.675 | | | | | ND | |
| 119 2,4- & 2,5- & 2,6- Dichlor | 125 | | 13.815 | | | | | ND | |
| 120 1,3,5-Trichlorobenzene | 180 | | 13.865 | | | | | ND | |
| 121 2,3- & 3,4- Dichlorotoluen | 125 | | 14.235 | | | | | ND | |
| 122 1,2,4-Trichlorobenzene | 180 | | 14.497 | | | | | ND | |
| 123 Hexachlorobutadiene | 225 | | 14.643 | | | | | ND | |
| 124 Naphthalene | 128 | | 14.758 | | | | | ND | |
| 125 1,2,3-Trichlorobenzene | 180 | | 14.977 | | | | | ND | |
| 126 2,4,5-Trichlorotoluene | 159 | | 15.774 | | | | | ND | |
| 127 2,3,6-Trichlorotoluene | 159 | | 15.877 | | | | | ND | |
| 150 Tert-butyl ethyl ether (TI) | 1 | | 0.000 | | | | | ND | |
| 153 1,2 Epoxybutane TIC | 1 | | 0.000 | | | | | ND | |
| 151 Tert-amyl methyl ether (TI) | 1 | | 0.000 | | | | | ND | |
| 146 3,4-Dichlorotoluene | 1 | | 0.000 | | | | | ND | |
| 152 Formaldehyde TIC | 1 | | 0.000 | | | | | ND | |
| 149 Isopropyl ether TIC | 1 | | 0.000 | | | | | ND | |
| S 131 Xylenes, Total | 106 | | 1.000 | | | | | ND | |
| S 130 1,2-Dichloroethene, Total | 96 | | 1.000 | | | | | ND | |
| S 132 1,3-Dichloropropene, Total | 1 | | 0.000 | | | | | ND | |
| S 154 Total BTEX | 1 | | 0.000 | | | | | ND | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|--------------|------------------|------------------|---|----------|---------------|-----------------|-------|
| T 133 Tetrahydrofuran TIC | 42 | | 0.000 | | | | | | ND |
| T 135 Mesityl oxide TIC | 83 | | 0.000 | | | | | | ND |
| T 134 Methyl n-amyl ketone TIC | 43 | | 0.000 | | | | | | ND |

Reagents:

VOA8260INT_00074

Amount Added: 2.00

Units: uL

Run Reagent

VOA8260SURR_00073

Amount Added: 2.00

Units: uL

Run Reagent

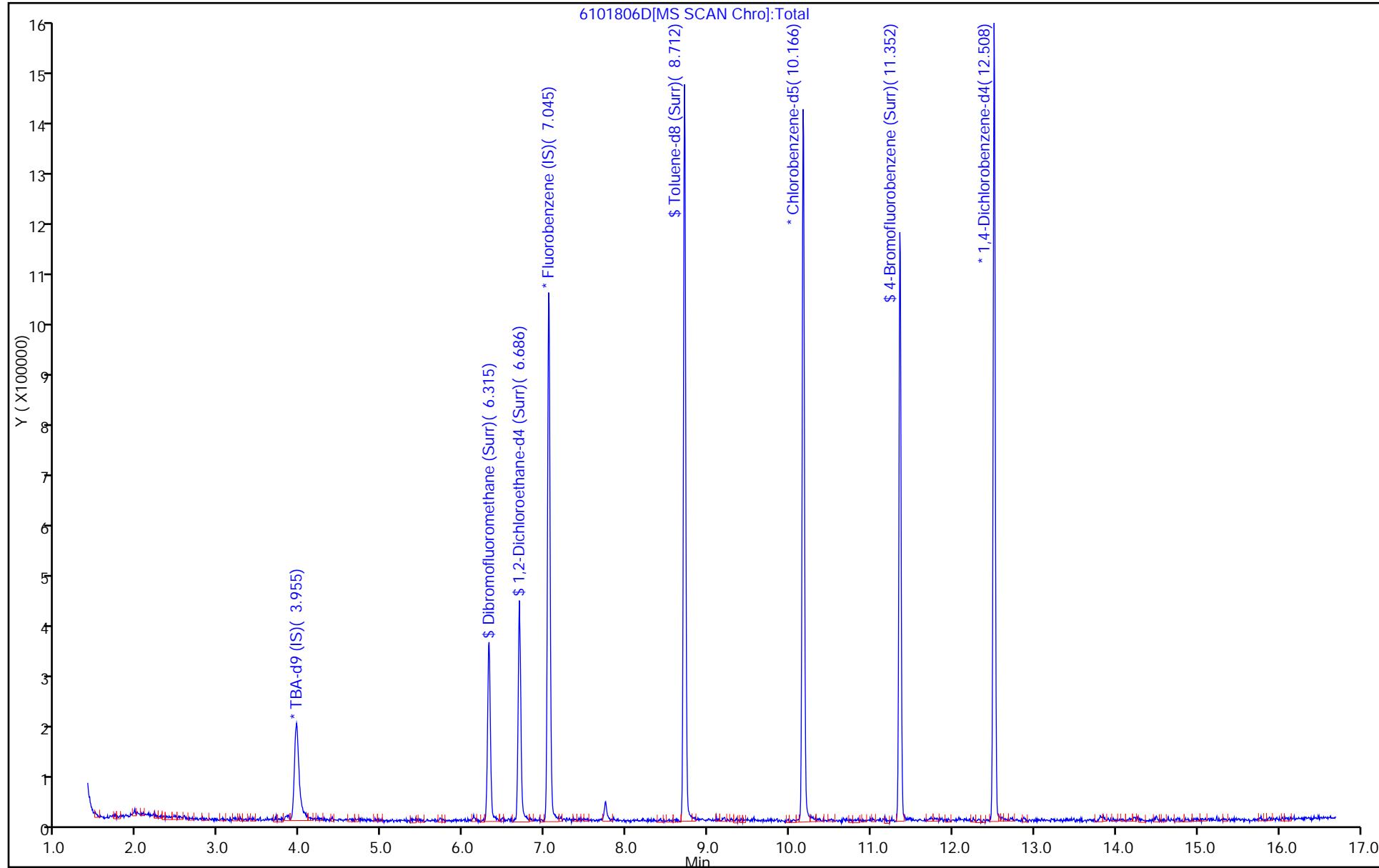
Report Date: 18-Oct-2017 20:30:55

Chrom Revision: 2.2 16-Aug-2017 16:24:46

TestAmerica Pittsburgh

Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101806D.D
Injection Date: 18-Oct-2017 02:09:30 Instrument ID: CHHP6
Lims ID: MB Operator ID: 034635
Client ID:
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 6
Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm)

