

ANALYTICAL REPORT

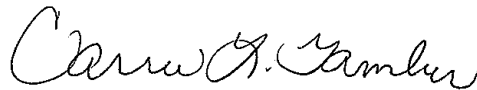
Job Number: 180-70873-1

Job Description: Harley Davidson

For:

Groundwater Sciences Corporation
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Attention: Christopher O'Neil



Approved for release.
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10/10/2017 7:22 AM

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10/10/2017

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Definitions/Glossary

Client: Groundwater Sciences Corporation
Project/Site: Harley Davidson

TestAmerica Job ID: 180-70873-1

Qualifiers

GC/MS VOA

| Qualifier | Qualifier Description |
|-----------|--|
| U | Indicates the analyte was analyzed for but not detected. |
| ^c | CCV Recovery is outside acceptance limits. |
| E | Result exceeded calibration range. |
| 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. |
|----------------|---|
| α | 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: 180-70873-1

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 10/3/2017 9:00 AM; the samples arrived in good condition, properly preserved and, where required, on ice. The temperature of the cooler at receipt was 2.4° C.

The following Trip Blank sample was received with headspace in 2 of 2 vials. The client was emailed and the analysis proceeded with this narration.

VOLATILES

The following sample was diluted to bring the concentration of target analytes within the calibration range: HD-SPBA-CW-22-0/1-0 (180-70873-1). Elevated reporting limits (RLs) are provided.

The continuing calibration verification (CCV) analyzed in batch 180-224792 was outside the method criteria for the following analytes: 1,4-Dioxane, Bromoform, Chloromethane, cis-1,3-Dichloropropene and Dibromochloromethane. 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.

The continuing calibration verification (CCV) analyzed in batch 180-224919 was outside the method criteria for the following analytes: 1,4-Dioxane, Chloroethane, Chloromethane and Vinyl chloride. 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: 180-70873-1

Client Sample ID: HD-SPBA-CW-22-0/1-0

Lab Sample ID: 180-70873-1

| Analyte | Result | Qualifier | RL | MDL | Unit | Dil Fac | D | Method | Prep Type |
|------------------------|--------|-----------|-----|------|------|---------|---|--------|-----------|
| cis-1,2-Dichloroethene | 1.7 | | 1.0 | 0.30 | ug/L | 1 | | 8260C | Total/NA |
| Trichloroethene | 220 | E | 1.0 | 0.20 | ug/L | 1 | | 8260C | Total/NA |
| Tetrachloroethene | 320 | E | 1.0 | 0.24 | ug/L | 1 | | 8260C | Total/NA |
| Trichloroethene - DL | 190 | | 13 | 2.5 | ug/L | 12.5 | | 8260C | Total/NA |
| Tetrachloroethene - DL | 290 | | 13 | 3.1 | ug/L | 12.5 | | 8260C | Total/NA |

Client Sample ID: HD-QC4-0/1-2

Lab Sample ID: 180-70873-2

No Detections.

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: 180-70873-1

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

Client Sample ID: HD-SPBA-CW-22-0/1-0

Date Collected: 09/29/17 09:40

Date Received: 10/03/17 09:00

Lab Sample ID: 180-70873-1

Matrix: Water

| Analyte | Result | Qualifier | RL | MDL | Unit | D | Prepared | Analyzed | Dil Fac |
|-------------------------------|------------|-----------|-----|------|------|---|----------|----------------|---------|
| Chloromethane | 1.0 | U ^c | 1.0 | 0.38 | ug/L | | | 10/04/17 05:23 | 1 |
| Vinyl chloride | 1.0 | U | 1.0 | 0.17 | ug/L | | | 10/04/17 05:23 | 1 |
| Bromomethane | 1.0 | U | 1.0 | 0.59 | ug/L | | | 10/04/17 05:23 | 1 |
| Chloroethane | 1.0 | U | 1.0 | 0.58 | ug/L | | | 10/04/17 05:23 | 1 |
| 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.32 | ug/L | | | 10/04/17 05:23 | 1 |
| Acetone | 5.0 | U | 5.0 | 3.1 | ug/L | | | 10/04/17 05:23 | 1 |
| Carbon disulfide | 1.0 | U | 1.0 | 0.53 | ug/L | | | 10/04/17 05:23 | 1 |
| Methylene Chloride | 1.0 | U | 1.0 | 0.94 | ug/L | | | 10/04/17 05:23 | 1 |
| trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.20 | ug/L | | | 10/04/17 05:23 | 1 |
| Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.20 | ug/L | | | 10/04/17 05:23 | 1 |
| 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.34 | ug/L | | | 10/04/17 05:23 | 1 |
| cis-1,2-Dichloroethene | 1.7 | | 1.0 | 0.30 | ug/L | | | 10/04/17 05:23 | 1 |
| Bromochloromethane | 1.0 | U | 1.0 | 0.36 | ug/L | | | 10/04/17 05:23 | 1 |
| 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 | ug/L | | | 10/04/17 05:23 | 1 |
| Chloroform | 1.0 | U | 1.0 | 0.27 | ug/L | | | 10/04/17 05:23 | 1 |
| 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.27 | ug/L | | | 10/04/17 05:23 | 1 |
| Carbon tetrachloride | 1.0 | U | 1.0 | 0.56 | ug/L | | | 10/04/17 05:23 | 1 |
| Benzene | 1.0 | U | 1.0 | 0.18 | ug/L | | | 10/04/17 05:23 | 1 |
| 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.24 | ug/L | | | 10/04/17 05:23 | 1 |
| Trichloroethene | 220 | E | 1.0 | 0.20 | ug/L | | | 10/04/17 05:23 | 1 |
| 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.35 | ug/L | | | 10/04/17 05:23 | 1 |
| Bromodichloromethane | 1.0 | U ^c | 1.0 | 0.57 | ug/L | | | 10/04/17 05:23 | 1 |
| cis-1,3-Dichloropropene | 1.0 | U ^c | 1.0 | 0.32 | ug/L | | | 10/04/17 05:23 | 1 |
| 4-Methyl-2-pentanone (MIBK) | 5.0 | U | 5.0 | 2.2 | ug/L | | | 10/04/17 05:23 | 1 |
| Toluene | 1.0 | U | 1.0 | 0.16 | ug/L | | | 10/04/17 05:23 | 1 |
| trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.22 | ug/L | | | 10/04/17 05:23 | 1 |
| 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.31 | ug/L | | | 10/04/17 05:23 | 1 |
| Tetrachloroethene | 320 | E | 1.0 | 0.24 | ug/L | | | 10/04/17 05:23 | 1 |
| 2-Hexanone | 5.0 | U | 5.0 | 2.0 | ug/L | | | 10/04/17 05:23 | 1 |
| Dibromochloromethane | 1.0 | U | 1.0 | 0.44 | ug/L | | | 10/04/17 05:23 | 1 |
| 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.51 | ug/L | | | 10/04/17 05:23 | 1 |
| Chlorobenzene | 1.0 | U | 1.0 | 0.15 | ug/L | | | 10/04/17 05:23 | 1 |
| 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.49 | ug/L | | | 10/04/17 05:23 | 1 |
| Ethylbenzene | 1.0 | U | 1.0 | 0.25 | ug/L | | | 10/04/17 05:23 | 1 |
| Xylenes, Total | 2.0 | U | 2.0 | 0.27 | ug/L | | | 10/04/17 05:23 | 1 |
| Styrene | 1.0 | U | 1.0 | 0.22 | ug/L | | | 10/04/17 05:23 | 1 |
| Bromoform | 1.0 | U ^c | 1.0 | 0.76 | ug/L | | | 10/04/17 05:23 | 1 |
| 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.37 | ug/L | | | 10/04/17 05:23 | 1 |
| Acrylonitrile | 20 | U | 20 | 3.3 | ug/L | | | 10/04/17 05:23 | 1 |
| 1,4-Dioxane | 200 | U | 200 | 16 | ug/L | | | 10/04/17 05:23 | 1 |

| Surrogate | %Recovery | Qualifier | Limits | Prepared | Analyzed | Dil Fac |
|------------------------------|-----------|-----------|----------|----------|----------------|---------|
| 1,2-Dichloroethane-d4 (Surr) | 113 | | 65 - 121 | | 10/04/17 05:23 | 1 |
| Toluene-d8 (Surr) | 97 | | 73 - 120 | | 10/04/17 05:23 | 1 |
| 4-Bromofluorobenzene (Surr) | 94 | | 80 - 120 | | 10/04/17 05:23 | 1 |
| Dibromofluoromethane (Surr) | 101 | | 73 - 120 | | 10/04/17 05:23 | 1 |

Client Sample Results

Client: Groundwater Sciences Corporation
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-70873-1