Worklist Smp#: 6



TestAmerica Pittsburgh
Recovery Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\6101806.D.D
 Lims ID: MB
 Client ID:
 Sample Type: MB
 Inject. Date: 18-Oct-2017 02:09:30 ALS Bottle#: 6 Worklist Smp#: 6
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018914-006
 Misc. Info.: MB
 Operator ID: 034635 Instrument ID: CHHP6
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 18-Oct-2017 20:30:42 Calib Date: 28-Sep-2017 15:13:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170928-18631.b\60928P06.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK031

First Level Reviewer: bungardf Date: 18-Oct-2017 02:37:50

| Compound | Amount Added | Amount Recovered | % Rec. |
|-----------------------------------|--------------|------------------|--------|
| \$ 5 Dibromofluoromethane (Surr) | 50.0 | 45.5 | 91.03 |
| \$ 6 1,2-Dichloroethane-d4 (Surr) | 50.0 | 46.7 | 93.44 |
| \$ 7 Toluene-d8 (Surr) | 50.0 | 44.7 | 89.44 |
| \$ 8 4-Bromofluorobenzene (Surr) | 50.0 | 44.9 | 89.73 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.: _____

Client Sample ID: _____

Lab Sample ID: LCS 180-226148/4

Matrix: Water

Lab File ID: 6101804D.D

Analysis Method: 8260C

Date Collected: _____

Sample wt/vol: 5 (mL)

Date Analyzed: 10/18/2017 01:04

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.: 226148

Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|------------|-----------------------------|--------|---|-----|------|
| 74-87-3 | Chloromethane | 7.89 | | 1.0 | 0.90 |
| 75-01-4 | Vinyl chloride | 8.87 | | 1.0 | 0.88 |
| 74-83-9 | Bromomethane | 7.90 | | 1.0 | 0.89 |
| 75-00-3 | Chloroethane | 9.09 | | 1.0 | 0.90 |
| 75-35-4 | 1,1-Dichloroethene | 9.66 | | 1.0 | 0.55 |
| 67-64-1 | Acetone | 22.4 | | 5.0 | 3.4 |
| 75-15-0 | Carbon disulfide | 8.89 | | 1.0 | 0.88 |
| 75-09-2 | Methylene Chloride | 8.98 | | 1.0 | 0.36 |
| 156-60-5 | trans-1,2-Dichloroethene | 9.35 | | 1.0 | 0.67 |
| 1634-04-4 | Methyl tert-butyl ether | 8.28 | | 1.0 | 0.59 |
| 75-34-3 | 1,1-Dichloroethane | 9.00 | | 1.0 | 0.63 |
| 156-59-2 | cis-1,2-Dichloroethene | 9.03 | | 1.0 | 0.71 |
| 74-97-5 | Bromochloromethane | 9.10 | | 1.0 | 0.63 |
| 78-93-3 | 2-Butanone (MEK) | 18.3 | | 5.0 | 2.6 |
| 67-66-3 | Chloroform | 9.63 | | 1.0 | 0.60 |
| 71-55-6 | 1,1,1-Trichloroethane | 10.4 | | 1.0 | 0.60 |
| 56-23-5 | Carbon tetrachloride | 12.2 | | 1.0 | 0.88 |
| 71-43-2 | Benzene | 9.47 | | 1.0 | 0.60 |
| 107-06-2 | 1,2-Dichloroethane | 9.07 | | 1.0 | 0.57 |
| 79-01-6 | Trichloroethene | 9.21 | | 1.0 | 0.69 |
| 78-87-5 | 1,2-Dichloropropane | 8.45 | | 1.0 | 0.66 |
| 75-27-4 | Bromodichloromethane | 9.26 | | 1.0 | 0.64 |
| 10061-01-5 | cis-1,3-Dichloropropene | 9.17 | | 1.0 | 0.59 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 13.2 | | 5.0 | 3.1 |
| 108-88-3 | Toluene | 9.31 | | 1.0 | 0.46 |
| 10061-02-6 | trans-1,3-Dichloropropene | 9.36 | | 1.0 | 0.58 |
| 79-00-5 | 1,1,2-Trichloroethane | 8.74 | | 1.0 | 0.45 |
| 127-18-4 | Tetrachloroethene | 9.13 | | 1.0 | 0.47 |
| 591-78-6 | 2-Hexanone | 18.9 | | 5.0 | 3.3 |
| 124-48-1 | Dibromochloromethane | 9.51 | | 1.0 | 0.84 |
| 106-93-4 | 1,2-Dibromoethane (EDB) | 8.97 | | 1.0 | 0.50 |
| 108-90-7 | Chlorobenzene | 9.44 | | 1.0 | 0.50 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 10.1 | | 1.0 | 0.57 |
| 100-41-4 | Ethylbenzene | 9.18 | | 1.0 | 0.51 |
| 1330-20-7 | Xylenes, Total | 18.2 | | 2.0 | 0.89 |
| 100-42-5 | Styrene | 9.53 | | 1.0 | 0.47 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.: _____

Client Sample ID: _____

Lab Sample ID: LCS 180-226148/4

Matrix: Water

Lab File ID: 6101804D.D

Analysis Method: 8260C

Date Collected: _____

Sample wt/vol: 5 (mL)