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

Client Sample ID: HD-QC4-0/1-2

Date Collected: 09/29/17 12:00

Date Received: 10/03/17 09:00

Lab Sample ID: 180-70873-2

Matrix: Water

| Analyte | Result | Qualifier | RL | MDL | Unit | D | Prepared | Analyzed | Dil Fac |
|-----------------------------|--------|-----------|-----|------|------|---|----------|----------------|---------|
| Chloromethane | 1.0 | U ^c | 1.0 | 0.38 | ug/L | | | 10/04/17 04:59 | 1 |
| Vinyl chloride | 1.0 | U | 1.0 | 0.17 | ug/L | | | 10/04/17 04:59 | 1 |
| Bromomethane | 1.0 | U | 1.0 | 0.59 | ug/L | | | 10/04/17 04:59 | 1 |
| Chloroethane | 1.0 | U | 1.0 | 0.58 | ug/L | | | 10/04/17 04:59 | 1 |
| 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.32 | ug/L | | | 10/04/17 04:59 | 1 |
| Acetone | 5.0 | U | 5.0 | 3.1 | ug/L | | | 10/04/17 04:59 | 1 |
| Carbon disulfide | 1.0 | U | 1.0 | 0.53 | ug/L | | | 10/04/17 04:59 | 1 |
| Methylene Chloride | 1.0 | U | 1.0 | 0.94 | ug/L | | | 10/04/17 04:59 | 1 |
| trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.20 | ug/L | | | 10/04/17 04:59 | 1 |
| Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.20 | ug/L | | | 10/04/17 04:59 | 1 |
| 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.34 | ug/L | | | 10/04/17 04:59 | 1 |
| cis-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.30 | ug/L | | | 10/04/17 04:59 | 1 |
| Bromochloromethane | 1.0 | U | 1.0 | 0.36 | ug/L | | | 10/04/17 04:59 | 1 |
| 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 | ug/L | | | 10/04/17 04:59 | 1 |
| Chloroform | 1.0 | U | 1.0 | 0.27 | ug/L | | | 10/04/17 04:59 | 1 |
| 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.27 | ug/L | | | 10/04/17 04:59 | 1 |
| Carbon tetrachloride | 1.0 | U | 1.0 | 0.56 | ug/L | | | 10/04/17 04:59 | 1 |
| Benzene | 1.0 | U | 1.0 | 0.18 | ug/L | | | 10/04/17 04:59 | 1 |
| 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.24 | ug/L | | | 10/04/17 04:59 | 1 |
| Trichloroethene | 1.0 | U | 1.0 | 0.20 | ug/L | | | 10/04/17 04:59 | 1 |
| 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.35 | ug/L | | | 10/04/17 04:59 | 1 |
| Bromodichloromethane | 1.0 | U ^c | 1.0 | 0.57 | ug/L | | | 10/04/17 04:59 | 1 |
| cis-1,3-Dichloropropene | 1.0 | U ^c | 1.0 | 0.32 | ug/L | | | 10/04/17 04:59 | 1 |
| 4-Methyl-2-pentanone (MIBK) | 5.0 | U | 5.0 | 2.2 | ug/L | | | 10/04/17 04:59 | 1 |
| Toluene | 1.0 | U | 1.0 | 0.16 | ug/L | | | 10/04/17 04:59 | 1 |
| trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.22 | ug/L | | | 10/04/17 04:59 | 1 |
| 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.31 | ug/L | | | 10/04/17 04:59 | 1 |
| Tetrachloroethene | 1.0 | U | 1.0 | 0.24 | ug/L | | | 10/04/17 04:59 | 1 |
| 2-Hexanone | 5.0 | U | 5.0 | 2.0 | ug/L | | | 10/04/17 04:59 | 1 |
| Dibromochloromethane | 1.0 | U | 1.0 | 0.44 | ug/L | | | 10/04/17 04:59 | 1 |
| 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.51 | ug/L | | | 10/04/17 04:59 | 1 |
| Chlorobenzene | 1.0 | U | 1.0 | 0.15 | ug/L | | | 10/04/17 04:59 | 1 |
| 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.49 | ug/L | | | 10/04/17 04:59 | 1 |
| Ethylbenzene | 1.0 | U | 1.0 | 0.25 | ug/L | | | 10/04/17 04:59 | 1 |
| Xylenes, Total | 2.0 | U | 2.0 | 0.27 | ug/L | | | 10/04/17 04:59 | 1 |
| Styrene | 1.0 | U | 1.0 | 0.22 | ug/L | | | 10/04/17 04:59 | 1 |
| Bromoform | 1.0 | U ^c | 1.0 | 0.76 | ug/L | | | 10/04/17 04:59 | 1 |
| 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.37 | ug/L | | | 10/04/17 04:59 | 1 |
| Acrylonitrile | 20 | U | 20 | 3.3 | ug/L | | | 10/04/17 04:59 | 1 |
| 1,4-Dioxane | 200 | U | 200 | 16 | ug/L | | | 10/04/17 04:59 | 1 |

| Surrogate | %Recovery | Qualifier | Limits | Prepared | Analyzed | Dil Fac |
|------------------------------|-----------|-----------|----------|----------|----------------|---------|
| 1,2-Dichloroethane-d4 (Surr) | 113 | | 65 - 121 | | 10/04/17 04:59 | 1 |
| Toluene-d8 (Surr) | 101 | | 73 - 120 | | 10/04/17 04:59 | 1 |
| 4-Bromofluorobenzene (Surr) | 92 | | 80 - 120 | | 10/04/17 04:59 | 1 |
| Dibromofluoromethane (Surr) | 101 | | 73 - 120 | | 10/04/17 04:59 | 1 |

Client Sample Results

Client: Groundwater Sciences Corporation
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-70873-1

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

Client Sample ID: HD-SPBA-CW-22-0/1-0

Lab Sample ID: 180-70873-1

Date Collected: 09/29/17 09:40

Matrix: Water

Date Received: 10/03/17 09:00

| Analyte | Result | Qualifier | RL | MDL | Unit | D | Prepared | Analyzed | Dil Fac |
|-----------------------------|------------|-----------|------|-----|------|---|----------|----------------|---------|
| Chloromethane | 13 | U ^c | 13 | 4.8 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Vinyl chloride | 13 | U ^c | 13 | 2.1 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Bromomethane | 13 | U | 13 | 7.3 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Chloroethane | 13 | U ^c | 13 | 7.2 | ug/L | | | 10/05/17 06:31 | 12.5 |
| 1,1-Dichloroethene | 13 | U | 13 | 4.0 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Acetone | 63 | U | 63 | 39 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Carbon disulfide | 13 | U | 13 | 6.6 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Methylene Chloride | 13 | U | 13 | 12 | ug/L | | | 10/05/17 06:31 | 12.5 |
| trans-1,2-Dichloroethene | 13 | U | 13 | 2.5 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Methyl tert-butyl ether | 13 | U | 13 | 2.4 | ug/L | | | 10/05/17 06:31 | 12.5 |
| 1,1-Dichloroethane | 13 | U | 13 | 4.2 | ug/L | | | 10/05/17 06:31 | 12.5 |
| cis-1,2-Dichloroethene | 13 | U | 13 | 3.8 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Bromochloromethane | 13 | U | 13 | 4.5 | ug/L | | | 10/05/17 06:31 | 12.5 |
| 2-Butanone (MEK) | 63 | U | 63 | 32 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Chloroform | 13 | U | 13 | 3.3 | ug/L | | | 10/05/17 06:31 | 12.5 |
| 1,1,1-Trichloroethane | 13 | U | 13 | 3.4 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Carbon tetrachloride | 13 | U | 13 | 7.0 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Benzene | 13 | U | 13 | 2.3 | ug/L | | | 10/05/17 06:31 | 12.5 |
| 1,2-Dichloroethane | 13 | U | 13 | 3.0 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Trichloroethene | 190 | | 13 | 2.5 | ug/L | | | 10/05/17 06:31 | 12.5 |
| 1,2-Dichloropropane | 13 | U | 13 | 4.3 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Bromodichloromethane | 13 | U | 13 | 7.1 | ug/L | | | 10/05/17 06:31 | 12.5 |
| cis-1,3-Dichloropropene | 13 | U | 13 | 4.0 | ug/L | | | 10/05/17 06:31 | 12.5 |
| 4-Methyl-2-pentanone (MIBK) | 63 | U | 63 | 27 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Toluene | 13 | U | 13 | 2.0 | ug/L | | | 10/05/17 06:31 | 12.5 |
| trans-1,3-Dichloropropene | 13 | U | 13 | 2.8 | ug/L | | | 10/05/17 06:31 | 12.5 |
| 1,1,2-Trichloroethane | 13 | U | 13 | 3.8 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Tetrachloroethene | 290 | | 13 | 3.1 | ug/L | | | 10/05/17 06:31 | 12.5 |
| 2-Hexanone | 63 | U | 63 | 25 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Dibromochloromethane | 13 | U | 13 | 5.5 | ug/L | | | 10/05/17 06:31 | 12.5 |
| 1,2-Dibromoethane (EDB) | 13 | U | 13 | 6.4 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Chlorobenzene | 13 | U | 13 | 1.8 | ug/L | | | 10/05/17 06:31 | 12.5 |
| 1,1,1,2-Tetrachloroethane | 13 | U | 13 | 6.2 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Ethylbenzene | 13 | U | 13 | 3.2 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Xylenes, Total | 25 | U | 25 | 3.4 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Styrene | 13 | U | 13 | 2.7 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Bromoform | 13 | U | 13 | 9.5 | ug/L | | | 10/05/17 06:31 | 12.5 |
| 1,1,2,2-Tetrachloroethane | 13 | U | 13 | 4.6 | ug/L | | | 10/05/17 06:31 | 12.5 |
| Acrylonitrile | 250 | U | 250 | 42 | ug/L | | | 10/05/17 06:31 | 12.5 |
| 1,4-Dioxane | 2500 | U | 2500 | 200 | ug/L | | | 10/05/17 06:31 | 12.5 |

| Surrogate | %Recovery | Qualifier | Limits | Prepared | Analyzed | Dil Fac |
|------------------------------|-----------|-----------|----------|----------|----------------|---------|
| 1,2-Dichloroethane-d4 (Surr) | 117 | | 65 - 121 | | 10/05/17 06:31 | 12.5 |
| Toluene-d8 (Surr) | 99 | | 73 - 120 | | 10/05/17 06:31 | 12.5 |
| 4-Bromofluorobenzene (Surr) | 94 | | 80 - 120 | | 10/05/17 06:31 | 12.5 |
| Dibromofluoromethane (Surr) | 109 | | 73 - 120 | | 10/05/17 06:31 | 12.5 |

Default Detection Limits

Client: Groundwater Sciences Corporation
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-70873-1

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

| Analyte | RL | MDL | Units | Method |
|-----------------------------|-----|------|-------|--------|
| 1,1,1,2-Tetrachloroethane | 1.0 | 0.49 | ug/L | 8260C |
| 1,1,1-Trichloroethane | 1.0 | 0.27 | ug/L | 8260C |
| 1,1,2,2-Tetrachloroethane | 1.0 | 0.37 | ug/L | 8260C |
| 1,1,2-Trichloroethane | 1.0 | 0.31 | ug/L | 8260C |
| 1,1-Dichloroethane | 1.0 | 0.34 | ug/L | 8260C |
| 1,1-Dichloroethene | 1.0 | 0.32 | ug/L | 8260C |
| 1,2-Dibromoethane (EDB) | 1.0 | 0.51 | ug/L | 8260C |
| 1,2-Dichloroethane | 1.0 | 0.24 | ug/L | 8260C |
| 1,2-Dichloropropane | 1.0 | 0.35 | ug/L | 8260C |
| 1,4-Dioxane | 200 | 16 | ug/L | 8260C |
| 2-Butanone (MEK) | 5.0 | 2.6 | ug/L | 8260C |
| 2-Hexanone | 5.0 | 2.0 | ug/L | 8260C |
| 4-Methyl-2-pentanone (MIBK) | 5.0 | 2.2 | ug/L | 8260C |
| Acetone | 5.0 | 3.1 | ug/L | 8260C |
| Acrylonitrile | 20 | 3.3 | ug/L | 8260C |
| Benzene | 1.0 | 0.18 | ug/L | 8260C |
| Bromochloromethane | 1.0 | 0.36 | ug/L | 8260C |
| Bromodichloromethane | 1.0 | 0.57 | ug/L | 8260C |
| Bromoform | 1.0 | 0.76 | ug/L | 8260C |
| Bromomethane | 1.0 | 0.59 | ug/L | 8260C |
| Carbon disulfide | 1.0 | 0.53 | ug/L | 8260C |
| Carbon tetrachloride | 1.0 | 0.56 | ug/L | 8260C |
| Chlorobenzene | 1.0 | 0.15 | ug/L | 8260C |
| Chloroethane | 1.0 | 0.58 | ug/L | 8260C |
| Chloroform | 1.0 | 0.27 | ug/L | 8260C |
| Chloromethane | 1.0 | 0.38 | ug/L | 8260C |
| cis-1,2-Dichloroethene | 1.0 | 0.30 | ug/L | 8260C |
| cis-1,3-Dichloropropene | 1.0 | 0.32 | ug/L | 8260C |
| Dibromochloromethane | 1.0 | 0.44 | ug/L | 8260C |
| Ethylbenzene | 1.0 | 0.25 | ug/L | 8260C |
| Methyl tert-butyl ether | 1.0 | 0.20 | ug/L | 8260C |
| Methylene Chloride | 1.0 | 0.94 | ug/L | 8260C |
| Styrene | 1.0 | 0.22 | ug/L | 8260C |
| Tetrachloroethene | 1.0 | 0.24 | ug/L | 8260C |
| Toluene | 1.0 | 0.16 | ug/L | 8260C |
| trans-1,2-Dichloroethene | 1.0 | 0.20 | ug/L | 8260C |
| trans-1,3-Dichloropropene | 1.0 | 0.22 | ug/L | 8260C |
| Trichloroethene | 1.0 | 0.20 | ug/L | 8260C |
| Vinyl chloride | 1.0 | 0.17 | ug/L | 8260C |
| Xylenes, Total | 2.0 | 0.27 | ug/L | 8260C |

Surrogate Summary

Client: Groundwater Sciences Corporation
Project/Site: Harley Davidson

TestAmerica Job ID: 180-70873-1

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) |
| 180-70873-1 | HD-SPBA-CW-22-0/1-0 | 113 | 97 | 94 | 101 |
| 180-70873-1 - DL | HD-SPBA-CW-22-0/1-0 | 117 | 99 | 94 | 109 |
| 180-70873-2 | HD-QC4-0/1-2 | 113 | 101 | 92 | 101 |
| LCS 180-224792/4 | Lab Control Sample | 106 | 113 | 107 | 98 |
| LCS 180-224919/4 | Lab Control Sample | 102 | 110 | 102 | 90 |
| MB 180-224792/6 | Method Blank | 107 | 98 | 93 | 94 |
| MB 180-224919/7 | Method Blank | 111 | 101 | 95 | 96 |

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: 180-70873-1

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

Lab Sample ID: MB 180-224792/6

Matrix: Water

Analysis Batch: 224792

Client Sample ID: Method Blank

Prep Type: Total/NA

| Analyte | MB | MB | RL | MDL | Unit | D | Prepared | Analyzed | Dil Fac |
|-----------------------------|--------|-----------|-----|------|------|---|----------|----------------|---------|
| | Result | Qualifier | | | | | | | |
| Chloromethane | 1.0 | U | 1.0 | 0.38 | ug/L | | | 10/04/17 02:21 | 1 |
| Vinyl chloride | 1.0 | U | 1.0 | 0.17 | ug/L | | | 10/04/17 02:21 | 1 |
| Bromomethane | 1.0 | U | 1.0 | 0.59 | ug/L | | | 10/04/17 02:21 | 1 |
| Chloroethane | 1.0 | U | 1.0 | 0.58 | ug/L | | | 10/04/17 02:21 | 1 |
| 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.32 | ug/L | | | 10/04/17 02:21 | 1 |
| Acetone | 5.0 | U | 5.0 | 3.1 | ug/L | | | 10/04/17 02:21 | 1 |
| Carbon disulfide | 1.0 | U | 1.0 | 0.53 | ug/L | | | 10/04/17 02:21 | 1 |
| Methylene Chloride | 1.0 | U | 1.0 | 0.94 | ug/L | | | 10/04/17 02:21 | 1 |
| trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.20 | ug/L | | | 10/04/17 02:21 | 1 |
| Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.20 | ug/L | | | 10/04/17 02:21 | 1 |
| 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.34 | ug/L | | | 10/04/17 02:21 | 1 |
| cis-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.30 | ug/L | | | 10/04/17 02:21 | 1 |
| Bromochloromethane | 1.0 | U | 1.0 | 0.36 | ug/L | | | 10/04/17 02:21 | 1 |
| 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 | ug/L | | | 10/04/17 02:21 | 1 |
| Chloroform | 1.0 | U | 1.0 | 0.27 | ug/L | | | 10/04/17 02:21 | 1 |
| 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.27 | ug/L | | | 10/04/17 02:21 | 1 |
| Carbon tetrachloride | 1.0 | U | 1.0 | 0.56 | ug/L | | | 10/04/17 02:21 | 1 |
| Benzene | 1.0 | U | 1.0 | 0.18 | ug/L | | | 10/04/17 02:21 | 1 |
| 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.24 | ug/L | | | 10/04/17 02:21 | 1 |
| Trichloroethene | 1.0 | U | 1.0 | 0.20 | ug/L | | | 10/04/17 02:21 | 1 |
| 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.35 | ug/L | | | 10/04/17 02:21 | 1 |
| Bromodichloromethane | 1.0 | U | 1.0 | 0.57 | ug/L | | | 10/04/17 02:21 | 1 |
| cis-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.32 | ug/L | | | 10/04/17 02:21 | 1 |
| 4-Methyl-2-pentanone (MIBK) | 5.0 | U | 5.0 | 2.2 | ug/L | | | 10/04/17 02:21 | 1 |
| Toluene | 1.0 | U | 1.0 | 0.16 | ug/L | | | 10/04/17 02:21 | 1 |
| trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.22 | ug/L | | | 10/04/17 02:21 | 1 |
| 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.31 | ug/L | | | 10/04/17 02:21 | 1 |
| Tetrachloroethene | 1.0 | U | 1.0 | 0.24 | ug/L | | | 10/04/17 02:21 | 1 |
| 2-Hexanone | 5.0 | U | 5.0 | 2.0 | ug/L | | | 10/04/17 02:21 | 1 |
| Dibromochloromethane | 1.0 | U | 1.0 | 0.44 | ug/L | | | 10/04/17 02:21 | 1 |
| 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.51 | ug/L | | | 10/04/17 02:21 | 1 |
| Chlorobenzene | 1.0 | U | 1.0 | 0.15 | ug/L | | | 10/04/17 02:21 | 1 |
| 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.49 | ug/L | | | 10/04/17 02:21 | 1 |
| Ethylbenzene | 1.0 | U | 1.0 | 0.25 | ug/L | | | 10/04/17 02:21 | 1 |
| Xylenes, Total | 2.0 | U | 2.0 | 0.27 | ug/L | | | 10/04/17 02:21 | 1 |
| Styrene | 1.0 | U | 1.0 | 0.22 | ug/L | | | 10/04/17 02:21 | 1 |
| Bromoform | 1.0 | U | 1.0 | 0.76 | ug/L | | | 10/04/17 02:21 | 1 |
| 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.37 | ug/L | | | 10/04/17 02:21 | 1 |
| Acrylonitrile | 20 | U | 20 | 3.3 | ug/L | | | 10/04/17 02:21 | 1 |
| 1,4-Dioxane | 200 | U | 200 | 16 | ug/L | | | 10/04/17 02:21 | 1 |

| Surrogate | MB | MB | Limits | Prepared | Analyzed | Dil Fac |
|------------------------------|-----------|-----------|----------|----------|----------------|---------|
| | %Recovery | Qualifier | | | | |
| 1,2-Dichloroethane-d4 (Surr) | 107 | | 65 - 121 | | 10/04/17 02:21 | 1 |
| Toluene-d8 (Surr) | 98 | | 73 - 120 | | 10/04/17 02:21 | 1 |
| 4-Bromofluorobenzene (Surr) | 93 | | 80 - 120 | | 10/04/17 02:21 | 1 |
| Dibromofluoromethane (Surr) | 94 | | 73 - 120 | | 10/04/17 02:21 | 1 |