Date Analyzed: 10/18/2017 01:04

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.: 226148

Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|----------|---------------------------|--------|---|-----|------|
| 75-25-2 | Bromoform | 8.29 | | 1.0 | 0.98 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 8.83 | | 1.0 | 0.60 |
| 107-13-1 | Acrylonitrile | 65.5 | | 20 | 7.8 |
| 123-91-1 | 1,4-Dioxane | 144 | J | 200 | 14 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|------------|------------------------------|------|---|--------|
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 92 | | 65-121 |
| 2037-26-5 | Toluene-d8 (Surr) | 93 | | 73-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 94 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 92 | | 73-120 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\6101804.D.D
 Lims ID: LCS
 Client ID:
 Sample Type: LCS
 Inject. Date: 18-Oct-2017 01:04:30 ALS Bottle#: 4 Worklist Smp#: 4
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018914-004
 Misc. Info.: LCS
 Operator ID: 034635 Instrument ID: CHHP6
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 18-Oct-2017 20:30:42 Calib Date: 28-Sep-2017 15:13:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICAL File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170928-18631.b\60928P06.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK031

First Level Reviewer: bungardf Date: 18-Oct-2017 01:25:42

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 3.957 | 3.949 | 0.008 | 97 | 373208 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.047 | 7.045 | 0.002 | 98 | 1246060 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.168 | 10.166 | 0.002 | 87 | 301504 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.510 | 12.508 | 0.002 | 96 | 469268 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr) | 113 | 6.311 | 6.314 | -0.003 | 92 | 298392 | 50.0 | 46.1 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur) | 65 | 6.682 | 6.685 | -0.003 | 76 | 423400 | 50.0 | 45.8 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.714 | 8.711 | 0.003 | 93 | 1128430 | 50.0 | 46.5 | |
| \$ 8 4-Bromofluorobenzene (Surr) | 95 | 11.355 | 11.351 | 0.003 | 82 | 487330 | 50.0 | 46.8 | |
| 11 Dichlorodifluoromethane | 85 | 1.469 | 1.472 | -0.003 | 99 | 389857 | 50.0 | 49.0 | |
| 12 Chloromethane | 50 | 1.627 | 1.624 | 0.003 | 99 | 281976 | 50.0 | 39.4 | |
| 13 Vinyl chloride | 62 | 1.743 | 1.740 | 0.003 | 97 | 332607 | 50.0 | 44.4 | |
| 14 Butadiene | 39 | 1.785 | 1.788 | -0.003 | 91 | 272714 | 50.0 | 43.9 | |
| 15 Bromomethane | 94 | 2.077 | 2.068 | 0.009 | 92 | 138063 | 50.0 | 39.5 | |
| 16 Chloroethane | 64 | 2.205 | 2.202 | 0.003 | 98 | 184081 | 50.0 | 45.5 | |
| 17 Dichlorofluoromethane | 67 | 2.467 | 2.457 | 0.010 | 95 | 396925 | 50.0 | 45.5 | |
| 18 Trichlorofluoromethane | 101 | 2.467 | 2.470 | -0.003 | 77 | 408053 | 50.0 | 55.3 | |
| 20 Ethyl ether | 59 | 2.819 | 2.816 | 0.003 | 84 | 284888 | 50.0 | 45.1 | |
| 21 Acrolein | 56 | 2.984 | 2.981 | 0.003 | 97 | 168069 | 150.0 | 124.2 | |
| 22 1,1-Dichloroethene | 96 | 3.099 | 3.090 | 0.009 | 98 | 312717 | 50.0 | 48.3 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.154 | 3.151 | 0.003 | 95 | 344025 | 50.0 | 55.4 | |
| 24 Acetone | 43 | 3.184 | 3.181 | 0.003 | 100 | 300023 | 100.0 | 112.2 | |
| 25 Iodomethane | 142 | 3.282 | 3.273 | 0.009 | 99 | 415431 | 50.0 | 45.6 | |
| 26 Carbon disulfide | 76 | 3.361 | 3.364 | -0.003 | 98 | 640921 | 50.0 | 44.4 | |
| 29 3-Chloro-1-propene | 76 | 3.629 | 3.631 | -0.003 | 82 | 167178 | 50.0 | 44.1 | |
| 30 Methyl acetate | 43 | 3.641 | 3.644 | -0.003 | 95 | 399082 | 100.0 | 68.1 | |
| 31 Methylene Chloride | 84 | 3.835 | 3.838 | -0.003 | 80 | 385298 | 50.0 | 44.9 | |
| 32 2-Methyl-2-propanol | 59 | 4.097 | 4.094 | 0.003 | 91 | 197945 | 500.0 | 479.2 | |
| 33 Acrylonitrile | 53 | 4.225 | 4.222 | 0.003 | 98 | 1024713 | 500.0 | 327.4 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.273 | 4.270 | 0.003 | 97 | 343681 | 50.0 | 46.8 | |
| 35 Methyl tert-butyl ether | 73 | 4.286 | 4.276 | 0.010 | 94 | 972508 | 50.0 | 41.4 | |
| 36 Hexane | 57 | 4.705 | 4.696 | 0.009 | 90 | 365510 | 50.0 | 42.0 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|--------------|------------------|------------------|-----|----------|---------------|-----------------|-------|
| 37 1,1-Dichloroethane | 63 | 4.924 | 4.921 | 0.003 | 95 | 537980 | 50.0 | 45.0 | |
| 38 Vinyl acetate | 43 | 4.973 | 4.976 | -0.003 | 97 | 439271 | 50.0 | 31.6 | |
| 42 2,2-Dichloropropane | 97 | 5.685 | 5.676 | 0.009 | 69 | 67110 | 50.0 | 54.2 | |
| 43 cis-1,2-Dichloroethene | 96 | 5.691 | 5.688 | 0.003 | 88 | 387487 | 50.0 | 45.1 | |
| 44 2-Butanone (MEK) | 43 | 5.703 | 5.694 | 0.009 | 95 | 350577 | 100.0 | 91.5 | |
| 48 Chlorobromomethane | 128 | 5.977 | 5.974 | 0.003 | 87 | 170416 | 50.0 | 45.5 | |
| 49 Tetrahydrofuran | 42 | 5.983 | 5.992 | -0.009 | 79 | 144858 | 100.0 | 55.8 | |
| 50 Chloroform | 83 | 6.129 | 6.126 | 0.003 | 92 | 620509 | 50.0 | 48.1 | |
| 51 1,1,1-Trichloroethane | 97 | 6.287 | 6.290 | -0.003 | 98 | 424391 | 50.0 | 51.8 | |
| 52 Cyclohexane | 56 | 6.354 | 6.357 | -0.003 | 83 | 472230 | 50.0 | 40.6 | |
| 53 Carbon tetrachloride | 117 | 6.457 | 6.454 | 0.003 | 96 | 358465 | 50.0 | 60.8 | |
| 54 1,1-Dichloropropene | 75 | 6.476 | 6.479 | -0.003 | 98 | 486185 | 50.0 | 49.3 | |
| 56 Benzene | 78 | 6.695 | 6.691 | 0.004 | 97 | 1359299 | 50.0 | 47.3 | |
| 55 Isobutyl alcohol | 41 | 6.695 | 6.691 | 0.004 | 42 | 173150 | 1250.0 | 1001.8 | |
| 57 1,2-Dichloroethane | 62 | 6.774 | 6.777 | -0.003 | 98 | 481251 | 50.0 | 45.4 | |
| 59 n-Heptane | 43 | 7.072 | 7.069 | 0.003 | 83 | 255785 | 50.0 | 38.6 | |
| 61 Trichloroethene | 130 | 7.437 | 7.440 | -0.003 | 95 | 321208 | 50.0 | 46.0 | |
| 63 Methylcyclohexane | 83 | 7.674 | 7.671 | 0.003 | 84 | 569359 | 50.0 | 46.7 | |
| 64 1,2-Dichloropropane | 63 | 7.717 | 7.713 | 0.004 | 92 | 303214 | 50.0 | 42.2 | |
| 65 1,4-Dioxane | 88 | 7.796 | 7.793 | 0.003 | 38 | 47150 | 1000.0 | 720.1 | |
| 67 Dibromomethane | 93 | 7.802 | 7.799 | 0.003 | 91 | 220429 | 50.0 | 46.9 | |
| 68 Dichlorobromomethane | 83 | 8.003 | 7.999 | 0.004 | 99 | 370300 | 50.0 | 46.3 | |
| 70 2-Chloroethyl vinyl ether | 63 | 8.313 | 8.310 | 0.003 | 93 | 414717 | 100.0 | 87.2 | |
| 71 cis-1,3-Dichloropropene | 75 | 8.453 | 8.450 | 0.003 | 96 | 415674 | 50.0 | 45.8 | |
| 72 4-Methyl-2-pentanone (MIBK) | 43 | 8.611 | 8.608 | 0.003 | 92 | 508907 | 100.0 | 66.1 | |
| 73 Toluene | 91 | 8.781 | 8.778 | 0.003 | 99 | 1382578 | 50.0 | 46.5 | |
| 74 trans-1,3-Dichloropropene | 75 | 9.037 | 9.034 | 0.003 | 92 | 372525 | 50.0 | 46.8 | |
| 75 Ethyl methacrylate | 69 | 9.104 | 9.100 | 0.004 | 85 | 438507 | 50.0 | 43.7 | |
| 76 1,1,2-Trichloroethane | 97 | 9.231 | 9.222 | 0.009 | 92 | 300412 | 50.0 | 43.7 | |
| 77 Tetrachloroethene | 164 | 9.298 | 9.295 | 0.003 | 94 | 242848 | 50.0 | 45.7 | |
| 78 1,3-Dichloropropane | 76 | 9.383 | 9.380 | 0.003 | 87 | 569886 | 50.0 | 45.7 | |
| 79 2-Hexanone | 43 | 9.450 | 9.447 | 0.003 | 94 | 467979 | 100.0 | 94.4 | |
| 81 Chlorodibromomethane | 129 | 9.602 | 9.593 | 0.009 | 89 | 221436 | 50.0 | 47.5 | |
| 82 Ethylene Dibromide | 107 | 9.706 | 9.709 | -0.003 | 100 | 296761 | 50.0 | 44.9 | |
| 83 3-Chlorobenzotrifluoride | 180 | 10.180 | 10.183 | -0.003 | 90 | 489724 | 50.0 | 56.7 | |
| 84 Chlorobenzene | 112 | 10.199 | 10.196 | 0.003 | 93 | 911982 | 50.0 | 47.2 | |
| 85 4-Chlorobenzotrifluoride | 180 | 10.266 | 10.269 | -0.003 | 96 | 464828 | 50.0 | 58.1 | |
| 86 1,1,1,2-Tetrachloroethane | 131 | 10.296 | 10.293 | 0.003 | 89 | 277135 | 50.0 | 50.4 | |
| 87 Ethylbenzene | 106 | 10.302 | 10.299 | 0.003 | 98 | 507318 | 50.0 | 45.9 | |
| 88 m-Xylene & p-Xylene | 106 | 10.430 | 10.433 | -0.003 | 99 | 627395 | 50.0 | 46.3 | |
| 89 o-Xylene | 106 | 10.813 | 10.816 | -0.003 | 96 | 609176 | 50.0 | 44.9 | |
| 90 Styrene | 104 | 10.831 | 10.834 | -0.003 | 93 | 1046570 | 50.0 | 47.7 | |
| 91 Bromoform | 173 | 11.014 | 11.011 | 0.003 | 90 | 108624 | 50.0 | 41.4 | |
| 92 2-Chlorobenzotrifluoride | 180 | 11.087 | 11.084 | 0.003 | 91 | 492243 | 50.0 | 56.2 | |
| 93 Isopropylbenzene | 105 | 11.184 | 11.181 | 0.003 | 97 | 1450706 | 50.0 | 47.7 | |
| 95 Bromobenzene | 156 | 11.488 | 11.485 | 0.003 | 95 | 373442 | 50.0 | 42.8 | |
| 96 1,1,2,2-Tetrachloroethane | 83 | 11.494 | 11.497 | -0.003 | 95 | 427212 | 50.0 | 44.2 | |
| 97 trans-1,4-Dichloro-2-butene | 53 | 11.531 | 11.540 | -0.009 | 76 | 90871 | 50.0 | 35.6 | |
| 98 1,2,3-Trichloropropane | 110 | 11.549 | 11.546 | 0.003 | 85 | 150605 | 50.0 | 41.6 | |
| 99 N-Propylbenzene | 120 | 11.598 | 11.595 | 0.003 | 98 | 428743 | 50.0 | 45.2 | |
| 100 2-Chlorotoluene | 126 | 11.683 | 11.680 | 0.003 | 95 | 362595 | 50.0 | 44.1 | |
| 101 3-Chlorotoluene | 126 | 11.750 | 11.747 | 0.003 | 96 | 463181 | 50.0 | 53.7 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|--------------|------------------|------------------|----|----------|---------------|-----------------|-------|
| 102 1,3,5-Trimethylbenzene | 105 | 11.780 | 11.783 | -0.003 | 93 | 1273063 | 50.0 | 47.2 | |
| 103 4-Chlorotoluene | 126 | 11.805 | 11.802 | 0.003 | 99 | 406783 | 50.0 | 45.3 | |
| 104 tert-Butylbenzene | 119 | 12.097 | 12.094 | 0.003 | 91 | 1005579 | 50.0 | 47.0 | |
| 106 1,2,4-Trimethylbenzene | 105 | 12.151 | 12.154 | -0.003 | 98 | 1325719 | 50.0 | 47.2 | |
| 107 1,2-dichloro-4-(trifluorom | 214 | 12.200 | 12.203 | -0.003 | 94 | 347599 | 50.0 | 55.7 | |
| 108 sec-Butylbenzene | 105 | 12.316 | 12.319 | -0.003 | 96 | 1437631 | 50.0 | 47.9 | |
| 109 1,3-Dichlorobenzene | 146 | 12.431 | 12.428 | 0.003 | 94 | 710798 | 50.0 | 45.4 | |
| 110 4-Isopropyltoluene | 119 | 12.474 | 12.471 | 0.003 | 95 | 1224129 | 50.0 | 49.4 | |
| 111 1,4-Dichlorobenzene | 146 | 12.535 | 12.532 | 0.003 | 92 | 733710 | 50.0 | 45.3 | |
| 113 2,4-Dichloro-1-(trifluorom | 214 | 12.565 | 12.568 | -0.003 | 93 | 338043 | 50.0 | 57.5 | |
| 114 2,5-Dichlorobenzotrifluori | 214 | 12.608 | 12.611 | -0.003 | 95 | 371128 | 50.0 | 57.2 | |
| 116 n-Butylbenzene | 91 | 12.881 | 12.878 | 0.003 | 97 | 1105430 | 50.0 | 48.9 | |
| 117 1,2-Dichlorobenzene | 146 | 12.888 | 12.890 | -0.002 | 93 | 683664 | 50.0 | 45.9 | |
| 118 1,2-Dibromo-3-Chloropropan | 75 | 13.678 | 13.675 | 0.003 | 77 | 54373 | 50.0 | 40.9 | |
| 119 2,4- & 2,5- & 2,6- Dichlor | 125 | 13.818 | 13.815 | 0.003 | 99 | 1803892 | 150.0 | 190.1 | |
| 121 2,3- & 3,4- Dichlorotoluen | 125 | 14.232 | 14.235 | -0.003 | 99 | 1317350 | 100.0 | 127.1 | |
| 122 1,2,4-Trichlorobenzene | 180 | 14.494 | 14.497 | -0.003 | 93 | 401249 | 50.0 | 51.5 | |
| 123 Hexachlorobutadiene | 225 | 14.646 | 14.643 | 0.003 | 93 | 135785 | 50.0 | 55.1 | |
| 124 Naphthalene | 128 | 14.761 | 14.758 | 0.003 | 98 | 1047608 | 50.0 | 44.2 | |
| 125 1,2,3-Trichlorobenzene | 180 | 14.980 | 14.977 | 0.003 | 92 | 360452 | 50.0 | 51.6 | |
| 126 2,4,5-Trichlorotoluene | 159 | 15.777 | 15.774 | 0.003 | 0 | 224866 | 50.0 | 55.9 | |
| 127 2,3,6-Trichlorotoluene | 159 | 15.875 | 15.877 | -0.002 | 95 | 223918 | 50.0 | 62.1 | |
| 146 3,4-Dichlorotoluene | 1 | | 0.000 | | | | ND | ND | |
| S 131 Xylenes, Total | 106 | | | | | 0 | 100.0 | 91.2 | |
| S 130 1,2-Dichloroethene, Total | 96 | | | | | 0 | 100.0 | 91.9 | |
| S 132 1,3-Dichloropropene, Total | 1 | | | | | 0 | 100.0 | 92.7 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