TestAmerica Pittsburgh

QC Sample Results

Client: Groundwater Sciences Corporation
Project/Site: Harley Davidson

TestAmerica Job ID: 180-70873-1

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

Lab Sample ID: LCS 180-224792/4

Matrix: Water

Analysis Batch: 224792

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 | 11.6 | | ug/L | | 116 | 49 - 135 |
| Vinyl chloride | 10.0 | 11.8 | | ug/L | | 118 | 52 - 136 |
| Bromomethane | 10.0 | 10.1 | | ug/L | | 101 | 37 - 150 |
| Chloroethane | 10.0 | 13.5 | | ug/L | | 135 | 44 - 139 |
| 1,1-Dichloroethene | 10.0 | 10.3 | | ug/L | | 103 | 64 - 131 |
| Acetone | 20.0 | 22.4 | | ug/L | | 112 | 24 - 150 |
| Carbon disulfide | 10.0 | 9.37 | | ug/L | | 94 | 20 - 150 |
| Methylene Chloride | 10.0 | 9.77 | | ug/L | | 98 | 66 - 123 |
| trans-1,2-Dichloroethene | 10.0 | 10.0 | | ug/L | | 100 | 70 - 123 |
| Methyl tert-butyl ether | 10.0 | 9.62 | | ug/L | | 96 | 66 - 130 |
| 1,1-Dichloroethane | 10.0 | 9.85 | | ug/L | | 98 | 66 - 122 |
| cis-1,2-Dichloroethene | 10.0 | 9.81 | | ug/L | | 98 | 73 - 120 |
| Bromochloromethane | 10.0 | 9.53 | | ug/L | | 95 | 73 - 122 |
| 2-Butanone (MEK) | 20.0 | 21.0 | | ug/L | | 105 | 37 - 150 |
| Chloroform | 10.0 | 9.83 | | ug/L | | 98 | 72 - 123 |
| 1,1,1-Trichloroethane | 10.0 | 10.4 | | ug/L | | 104 | 66 - 129 |
| Carbon tetrachloride | 10.0 | 9.98 | | ug/L | | 100 | 58 - 145 |
| Benzene | 10.0 | 9.68 | | ug/L | | 97 | 75 - 123 |
| 1,2-Dichloroethane | 10.0 | 10.6 | | ug/L | | 106 | 63 - 130 |
| Trichloroethene | 10.0 | 9.28 | | ug/L | | 93 | 74 - 121 |
| 1,2-Dichloropropane | 10.0 | 9.20 | | ug/L | | 92 | 67 - 119 |
| Bromodichloromethane | 10.0 | 9.11 | | ug/L | | 91 | 62 - 127 |
| cis-1,3-Dichloropropene | 10.0 | 8.97 | | ug/L | | 90 | 61 - 127 |
| 4-Methyl-2-pentanone (MIBK) | 20.0 | 21.2 | | ug/L | | 106 | 41 - 135 |
| Toluene | 10.0 | 10.9 | | ug/L | | 109 | 76 - 129 |
| trans-1,3-Dichloropropene | 10.0 | 9.70 | | ug/L | | 97 | 61 - 136 |
| 1,1,2-Trichloroethane | 10.0 | 10.8 | | ug/L | | 108 | 74 - 126 |
| Tetrachloroethene | 10.0 | 10.5 | | ug/L | | 105 | 76 - 128 |
| 2-Hexanone | 20.0 | 21.2 | | ug/L | | 106 | 37 - 150 |
| Dibromochloromethane | 10.0 | 9.60 | | ug/L | | 96 | 63 - 131 |
| 1,2-Dibromoethane (EDB) | 10.0 | 9.86 | | ug/L | | 99 | 76 - 128 |
| Chlorobenzene | 10.0 | 10.3 | | ug/L | | 103 | 79 - 124 |
| 1,1,1,2-Tetrachloroethane | 10.0 | 10.6 | | ug/L | | 106 | 70 - 130 |
| Ethylbenzene | 10.0 | 9.83 | | ug/L | | 98 | 77 - 124 |
| Xylenes, Total | 20.0 | 20.2 | | ug/L | | 101 | 76 - 124 |
| Styrene | 10.0 | 10.0 | | ug/L | | 100 | 80 - 125 |
| Bromoform | 10.0 | 8.98 | | ug/L | | 90 | 54 - 136 |
| 1,1,2,2-Tetrachloroethane | 10.0 | 10.4 | | ug/L | | 104 | 72 - 128 |
| Acrylonitrile | 100 | 108 | | ug/L | | 108 | 60 - 130 |
| 1,4-Dioxane | 200 | 271 | | ug/L | | 135 | 26 - 150 |

| Surrogate | LCS %Recovery | LCS Qualifier | Limits |
|------------------------------|---------------|---------------|----------|
| 1,2-Dichloroethane-d4 (Surr) | 106 | | 65 - 121 |
| Toluene-d8 (Surr) | 113 | | 73 - 120 |
| 4-Bromofluorobenzene (Surr) | 107 | | 80 - 120 |
| Dibromofluoromethane (Surr) | 98 | | 73 - 120 |

QC Sample Results

Client: Groundwater Sciences Corporation
 Project/Site: Harley Davidson

TestAmerica Job ID: 180-70873-1

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

Lab Sample ID: MB 180-224919/7
Matrix: Water
Analysis Batch: 224919

Client Sample ID: Method Blank
Prep Type: Total/NA

| Analyte | MB Result | MB Qualifier | RL | MDL | Unit | D | Prepared | Analyzed | Dil Fac |
|-----------------------------|-----------|--------------|-----|------|------|---|----------|----------------|---------|
| Chloromethane | 1.0 | U | 1.0 | 0.38 | ug/L | | | 10/05/17 02:41 | 1 |
| Vinyl chloride | 1.0 | U | 1.0 | 0.17 | ug/L | | | 10/05/17 02:41 | 1 |
| Bromomethane | 1.0 | U | 1.0 | 0.59 | ug/L | | | 10/05/17 02:41 | 1 |
| Chloroethane | 1.0 | U | 1.0 | 0.58 | ug/L | | | 10/05/17 02:41 | 1 |
| 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.32 | ug/L | | | 10/05/17 02:41 | 1 |
| Acetone | 5.0 | U | 5.0 | 3.1 | ug/L | | | 10/05/17 02:41 | 1 |
| Carbon disulfide | 1.0 | U | 1.0 | 0.53 | ug/L | | | 10/05/17 02:41 | 1 |
| Methylene Chloride | 1.0 | U | 1.0 | 0.94 | ug/L | | | 10/05/17 02:41 | 1 |
| trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.20 | ug/L | | | 10/05/17 02:41 | 1 |
| Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.20 | ug/L | | | 10/05/17 02:41 | 1 |
| 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.34 | ug/L | | | 10/05/17 02:41 | 1 |
| cis-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.30 | ug/L | | | 10/05/17 02:41 | 1 |
| Bromochloromethane | 1.0 | U | 1.0 | 0.36 | ug/L | | | 10/05/17 02:41 | 1 |
| 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 | ug/L | | | 10/05/17 02:41 | 1 |
| Chloroform | 1.0 | U | 1.0 | 0.27 | ug/L | | | 10/05/17 02:41 | 1 |
| 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.27 | ug/L | | | 10/05/17 02:41 | 1 |
| Carbon tetrachloride | 1.0 | U | 1.0 | 0.56 | ug/L | | | 10/05/17 02:41 | 1 |
| Benzene | 1.0 | U | 1.0 | 0.18 | ug/L | | | 10/05/17 02:41 | 1 |
| 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.24 | ug/L | | | 10/05/17 02:41 | 1 |
| Trichloroethene | 1.0 | U | 1.0 | 0.20 | ug/L | | | 10/05/17 02:41 | 1 |
| 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.35 | ug/L | | | 10/05/17 02:41 | 1 |
| Bromodichloromethane | 1.0 | U | 1.0 | 0.57 | ug/L | | | 10/05/17 02:41 | 1 |
| cis-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.32 | ug/L | | | 10/05/17 02:41 | 1 |
| 4-Methyl-2-pentanone (MIBK) | 5.0 | U | 5.0 | 2.2 | ug/L | | | 10/05/17 02:41 | 1 |
| Toluene | 1.0 | U | 1.0 | 0.16 | ug/L | | | 10/05/17 02:41 | 1 |
| trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.22 | ug/L | | | 10/05/17 02:41 | 1 |
| 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.31 | ug/L | | | 10/05/17 02:41 | 1 |
| Tetrachloroethene | 1.0 | U | 1.0 | 0.24 | ug/L | | | 10/05/17 02:41 | 1 |
| 2-Hexanone | 5.0 | U | 5.0 | 2.0 | ug/L | | | 10/05/17 02:41 | 1 |
| Dibromochloromethane | 1.0 | U | 1.0 | 0.44 | ug/L | | | 10/05/17 02:41 | 1 |
| 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.51 | ug/L | | | 10/05/17 02:41 | 1 |
| Chlorobenzene | 1.0 | U | 1.0 | 0.15 | ug/L | | | 10/05/17 02:41 | 1 |
| 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.49 | ug/L | | | 10/05/17 02:41 | 1 |
| Ethylbenzene | 1.0 | U | 1.0 | 0.25 | ug/L | | | 10/05/17 02:41 | 1 |
| Xylenes, Total | 2.0 | U | 2.0 | 0.27 | ug/L | | | 10/05/17 02:41 | 1 |
| Styrene | 1.0 | U | 1.0 | 0.22 | ug/L | | | 10/05/17 02:41 | 1 |
| Bromoform | 1.0 | U | 1.0 | 0.76 | ug/L | | | 10/05/17 02:41 | 1 |
| 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.37 | ug/L | | | 10/05/17 02:41 | 1 |
| Acrylonitrile | 20 | U | 20 | 3.3 | ug/L | | | 10/05/17 02:41 | 1 |
| 1,4-Dioxane | 200 | U | 200 | 16 | ug/L | | | 10/05/17 02:41 | 1 |

| Surrogate | MB %Recovery | MB Qualifier | Limits | Prepared | Analyzed | Dil Fac |
|------------------------------|--------------|--------------|----------|----------|----------------|---------|
| 1,2-Dichloroethane-d4 (Surr) | 111 | | 65 - 121 | | 10/05/17 02:41 | 1 |
| Toluene-d8 (Surr) | 101 | | 73 - 120 | | 10/05/17 02:41 | 1 |
| 4-Bromofluorobenzene (Surr) | 95 | | 80 - 120 | | 10/05/17 02:41 | 1 |
| Dibromofluoromethane (Surr) | 96 | | 73 - 120 | | 10/05/17 02:41 | 1 |