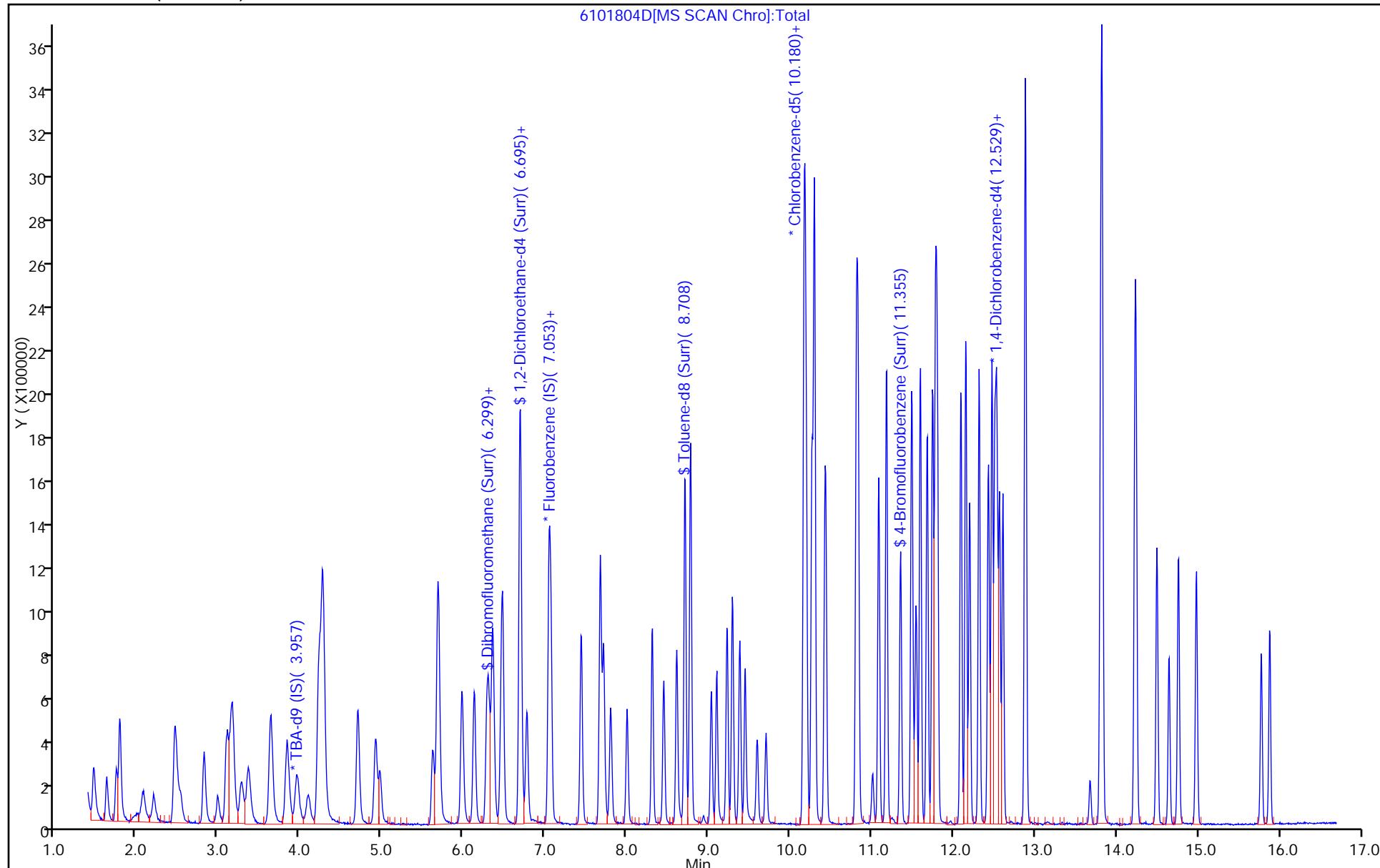
Reagents:

| | | | |
|---------------------|--------------------|-----------|-------------|
| voaWKetmix1st_00006 | Amount Added: 2.00 | Units: uL | |
| voaWVA1stRest_00021 | Amount Added: 2.00 | Units: uL | |
| voaWAcro1stRe_00021 | Amount Added: 6.00 | Units: uL | |
| voaWEEmix1stR_00014 | Amount Added: 2.00 | Units: uL | |
| VOA2CEVE2ND_00010 | Amount Added: 2.00 | Units: uL | |
| VOA8260VOAPRI_00266 | Amount Added: 2.00 | Units: uL | |
| VOA8260INT_00074 | Amount Added: 2.00 | Units: uL | Run Reagent |
| VOA8260SURR_00073 | Amount Added: 2.00 | Units: uL | Run Reagent |

Report Date: 18-Oct-2017 20:30:47

Chrom Revision: 2.2 16-Aug-2017 16:24:46

TestAmerica Pittsburgh
Data File: \\ChromNA\\Pittsburgh\\ChromData\\CHHP6\\20171017-18914.b\\6101804D.D
Injection Date: 18-Oct-2017 01:04:30 Instrument ID: CHHP6
Lims ID: LCS Operator ID: 034635
Client ID:
Purge Vol: 5.000 mL Dil. Factor: 1.0000 ALS Bottle#: 4
Method: MSVOA_LL_CHHP6 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Recovery Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\6101804.D.D
 Lims ID: LCS
 Client ID:
 Sample Type: LCS
 Inject. Date: 18-Oct-2017 01:04:30 ALS Bottle#: 4 Worklist Smp#: 4
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018914-004
 Misc. Info.: LCS
 Operator ID: 034635 Instrument ID: CHHP6
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP6\20171017-18914.b\MSVOA_LL_CHHP6.m
 Limit Group: VOA 8260C ICAL
 Last Update: 18-Oct-2017 20:30:42 Calib Date: 28-Sep-2017 15:13:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICAL File: \\ChromNA\Pittsburgh\ChromData\CHHP6\20170928-18631.b\60928P06.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK031

First Level Reviewer: bungardf Date: 18-Oct-2017 01:25:42

| Compound | Amount Added | Amount Recovered | % Rec. |
|-----------------------------------|--------------|------------------|--------|
| \$ 5 Dibromofluoromethane (Surr) | 50.0 | 46.1 | 92.16 |
| \$ 6 1,2-Dichloroethane-d4 (Surr) | 50.0 | 45.8 | 91.50 |
| \$ 7 Toluene-d8 (Surr) | 50.0 | 46.5 | 92.91 |
| \$ 8 4-Bromofluorobenzene (Surr) | 50.0 | 46.8 | 93.63 |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Pittsburgh Job No.: 180-71131-1

SDG No.: _____

Instrument ID: CHHP6 Start Date: 06/02/2017 06:02Analysis Batch Number: 213005 End Date: 06/02/2017 15:14

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|-------------------|------------------|------------------|-----------------|-------------|------------------|
| BFB 180-213005/3 | | 06/02/2017 06:02 | 1 | 60602003.D | DB-624 0.18 (mm) |
| IC 180-213005/14 | | 06/02/2017 11:37 | 1 | | DB-624 0.18 (mm) |
| IC 180-213005/15 | | 06/02/2017 12:01 | 1 | | DB-624 0.18 (mm) |
| IC 180-213005/16 | | 06/02/2017 12:25 | 1 | | DB-624 0.18 (mm) |
| IC 180-213005/17 | | 06/02/2017 12:49 | 1 | | DB-624 0.18 (mm) |
| IC 180-213005/18 | | 06/02/2017 13:13 | 1 | | DB-624 0.18 (mm) |
| IC 180-213005/19 | | 06/02/2017 13:37 | 1 | | DB-624 0.18 (mm) |
| IC 180-213005/20 | | 06/02/2017 14:02 | 1 | | DB-624 0.18 (mm) |
| IC 180-213005/21 | | 06/02/2017 14:26 | 1 | | DB-624 0.18 (mm) |
| ICV 180-213005/23 | | 06/02/2017 15:14 | 1 | 60602023.D | DB-624 0.18 (mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