QC Sample Results

Client: Groundwater Sciences Corporation
Project/Site: Harley Davidson

TestAmerica Job ID: 180-70873-1

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

Lab Sample ID: LCS 180-224919/4

Matrix: Water

Analysis Batch: 224919

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 | 12.0 | | ug/L | | 120 | 49 - 135 |
| Vinyl chloride | 10.0 | 12.6 | | ug/L | | 126 | 52 - 136 |
| Bromomethane | 10.0 | 12.4 | | ug/L | | 124 | 37 - 150 |
| Chloroethane | 10.0 | 12.5 | | ug/L | | 125 | 44 - 139 |
| 1,1-Dichloroethene | 10.0 | 9.95 | | ug/L | | 100 | 64 - 131 |
| Acetone | 20.0 | 21.7 | | ug/L | | 109 | 24 - 150 |
| Carbon disulfide | 10.0 | 9.04 | | ug/L | | 90 | 20 - 150 |
| Methylene Chloride | 10.0 | 8.90 | | ug/L | | 89 | 66 - 123 |
| trans-1,2-Dichloroethene | 10.0 | 9.52 | | ug/L | | 95 | 70 - 123 |
| Methyl tert-butyl ether | 10.0 | 9.23 | | ug/L | | 92 | 66 - 130 |
| 1,1-Dichloroethane | 10.0 | 9.51 | | ug/L | | 95 | 66 - 122 |
| cis-1,2-Dichloroethene | 10.0 | 8.65 | | ug/L | | 86 | 73 - 120 |
| Bromochloromethane | 10.0 | 8.75 | | ug/L | | 88 | 73 - 122 |
| 2-Butanone (MEK) | 20.0 | 18.3 | | ug/L | | 91 | 37 - 150 |
| Chloroform | 10.0 | 9.05 | | ug/L | | 91 | 72 - 123 |
| 1,1,1-Trichloroethane | 10.0 | 9.46 | | ug/L | | 95 | 66 - 129 |
| Carbon tetrachloride | 10.0 | 9.46 | | ug/L | | 95 | 58 - 145 |
| Benzene | 10.0 | 8.85 | | ug/L | | 88 | 75 - 123 |
| 1,2-Dichloroethane | 10.0 | 9.68 | | ug/L | | 97 | 63 - 130 |
| Trichloroethene | 10.0 | 8.20 | | ug/L | | 82 | 74 - 121 |
| 1,2-Dichloropropane | 10.0 | 8.44 | | ug/L | | 84 | 67 - 119 |
| Bromodichloromethane | 10.0 | 8.21 | | ug/L | | 82 | 62 - 127 |
| cis-1,3-Dichloropropene | 10.0 | 8.18 | | ug/L | | 82 | 61 - 127 |
| 4-Methyl-2-pentanone (MIBK) | 20.0 | 20.4 | | ug/L | | 102 | 41 - 135 |
| Toluene | 10.0 | 10.4 | | ug/L | | 104 | 76 - 129 |
| trans-1,3-Dichloropropene | 10.0 | 9.78 | | ug/L | | 98 | 61 - 136 |
| 1,1,2-Trichloroethane | 10.0 | 10.6 | | ug/L | | 106 | 74 - 126 |
| Tetrachloroethene | 10.0 | 9.77 | | ug/L | | 98 | 76 - 128 |
| 2-Hexanone | 20.0 | 19.4 | | ug/L | | 97 | 37 - 150 |
| Dibromochloromethane | 10.0 | 9.57 | | ug/L | | 96 | 63 - 131 |
| 1,2-Dibromoethane (EDB) | 10.0 | 9.26 | | ug/L | | 93 | 76 - 128 |
| Chlorobenzene | 10.0 | 9.72 | | ug/L | | 97 | 79 - 124 |
| 1,1,1,2-Tetrachloroethane | 10.0 | 9.84 | | ug/L | | 98 | 70 - 130 |
| Ethylbenzene | 10.0 | 9.31 | | ug/L | | 93 | 77 - 124 |
| Xylenes, Total | 20.0 | 18.9 | | ug/L | | 95 | 76 - 124 |
| Styrene | 10.0 | 9.11 | | ug/L | | 91 | 80 - 125 |
| Bromoform | 10.0 | 8.75 | | ug/L | | 88 | 54 - 136 |
| 1,1,2,2-Tetrachloroethane | 10.0 | 9.67 | | ug/L | | 97 | 72 - 128 |
| Acrylonitrile | 100 | 102 | | ug/L | | 102 | 60 - 130 |
| 1,4-Dioxane | 200 | 189 | J | ug/L | | 94 | 26 - 150 |

| Surrogate | LCS %Recovery | LCS Qualifier | Limits |
|------------------------------|---------------|---------------|----------|
| 1,2-Dichloroethane-d4 (Surr) | 102 | | 65 - 121 |
| Toluene-d8 (Surr) | 110 | | 73 - 120 |
| 4-Bromofluorobenzene (Surr) | 102 | | 80 - 120 |
| Dibromofluoromethane (Surr) | 90 | | 73 - 120 |

QC Association Summary

Client: Groundwater Sciences Corporation
Project/Site: Harley Davidson

TestAmerica Job ID: 180-70873-1

GC/MS VOA

Analysis Batch: 224792

| Lab Sample ID | Client Sample ID | Prep Type | Matrix | Method | Prep Batch |
|------------------|---------------------|-----------|--------|--------|------------|
| 180-70873-1 | HD-SPBA-CW-22-0/1-0 | Total/NA | Water | 8260C | |
| 180-70873-2 | HD-QC4-0/1-2 | Total/NA | Water | 8260C | |
| MB 180-224792/6 | Method Blank | Total/NA | Water | 8260C | |
| LCS 180-224792/4 | Lab Control Sample | Total/NA | Water | 8260C | |

Analysis Batch: 224919

| Lab Sample ID | Client Sample ID | Prep Type | Matrix | Method | Prep Batch |
|------------------|---------------------|-----------|--------|--------|------------|
| 180-70873-1 - DL | HD-SPBA-CW-22-0/1-0 | Total/NA | Water | 8260C | |
| MB 180-224919/7 | Method Blank | Total/NA | Water | 8260C | |
| LCS 180-224919/4 | Lab Control Sample | Total/NA | Water | 8260C | |

Lab Chronicle

Client: Groundwater Sciences Corporation
Project/Site: Harley Davidson

TestAmerica Job ID: 180-70873-1

Client Sample ID: HD-SPBA-CW-22-0/1-0

Lab Sample ID: 180-70873-1

Date Collected: 09/29/17 09:40

Matrix: Water

Date Received: 10/03/17 09:00

| 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 | 224792 | 10/04/17 05:23 | FBB | TAL PIT |
| | | Instrument ID: CHHP5 | | | | | | | | |
| Total/NA | Analysis | 8260C | DL | 12.5 | 5 mL | 5 mL | 224919 | 10/05/17 06:31 | FBB | TAL PIT |
| | | Instrument ID: CHHP5 | | | | | | | | |

Client Sample ID: HD-QC4-0/1-2

Lab Sample ID: 180-70873-2

Date Collected: 09/29/17 12:00

Matrix: Water

Date Received: 10/03/17 09:00

| 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 | 224792 | 10/04/17 04:59 | FBB | TAL PIT |
| | | Instrument ID: CHHP5 | | | | | | | | |

Laboratory References:

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

Analyst References:

Lab: TAL PIT

Batch Type: Analysis

FBB = Frank Bungard

Accreditation/Certification Summary

Client: Groundwater Sciences Corporation
Project/Site: Harley Davidson

TestAmerica Job ID: 180-70873-1

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: 180-70873-1

| 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: 180-70873-1

| Lab Sample ID | Client Sample ID | Matrix | Collected | Received |
|----------------------|-------------------------|---------------|------------------|-----------------|
| 180-70873-1 | HD-SPBA-CW-22-0/1-0 | Water | 09/29/17 09:40 | 10/03/17 09:00 |
| 180-70873-2 | HD-QC4-0/1-2 | Water | 09/29/17 12:00 | 10/03/17 09:00 |

GC/MS VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1

SDG No.: _____

Instrument ID: CHHP5 Analysis Batch Number: 218218Lab Sample ID: IC 180-218218/2 Client Sample ID: _____Date Analyzed: 07/27/17 00:51 Lab File ID: 50727D02.D GC Column: DB-624 ID: 0.18 (mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|------------------------|----------------|---------------------|----------|----------------|
| | | REASON | ANALYST | DATE |
| Trichlorofluoromethane | 2.76 | Poor chromatography | bungardf | 07/27/17 03:06 |

Lab Sample ID: IC 180-218218/3 Client Sample ID: _____Date Analyzed: 07/27/17 01:15 Lab File ID: 50727D03.D GC Column: DB-624 ID: 0.18 (mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|------------------------|----------------|---------------------|----------|----------------|
| | | REASON | ANALYST | DATE |
| Trichlorofluoromethane | 2.75 | Poor chromatography | bungardf | 07/27/17 03:13 |
| 1,4-Dioxane | 8.05 | Poor chromatography | bungardf | 07/27/17 03:14 |

Lab Sample ID: ICIS 180-218218/4 Client Sample ID: _____Date Analyzed: 07/27/17 01:39 Lab File ID: 50727D04.D GC Column: DB-624 ID: 0.18 (mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|------------------------|----------------|---------------------|----------|----------------|
| | | REASON | ANALYST | DATE |
| Trichlorofluoromethane | 2.75 | Poor chromatography | bungardf | 07/27/17 03:15 |
| 1,4-Dioxane | 8.05 | Poor chromatography | bungardf | 07/27/17 03:15 |

GC/MS VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1

SDG No.: _____

Instrument ID: CHHP5 Analysis Batch Number: 224792

Lab Sample ID: CCVIS 180-224792/2 Client Sample ID: _____

Date Analyzed: 10/04/17 00:22 Lab File ID: 51003D02.D GC Column: DB-624 ID: 0.18 (mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---------------------|----------------|---------------------|----------|----------------|
| | | REASON | ANALYST | DATE |
| 2,2-Dichloropropane | 6.01 | Poor chromatography | bungardf | 10/04/17 00:54 |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-70873-1