Instrument ID: CHHP6

Start Date: 07/24/2017 04:58

Analysis Batch Number: 217861

End Date: 07/24/2017 16:22

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|----------------------|------------------|------------------|-----------------|-------------|------------------|
| BFB 180-217861/1 | | 07/24/2017 04:58 | 1 | 60724D01.D | DB-624 0.18 (mm) |
| ZZZZZ | | 07/24/2017 05:31 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 07/24/2017 05:31 | 1 | | DB-624 0.18 (mm) |
| IC 180-217861/3 | | 07/24/2017 06:39 | 1 | 60724D03.D | DB-624 0.18 (mm) |
| IC 180-217861/4 | | 07/24/2017 07:03 | 1 | 60724D04.D | DB-624 0.18 (mm) |
| ICIS 180-217861/5 | | 07/24/2017 07:27 | 1 | 60724D05.D | DB-624 0.18 (mm) |
| IC 180-217861/6 | | 07/24/2017 07:52 | 1 | 60724D06.D | DB-624 0.18 (mm) |
| IC 180-217861/7 | | 07/24/2017 08:16 | 1 | 60724D07.D | DB-624 0.18 (mm) |
| IC 180-217861/8 | | 07/24/2017 08:40 | 1 | 60724D08.D | DB-624 0.18 (mm) |
| IC 180-217861/9 | | 07/24/2017 09:04 | 1 | 60724D09.D | DB-624 0.18 (mm) |
| IC 180-217861/10 | | 07/24/2017 09:28 | 1 | 60724D10.D | DB-624 0.18 (mm) |
| ZZZZZ | | 07/24/2017 09:52 | 1 | | DB-624 0.18 (mm) |
| MDLV 180-217861/1011 | | 07/24/2017 09:52 | 1 | | DB-624 0.18 (mm) |
| CCV 180-217861/12 | | 07/24/2017 10:16 | 1 | | DB-624 0.18 (mm) |
| ICV 180-217861/13 | | 07/24/2017 10:40 | 1 | 60724D13.D | DB-624 0.18 (mm) |
| ZZZZZ | | 07/24/2017 11:04 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 07/24/2017 11:53 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 07/24/2017 12:17 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 07/24/2017 12:42 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 07/24/2017 13:37 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 07/24/2017 14:14 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 07/24/2017 14:45 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 07/24/2017 15:09 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 07/24/2017 15:33 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 07/24/2017 16:22 | 1 | | DB-624 0.18 (mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Pittsburgh

Job No.: 180-71131-1

SDG No.:

Instrument ID: CHHP6

Start Date: 10/17/2017 21:58

Analysis Batch Number: 226148

End Date: 10/18/2017 09:49

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|--------------------|------------------|------------------|-----------------|-------------|------------------|
| BFB 180-226148/1 | | 10/17/2017 21:58 | 1 | 6101801D.D | DB-624 0.18 (mm) |
| CCVIS 180-226148/2 | | 10/17/2017 23:55 | 1 | 6101802D.D | DB-624 0.18 (mm) |
| ZZZZZ | | 10/17/2017 23:55 | 1 | | DB-624 0.18 (mm) |
| CCV 180-226148/3 | | 10/18/2017 00:36 | 1 | | DB-624 0.18 (mm) |
| LCS 180-226148/4 | | 10/18/2017 01:04 | 1 | 6101804D.D | DB-624 0.18 (mm) |
| ZZZZZ | | 10/18/2017 01:43 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/18/2017 01:43 | 1 | | DB-624 0.18 (mm) |
| MB 180-226148/6 | | 10/18/2017 02:09 | 1 | 6101806D.D | DB-624 0.18 (mm) |
| ZZZZZ | | 10/18/2017 02:43 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/18/2017 03:11 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/18/2017 03:43 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/18/2017 04:32 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/18/2017 04:56 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/18/2017 05:20 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/18/2017 05:45 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/18/2017 06:09 | 1 | | DB-624 0.18 (mm) |
| 180-71131-1 | | 10/18/2017 06:33 | 1 | 6101816D.D | DB-624 0.18 (mm) |
| 180-71131-2 | | 10/18/2017 06:57 | 1 | 6101817D.D | DB-624 0.18 (mm) |
| 180-71131-3 | | 10/18/2017 07:45 | 1 | 6101819D.D | DB-624 0.18 (mm) |
| ZZZZZ | | 10/18/2017 08:10 | 100 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/18/2017 08:34 | 100 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/18/2017 08:58 | 200 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/18/2017 09:25 | 200 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/18/2017 09:49 | 100 | | DB-624 0.18 (mm) |

Shipping and Receiving Documents

b = weight entered manually
d = weight read from scale
f = faxed file
For more information, see
FBI Laboratory Guide for details.
Visit us at: FEDEX.COM
or call 1-800-824-8639
1-800-463-3329
Oct 09, 2017 9:05:10 AM