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_00267 | 10/09/17 | 10/02/17 | Methanol, Lot 2469119 | 10 mL | VOA8260GAS2ND_00211 | 100 uL | Bromomethane | 25 ug/mL |
| | | | | | | | Chloroethane | 25 ug/mL |
| | | | | | | | Chloromethane | 25 ug/mL |
| | | | | | | | Vinyl chloride | 25 ug/mL |
| | | | | | VOA8260VOA2ND_00263 | 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.: 180-70873-1

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_00211 | 06/30/20 | | Restek, Lot A0128832 | | | (Purchased Reagent) | Bromomethane | 2500 ug/mL |
| | | | | | | | Chloroethane | 2500 ug/mL |
| | | | | | | | Chloromethane | 2500 ug/mL |
| | | | | | | | Vinyl chloride | 2500 ug/mL |
| .VOA8260VOA2ND_00263 | 10/09/17 | 09/09/17 | Methanol, Lot 2469125 | 10 mL | VOA8260MEGA2_00062 | 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.: 180-70873-1

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_00062 | 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.: 180-70873-1

SDG No.: _____

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|------------|----------|-----------|---------------|----------------------|---------------------|--------------|---------------------------------------|---------------|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Trichlorofluoromethane | 25 ug/mL |
| | | | | | | | Vinyl chloride | 25 ug/mL |
| | | | | | VOA8260VOAPRI_00260 | 1 mL | 2-Butanone (MEK) | 25 ug/mL |
| | | | | | | | 2-Hexanone | 25 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 25 ug/mL |
| | | | | | | | Acetone | 25 ug/mL |
| | | | | | | | 1,1,1,2-Tetrachloroethane | 25 ug/mL |
| | | | | | | | 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.: 180-70873-1

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 |
| | | | | | | | 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 |
| .VOA8260VOAPRI_00260 | 08/06/17 | 07/06/17 | Methanol, Lot 2019056 | 10 mL | VOA8260KET1ST_00100 | 0.2 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-trifluoroethane | 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.: 180-70873-1

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.: 180-70873-1

SDG No.: _____

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|-----------------------|----------|-----------|----------------------|----------------------|---------------------|--------------|---------------------------------------|---------------|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 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 |
| ..VOA8260KET1ST_00100 | 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 |
| ..VOA8260MEGA1_00065 | 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-Trichloro-1,2,2-trifluoroethane | 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.: 180-70873-1

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 |
| VOABFB25_00090 | | | | | | | 1,2-Dichloroethene, Total | |
| | | | | | | | 1,3-Dichloropropene, Total | |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-70873-1

SDG No.: _____

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|-----------------------------|----------|-----------|-----------------------|----------------------|-------------------|---------------------|--|---------------|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Tentatively Identified Compound | |
| | | | | | | | Total BTEX | |
| | | | | | | | Xylenes, Total | |
| .VOABFB50_00093 | 08/10/17 | 07/10/17 | Methanol, Lot 2019056 | 50 mL | VOABFB50_00093 | 5 mL | BFB | 25 ug/mL |
| ..VOABFBRES_00058 | 11/30/21 | | Restek, Lot A0122647 | | VOABFBRES_00058 | 1 mL | BFB | 50 ug/mL |
| | | | | | | (Purchased Reagent) | BFB | 2500 ug/mL |
| VOABFB25_00093 | | | | | | | 1,2-Dichloroethene, Total | |
| | | | | | | | 1,3-Dichloropropene, Total | |
| | | | | | | | Tentatively Identified Compound | |
| | | | | | | | Total BTEX | |
| | | | | | | | Xylenes, Total | |
| .VOABFB50_00095 | 10/09/17 | 09/09/17 | Methanol, Lot 2469125 | 50 mL | VOABFB50_00095 | 5 mL | BFB | 25 ug/mL |
| ..VOABFBRES_00056 | 11/30/21 | | Restek, Lot A0122647 | | VOABFBRES_00056 | 1 mL | BFB | 50 ug/mL |
| | | | | | | (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 |
| | | | | | | | 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 |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-70873-1

SDG No.: _____

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|----------------------------|----------|-----------|-----------------------|----------------------|---------------------|--------------|-----------------------------|---------------|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Chlorobenzotrifluoride | 5000 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_00017 | 07/31/17 | 07/24/16 | Methanol, Lot 2019067 | 25 mL | VOA8260VARES_00083 | 125 uL | Vinyl acetate | 25 ug/mL |
| .VOA8260VARES_00083 | 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

CERTIFIED VALUES

| Elution Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L.; K=2) | | | |
|---------------|----------------------------------|-----------------------------|--------------------------------------|-------|-------------|-------|
| | | | Value | Unit | Method | Notes |
| 1 | Dichlorodifluoromethane (CFC-12) | 2,500.5 µg/mL | +/- 16.7232 | µg/mL | Gravimetric | |
| | CAS # 75-71-8 (Lot Q167-08) | | +/- 140.4412 | µg/mL | Unstressed | |
| | Purity 99% | | +/- 143.7161 | µg/mL | Stressed | |
| 2 | Chloromethane (methyl chloride) | 2,498.7 µg/mL | +/- 17.4998 | µg/mL | Gravimetric | |
| | CAS # 74-87-3 (Lot SHBG7976V) | | +/- 140.4406 | µg/mL | Unstressed | |
| | Purity 99% | | +/- 143.7111 | µg/mL | Stressed | |
| 3 | Vinyl chloride | 2,498.4 µg/mL | +/- 16.6753 | µg/mL | Gravimetric | |
| | CAS # 75-01-4 (Lot 1026101231B1) | | +/- 140.3203 | µg/mL | Unstressed | |
| | Purity 99% | | +/- 143.5926 | µg/mL | Stressed | |
| 4 | 1,3-Butadiene | 2,496.9 µg/mL | +/- 17.0619 | µg/mL | Gravimetric | |
| | CAS # 106-99-0 (Lot SHBF3387V) | | +/- 140.2843 | µg/mL | Unstressed | |
| | Purity 99% | | +/- 143.5535 | µg/mL | Stressed | |
| 5 | Bromomethane (methyl bromide) | 2,500.5 µg/mL | +/- 17.3456 | µg/mL | Gravimetric | |
| | CAS # 74-83-9 (Lot 101604) | | +/- 140.5211 | µg/mL | Unstressed | |
| | Purity 99% | | +/- 143.7944 | µg/mL | Stressed | |
| 6 | Chloroethane (ethyl chloride) | 2,500.5 µg/mL | +/- 16.8189 | µg/mL | Gravimetric | |
| | CAS # 75-00-3 (Lot 23593) | | +/- 140.4526 | µg/mL | Unstressed | |
| | Purity 99% | | +/- 143.7272 | µg/mL | Stressed | |
| 7 | Dichlorofluoromethane (CFC-21) | 2,500.0 µg/mL | +/- 10.0499 | µg/mL | Gravimetric | |
| | CAS # 75-43-4 (Lot 4938100) | | +/- 139.7786 | µg/mL | Unstressed | |
| | Purity 99% | | +/- 143.0675 | µg/mL | Stressed | |

| | | | | | | |
|---|---------------------------------|--------------------------|-----|----------|------------------|-------------|
| 8 | Trichlorofluoromethane (CFC-11) | 2,501.5 $\mu\text{g/mL}$ | +/- | 16.5404 | $\mu\text{g/mL}$ | Gravimetric |
| | CAS # 75-69-4 (Lot SHBG7531V) | | +/- | 140.4793 | $\mu\text{g/mL}$ | Unstressed |
| | Purity 99% | | +/- | 143.7562 | $\mu\text{g/mL}$ | Stressed |

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

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

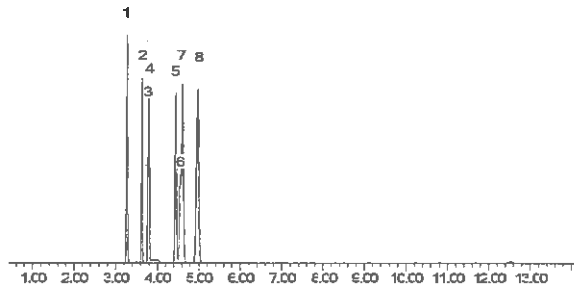
Carrier Gas:
 helium-constant flow 2.0 mL/min.

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 Jaglowski
 Joseph Jaglowski - Mix Technician

Date Mixed: 17-Jan-2017

Balance: 1125113331

Jennifer J Pollino
 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_00211



CERTIFIED REFERENCE MATERIAL

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

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Certificate of Analysis



FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

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

Catalog No.: 569722.SEC Lot No.: A0128832
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: June 30, 2020 Storage: 0°C or colder

CERTIFIED VALUES

Table with 7 rows and 7 columns: Elution Order, Compound, Grav. Conc. (weight/volume), Expanded Uncertainty (95% C.L.; K=2), and three additional columns for uncertainty components. Rows include Dichlorodifluoromethane (CFC-12), Chloromethane (methyl chloride), Vinyl chloride, 1,3-Butadiene, Bromomethane (methyl bromide), Chloroethane (ethyl chloride), and Dichlorofluoromethane (CFC-21).

General Certified Reference Material Notes

Expiration Notes:

- Expiration date valid for unopened ampul stored in compliance with the recommended conditions.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k \sqrt{U_{gravimetric}^2 + U_{homogeneity}^2 + U_{storage\ stability}^2 + U_{shipping\ stability}^2}$$

k is a coverage factor of 2, which gives a level of confidence of approximately 95%.

- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

| Label Conditions | Standard Conditions | Non-Standard Conditions |
|---------------------------------|---------------------|-------------------------|
| 25°C Nominal (Room Temperature) | < 60°C | ≥ 60°C up to 7 days |
| 10°C or colder (Refrigerate) | < 40°C | ≥ 40°C up to 7 days |
| 0°C or colder (Freezer) | < 25°C | ≥ 25°C up to 7 days |

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31861, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

Reagent

VOA8260INTRES_00123



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. : 568718 Lot No.: A0113246

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

CERTIFIED VALUES

| 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% (Lot I201P18) | 5,000.4 µg/mL | +/- | 29.0712 | µg/mL Gravimetric |
| | | | +/- | 106.0450 | µg/mL Unstressed |
| | | | +/- | 106.5155 | µg/mL Stressed |
| 2 | 2-Butanone-d5 CAS # 24313-50-6 Purity 99% (Lot M276P24) | 1,250.2 µg/mL | +/- | 7.2688 | µg/mL Gravimetric |
| | | | +/- | 26.5135 | µg/mL Unstressed |
| | | | +/- | 26.6311 | µg/mL Stressed |
| 3 | Fluorobenzene CAS # 462-06-6 Purity 99% (Lot BCBK8171V) | 250.2 µg/mL | +/- | 1.4580 | µg/mL Gravimetric |
| | | | +/- | 5.3070 | µg/mL Unstressed |
| | | | +/- | 5.3305 | µg/mL Stressed |
| 4 | 1,4-Dioxane-d8 CAS # 17647-74-4 Purity 98% (Lot I-19073) | 5,000.6 µg/mL | +/- | 29.0727 | µg/mL Gravimetric |
| | | | +/- | 106.0502 | µg/mL Unstressed |
| | | | +/- | 106.5208 | µg/mL Stressed |
| 5 | Chlorobenzene-d5 CAS # 3114-55-4 Purity 99% (Lot PR-23926) | 250.4 µg/mL | +/- | 1.4592 | µg/mL Gravimetric |
| | | | +/- | 5.3113 | µg/mL Unstressed |
| | | | +/- | 5.3348 | µg/mL Stressed |
| 6 | 1,4-Dichlorobenzene-d4 CAS # 3855-82-1 Purity 99% (Lot PR-18488) | 250.0 µg/mL | +/- | 1.4569 | µg/mL Gravimetric |
| | | | +/- | 5.3028 | µg/mL Unstressed |
| | | | +/- | 5.3263 | µg/mL Stressed |

Reagent

VOA8260INTRES_00135



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. : 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

CERTIFIED VALUES

| 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% (Lot I-201) | 5,050.0 µg/mL | +/- | 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% (Lot M-276) | 1,262.5 µg/mL | +/- | 7.3403 | µg/mL | Gravimetric |
| | | | +/- | 27.0303 | µg/mL | Unstressed |
| | | | +/- | 27.8161 | µg/mL | Stressed |
| 3 | Fluorobenzene CAS # 462-06-6 Purity 99% (Lot BCBK8171V) | 251.6 µg/mL | +/- | 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% (Lot I-19942) | 5,048.8 µg/mL | +/- | 29.3526 | µg/mL | Gravimetric |
| | | | +/- | 108.0950 | µg/mL | Unstressed |
| | | | +/- | 111.2375 | µg/mL | Stressed |
| 5 | Chlorobenzene-d5 CAS # 3114-55-4 Purity 99% (Lot PR-23926) | 251.5 µg/mL | +/- | 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% (Lot PR-18488) | 252.5 µg/mL | +/- | 1.4714 | µg/mL | Gravimetric |
| | | | +/- | 5.4070 | µg/mL | Unstressed |
| | | | +/- | 5.5641 | µg/mL | Stressed |

Reagent

VOA8260KET1ST_00099



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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. : 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

CERTIFIED VALUES

| Elution Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L., K=2) | | | |
|---------------|-----------------------------|-----------------------------------|--------------------------------------|----------|-------|-------------|
| 1 | Acetone | 12,517.5 µg/mL (Lot SHBH0922V) | +/- | 72.7778 | µg/mL | Gravimetric |
| | CAS # 67-64-1 | | +/- | 755.2362 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 757.0293 | µg/mL | Stressed |
| 2 | 2-Butanone (MEK) | 12,521.8 µg/mL (Lot SHBF2461V) | +/- | 72.8025 | µg/mL | Gravimetric |
| | CAS # 78-93-3 | | +/- | 755.4927 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 757.2863 | µg/mL | Stressed |
| 3 | 4-Methyl-2-pentanone (MIBK) | 12,519.8 µg/mL (Lot SHBG3630V) | +/- | 72.7909 | µg/mL | Gravimetric |
| | CAS # 108-10-1 | | +/- | 755.3720 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 757.1654 | µg/mL | Stressed |
| 4 | 2-Hexanone | 12,508.5 µg/mL (Lot MKBW0198V) | +/- | 72.7255 | µg/mL | Gravimetric |
| | CAS # 591-78-6 | | +/- | 754.6932 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 756.4850 | µg/mL | Stressed |

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

Reagent

VOA8260KET1ST_00100



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. : 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

CERTIFIED VALUES

| Elution Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L., K=2) | | | |
|---------------|-----------------------------|-----------------------------------|--------------------------------------|----------|-------|-------------|
| 1 | Acetone | 12,517.5 µg/mL (Lot SHBH0922V) | +/- | 72.7778 | µg/mL | Gravimetric |
| | CAS # 67-64-1 | | +/- | 755.2362 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 757.0293 | µg/mL | Stressed |
| 2 | 2-Butanone (MEK) | 12,521.8 µg/mL (Lot SHBF2461V) | +/- | 72.8025 | µg/mL | Gravimetric |
| | CAS # 78-93-3 | | +/- | 755.4927 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 757.2863 | µg/mL | Stressed |
| 3 | 4-Methyl-2-pentanone (MIBK) | 12,519.8 µg/mL (Lot SHBG3630V) | +/- | 72.7909 | µg/mL | Gravimetric |
| | CAS # 108-10-1 | | +/- | 755.3720 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 757.1654 | µg/mL | Stressed |
| 4 | 2-Hexanone | 12,508.5 µg/mL (Lot MKBW0198V) | +/- | 72.7255 | µg/mL | Gravimetric |
| | CAS # 591-78-6 | | +/- | 754.6932 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 756.4850 | µg/mL | Stressed |

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

Reagent

VOA8260KET1ST_00102



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. : 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

CERTIFIED VALUES

| Elution Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L.; K=2) | | | |
|---------------|--------------------------------|-----------------------------|--------------------------------------|----------|-------|-------------|
| 1 | Acetone | 12,517.5 µg/mL | +/- | 72.7778 | µg/mL | Gravimetric |
| | CAS # 67-64-1 (Lot SHBH0922V) | | +/- | 755.2362 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 757.0293 | µg/mL | Stressed |
| 2 | 2-Butanone (MEK) | 12,521.8 µg/mL | +/- | 72.8025 | µg/mL | Gravimetric |
| | CAS # 78-93-3 (Lot SHBF2461V) | | +/- | 755.4927 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 757.2863 | µg/mL | Stressed |
| 3 | 4-Methyl-2-pentanone (MIBK) | 12,519.8 µg/mL | +/- | 72.7909 | µg/mL | Gravimetric |
| | CAS # 108-10-1 (Lot SHBG3630V) | | +/- | 755.3720 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 757.1654 | µg/mL | Stressed |
| 4 | 2-Hexanone | 12,508.5 µg/mL | +/- | 72.7255 | µg/mL | Gravimetric |
| | CAS # 591-78-6 (Lot MKBW0198V) | | +/- | 754.6932 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 756.4850 | µg/mL | 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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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

CERTIFIED VALUES

| Elution Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L.; K=2) | | | |
|---------------|--|-----------------------------|--------------------------------------|------------------|-------------|--|
| 1 | Diethyl ether (ethyl ether) CAS # 60-29-7 (Lot SHBG1462V) Purity 99% | 2,501.3 µg/mL | +/- | 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 (Lot 00009482) Purity 99% | 2,505.1 µg/mL | +/- | 14.5650 µg/mL | Gravimetric | |
| | | | +/- | 151.1453 µg/mL | Unstressed | |
| | | | +/- | 151.5041 µg/mL | Stressed | |
| 3 | 1,1-dichloroethene CAS # 75-35-4 (Lot SHBG8609V) Purity 99% | 2,511.5 µg/mL | +/- | 14.6021 µg/mL | Gravimetric | |
| | | | +/- | 151.5299 µg/mL | Unstressed | |
| | | | +/- | 151.8897 µg/mL | Stressed | |
| 4 | tert-Butanol (TBA) CAS # 75-65-0 (Lot SHBF0688V) Purity 99% | 25,001.8 µg/mL | +/- | 145.3547 µg/mL | Gravimetric | |
| | | | +/- | 1,508.4656 µg/mL | Unstressed | |
| | | | +/- | 1,512.0470 µg/mL | Stressed | |
| 5 | Methyl acetate CAS # 79-20-9 (Lot SHBG4345V) Purity 99% | 5,000.5 µg/mL | +/- | 29.0733 µg/mL | Gravimetric | |
| | | | +/- | 301.7023 µg/mL | Unstressed | |
| | | | +/- | 302.4186 µg/mL | Stressed | |
| 6 | Iodomethane (methyl iodide) CAS # 74-88-4 (Lot SHBF2149V) Purity 99% | 2,502.9 µg/mL | +/- | 14.5519 µg/mL | Gravimetric | |
| | | | +/- | 151.0095 µg/mL | Unstressed | |
| | | | +/- | 151.3681 µg/mL | Stressed | |
| 7 | Allyl chloride (3-chloropropene) CAS # 107-05-1 (Lot SHBF8133V) Purity 99% | 2,517.1 µg/mL | +/- | 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 | Trichloroethene 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 | Tetrachloroethene 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 +/- 150.8497 +/- 151.2078 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 41 | 1,2-Dibromoethane (EDB) CAS # 106-93-4 Purity 99% | (Lot BCBH3877V) | 2,501.3 µg/mL | +/- 14.5425 +/- 150.9115 +/- 151.2698 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 42 | Chlorobenzene CAS # 108-90-7 Purity 99% | (Lot SHBF0505V) | 2,500.1 µg/mL | +/- 14.5359 +/- 150.8436 +/- 151.2017 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 43 | m-Xylene CAS # 108-38-3 Purity 99% | (Lot SHBG4347V) | 1,250.3 µg/mL | +/- 7.2691 +/- 75.4331 +/- 75.6122 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 44 | p-Xylene CAS # 106-42-3 Purity 99% | (Lot SHBG3928V) | 1,251.3 µg/mL | +/- 7.2749 +/- 75.4935 +/- 75.6727 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 45 | Ethylbenzene CAS # 100-41-4 Purity 99% | (Lot SHBG5920V) | 2,503.3 µg/mL | +/- 14.5541 +/- 151.0322 +/- 151.3907 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 46 | 1,1,1,2-Tetrachloroethane CAS # 630-20-6 Purity 99% | (Lot MKBS3769V) | 2,500.3 µg/mL | +/- 14.5367 +/- 150.8512 +/- 151.2093 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 47 | o-Xylene CAS # 95-47-6 Purity 99% | (Lot SHBH3432V) | 2,504.9 µg/mL | +/- 14.5636 +/- 151.1302 +/- 151.4890 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 48 | Styrene CAS # 100-42-5 Purity 99% | (Lot MKBS7097V) | 2,506.3 µg/mL | +/- 14.5716 +/- 151.2132 +/- 151.5722 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 49 | Isopropylbenzene (cumene) CAS # 98-82-8 Purity 99% | (Lot 10185056) | 2,501.6 µg/mL | +/- 14.5447 +/- 150.9341 +/- 151.2925 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 50 | bromoform CAS # 75-25-2 Purity 99% | (Lot SHBD8459V) | 2,502.9 µg/mL | +/- 14.5519 +/- 151.0095 +/- 151.3681 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 51 | bromodichloromethane CAS # 75-27-4 Purity 97% | (Lot MKBW5506V) | 2,506.8 µg/mL | +/- 14.5750 +/- 151.2490 +/- 151.6081 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 52 | 1,1,2,2-Tetrachloroethane CAS # 79-34-5 Purity 99% | (Lot CFA4D) | 2,501.3 µg/mL | +/- 14.5425 +/- 150.9115 +/- 151.2698 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 53 | 1,2,3-Trichloropropane CAS # 96-18-4 Purity 99% | (Lot BCBH8722V) | 2,508.5 µg/mL | +/- 14.5846 +/- 151.3489 +/- 151.7082 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 54 | trans-1,4-dichloro-2-butene CAS # 110-57-6 Purity 95% | (Lot MKBP6041V) | 2,500.8 µg/mL | +/- 14.5396 +/- 150.8817 +/- 151.2399 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 55 | n-Propylbenzene CAS # 103-65-1 Purity 99% | (Lot MKBJ0332V) | 2,501.9 µg/mL | +/- 14.5461 +/- 150.9492 +/- 151.3076 | µg/mL µg/mL µg/mL | Gravimetric Unstressed 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:
60m x 0.25mm x 1.4µm
Rtx-502.2 (cat.#10916)