Chain of Custody Record

| | | | | | | | | |
|---|--|--|-------------|---|--|--------------------------------------|-----------------------|---------------------|
| TestAmerica Pittsburgh 301 Alpha Drive | | Pittsburgh, PA 15238 Phone: 412.963.7010 Fax: 412.963.2470 | | Regulatory Program: <input type="checkbox"/> DW <input type="checkbox"/> NPDES <input type="checkbox"/> RORA <input type="checkbox"/> Other | | | | |
| Client Contact | | Project Manager: <u>Chasen</u> | | 170410 | | | | |
| Company Name: <u>Ground surface soil</u> Address: <u>7601 Market Place Ste 310</u> City/State/Zip: <u>Harrisburg, PA 17110</u> Phone: <u>217.672.0437</u> Fax: <u></u> Project Name: <u>Hockey-Park</u> Site: <u>SBA</u> PO # <u>1002.36</u> | | Tel/Fax: <input type="checkbox"/> ANALYSIS TURNAROUND TIME <input type="checkbox"/> CALENDAR DAYS <input type="checkbox"/> WORKING DAYS <input checked="" type="checkbox"/> TAT different from below 2 weeks 1 week 2 days 1 day | | Site Contact: <u>Eric Farmer</u> Lab Contact: <u>Chasen</u> Date: <u>10/10/17</u> Carrier: <u>FEAK</u> | | | | |
| Sample Identification | | Sample Date | Sample Time | Sample Type (O=Comp, G=Grab) | Matrix | # of Cont. | Filtered Sample (Y/N) | Perform MS/MS (Y/N) |
| HD-SpBA-CW-23-01-0 | | 10/10/17 | 0950 | G | 6 | 3 | Y | N |
| HD-CW-23-01-0 | | 10/10/17 | 0950 | G | 6 | 3 | Y | N |
| HD-QCG-01-2 | | 10/10/17 | 1200 | G | 2 | 2 | Y | N |
| Preservation Used: 1=Ice, 2=HCl; 3=H ₂ SO ₄ ; 4=HNO ₃ ; 5=NaOH; 6=Other | | | | | | | | |
| Possible Hazard Identification: Are any samples from a listed EPA Hazardous Waste? Please List any EPA Waste Codes for the sample in the Comments Section if the lab is to dispose of the sample. | | | | | | | | |
| <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"/> Return to Client <input type="checkbox"/> Disposal by Lab <input type="checkbox"/> Archive for _____ Months | | | |
| Special Instructions/QC Requirements & Comments: | | | | | | | | |
| Custody Seals Intact: <input type="checkbox"/> Yes <input type="checkbox"/> No | | Custody Seal No.: <u>G54</u> | | Cooler Temp. (°C): Obs'd: <u>10/10/17 0908</u> | | Com'd: <u></u> Therm ID No.: <u></u> | | |
| Relinquished by: <u>JM</u> | | Company: <u>G54</u> | | Received by: <u></u> | | Company: <u></u> Date/Time: <u></u> | | |
| Relinquished by: <u></u> | | Company: <u></u> | | Received by: <u></u> | | Company: <u></u> Date/Time: <u></u> | | |
| Relinquished by: <u></u> | | Company: <u></u> | | Received in Laboratory by: <u></u> | | Company: <u></u> Date/Time: <u></u> | | |

Chain of Custody Record

Pittsburgh, PA 15238
Phone: 412.963.7050 Fax: 412.963.2470

170410

| | | | | | | |
|---|---|---|-------------------------------------|-------------------------------------|-------------------------------------|---|
| Client Contact | Project Manager: | Regulatory Program: | <input type="checkbox"/> DW | <input type="checkbox"/> NPDES | <input type="checkbox"/> RCRA | <input type="checkbox"/> Other: |
| Company Name: Grant water services corp Address: 2601 Morris Park Ave Sec 310 City/State/Zip: Morrisburg PA 17110 Phone: 317-652-6432 Fax: _____ Project Name: Turkey Run Site: SPBA PO # 10022-36 | Tel/Fax: Analysis Turnaround Time <input type="checkbox"/> CALENDAR DAYS <input type="checkbox"/> WORKING DAYS TAT if different from Below <input checked="" type="checkbox"/> 2 weeks <input type="checkbox"/> 1 week <input type="checkbox"/> 2 days <input type="checkbox"/> 1 day | Lab Contact: Client Contact <input checked="" type="checkbox"/> Carrier: FedEx | Date: 10/19/11 | COC No: 1017100901 | | |
| Sample Identification | Sample Date | Sample Time | Sample Type (C=Comp., G=Grab) | Matrix | # of Cont. | Sample Specific Notes: (48 hrs sample) (72 hrs sample) top blank |
| HD-SpBA-cw/2310/1-0 HO-cw-23-01-0 HO-DCC-01-2 | 10/15/11 10/16/11 10/15/11 | 0950 0950 1200 | G G VW | 3 3 2 | 3 3 2 | |
| Preservation Used: 1=Ice, 2=HCl; 3=H2SO4; 4=HNO3; 5=NaOH; 6= Other | | | | | | Sample Disposal (A fee may be assessed if samples are retained longer than 1 month) Are any samples from a listed EPA Hazardous Waste? Please List any EPA Waste Codes for the sample in the Comments Section if the lab is to dispose of the sample. <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"/> Return to Client <input type="checkbox"/> Disposal by Lab <input type="checkbox"/> Disposal by Company <input type="checkbox"/> Archive for _____ Months |
| | | | | | | Special Instructions/QC Requirements & Comments: |
| Custody Seals Intact: <input type="checkbox"/> Yes <input type="checkbox"/> No | | Custody Seal No.: Company: | Cooler Temp. (°C): Obs'd: _____ | Corrd: _____ | Therm ID No.: _____ | |
| Relinquished by: _____ Relinquished by: _____ Relinquished by: _____ | | Date/Time: Received by: Company: | Date/Time: Received by: Company: | Date/Time: Received by: Company: | Date/Time: Received by: Company: | Date/Time: Received by: Company: |
| | | | | | | |

ORIGIN ID:THVA (631) 766-2976
KAITLIN FLEMING
GROUNDWATER SCIENCES CORPORATION
2601 MARKET PL STE 310
HARRISBURG, PA 17110
UNITED STATES US

SHIP DATE: 09OCT17
ACTWT: 10.90 LB
CAD: 006995074/SSFE1B22
DIMS: 12x11x11 IN
BILL RECIPIENT

DP # 1569974-35 RT12/15/15
34545

TO **ATTN: SAMPLE RECEIVING
TEST AMERICA
301 ALPHA DR**

PITTSBURGH PA 15238

(412) 963-7058
TRK#
PC:

REF#

DEPT#



TRK#
0201 7880 0479 0607

TUE - 10 OCT 10:30A
PRIORITY OVERNIGHT

E8 AGCA

**5238
PIT**

Uncorrected temp
Thermometer ID

CF O Initials JS

PT-WI-SR-001 effective 7/26/13

118 °C
13



180-71131 Waybill

Login Sample Receipt Checklist

Client: Groundwater Sciences Corporation

Job Number: 180-71131-1

Login Number: 71131

List Source: TestAmerica Pittsburgh

List Number: 1

Creator: Say, Thomas C

| 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 (excluding tests with immediate HTs) | 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 | |