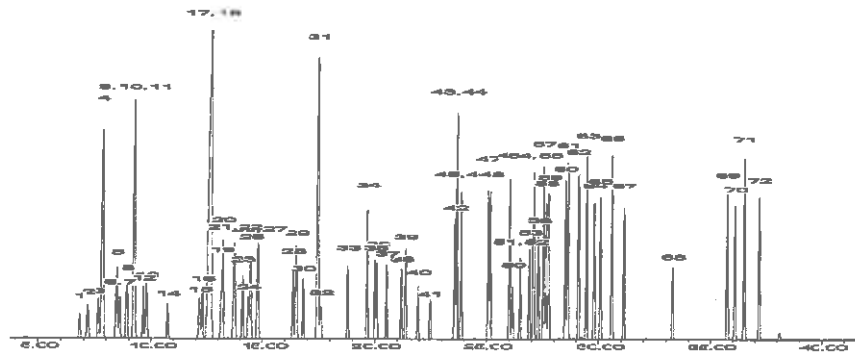
Carrier Gas:
helium-constant pressure 30 psi

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.

F. Joseph Tallon
F. Joseph Tallon - Mix Technician

Date Mixed: 22-Dec-2016 **Balance:** B251644995

Jennifer A. 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

VOA8260MEGA2_00062



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. : 571992.sec **Lot No.:** A0123775
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

CERTIFIED VALUES

| Elution Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L., K=2) | | | |
|---------------|--|-----------------------------|--------------------------------------|------------|-------|-------------|
| 1 | Diethyl ether (ethyl ether) | 2,501.2 µg/mL | +/- | 14.5422 | µg/mL | Gravimetric |
| | CAS # 60-29-7.SEC (Lot F23X068) | | +/- | 150.9088 | µg/mL | Unstressed |
| | Purity 98% | | +/- | 151.2671 | µg/mL | Stressed |
| 2 | 1,1,2-Trichlorotrifluoroethane (CFC-113) | 2,501.1 µg/mL | +/- | 14.5418 | µg/mL | Gravimetric |
| | CAS # 76-13-1.SEC (Lot 18342) | | +/- | 150.9040 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 151.2622 | µg/mL | Stressed |
| 3 | 1,1-Dichloroethene | 2,500.5 µg/mL | +/- | 14.5381 | µg/mL | Gravimetric |
| | CAS # 75-35-4.SEC (Lot 2767000) | | +/- | 150.8662 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 151.2244 | µg/mL | Stressed |
| 4 | tert-Butanol (TBA) | 25,003.1 µg/mL | +/- | 145.3626 | µg/mL | Gravimetric |
| | CAS # 75-65-0.SEC (Lot XYXDO) | | +/- | 1,508.5475 | µg/mL | Unstressed |
| | Purity 98% | | +/- | 1,512.1291 | µg/mL | Stressed |
| 5 | Methyl acetate | 5,000.4 µg/mL | +/- | 29.0726 | µg/mL | Gravimetric |
| | CAS # 79-20-9.SEC (Lot YDGVD) | | +/- | 301.6948 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 302.4111 | µg/mL | Stressed |
| 6 | Iodomethane (methyl iodide) | 2,500.4 µg/mL | +/- | 14.5374 | µg/mL | Gravimetric |
| | CAS # 74-88-4.SEC (Lot Y25A027) | | +/- | 150.8587 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 151.2169 | µg/mL | Stressed |
| 7 | Allyl chloride (3-chloropropene) | 2,500.1 µg/mL | +/- | 14.5358 | µg/mL | Gravimetric |
| | CAS # 107-05-1.SEC (Lot VEBOC) | | +/- | 150.8423 | µg/mL | Unstressed |
| | Purity 98% | | +/- | 151.2004 | µg/mL | Stressed |

| | | | | | | | |
|----|--------------------------------------|----------|-------|-----|------------|-------|-------------|
| 8 | Methylene chloride (dichloromethane) | 2,500.8 | µg/mL | +/- | 14.5396 | µg/mL | Gravimetric |
| | CAS # 75-09-2.SEC (Lot FGM02) | | | +/- | 150.8813 | µg/mL | Unstressed |
| | Purity 99% | | | +/- | 151.2395 | µg/mL | Stressed |
| 9 | Carbon disulfide | 2,500.9 | µg/mL | +/- | 14.5403 | µg/mL | Gravimetric |
| | CAS # 75-15-0.SEC (Lot MKBL1376V) | | | +/- | 150.8889 | µg/mL | Unstressed |
| | Purity 99% | | | +/- | 151.2471 | µg/mL | Stressed |
| 10 | Acrylonitrile | 25,000.9 | µg/mL | +/- | 145.3496 | µg/mL | Gravimetric |
| | CAS # 107-13-1.SEC (Lot UERIL) | | | +/- | 1,508.4128 | µg/mL | Unstressed |
| | Purity 99% | | | +/- | 1,511.9941 | µg/mL | Stressed |
| 11 | Methyl-tert-butyl ether (MTBE) | 2,500.0 | µg/mL | +/- | 14.5352 | µg/mL | Gravimetric |
| | CAS # 1634-04-4.SEC (Lot ZAQTA-MS) | | | +/- | 150.8361 | µg/mL | Unstressed |
| | Purity 99% | | | +/- | 151.1942 | µg/mL | Stressed |
| 12 | cis-1,2-Dichloroethene | 2,500.7 | µg/mL | +/- | 14.5394 | µg/mL | Gravimetric |
| | CAS # 156-59-2.SEC (Lot HGC01-BLKT) | | | +/- | 150.8792 | µg/mL | Unstressed |
| | Purity 98% | | | +/- | 151.2374 | µg/mL | Stressed |
| 13 | n-Hexane (C6) | 2,501.5 | µg/mL | +/- | 14.5439 | µg/mL | Gravimetric |
| | CAS # 110-54-3.SEC (Lot 10188491) | | | +/- | 150.9266 | µg/mL | Unstressed |
| | Purity 99% | | | +/- | 151.2849 | µg/mL | Stressed |
| 14 | 1,1-Dichloroethane | 2,500.3 | µg/mL | +/- | 14.5367 | µg/mL | Gravimetric |
| | CAS # 75-34-3.SEC (Lot 5379000) | | | +/- | 150.8512 | µg/mL | Unstressed |
| | Purity 99% | | | +/- | 151.2093 | µg/mL | Stressed |
| 15 | 2,2-Dichloropropane | 2,500.1 | µg/mL | +/- | 14.5358 | µg/mL | Gravimetric |
| | CAS # 594-20-7.SEC (Lot I7E8E) | | | +/- | 150.8423 | µg/mL | Unstressed |
| | Purity 98% | | | +/- | 151.2004 | µg/mL | Stressed |
| 16 | trans-1,2-Dichloroethene | 2,500.2 | µg/mL | +/- | 14.5362 | µg/mL | Gravimetric |
| | CAS # 156-60-5.SEC (Lot TS5UB) | | | +/- | 150.8466 | µg/mL | Unstressed |
| | Purity 97% | | | +/- | 151.2048 | µg/mL | Stressed |
| 17 | Isobutanol (2-Methyl-1-propanol) | 62,506.9 | µg/mL | +/- | 363.4014 | µg/mL | Gravimetric |
| | CAS # 78-83-1.SEC (Lot 83NHH) | | | +/- | 3,771.3149 | µg/mL | Unstressed |
| | Purity 99% | | | +/- | 3,780.2687 | µg/mL | Stressed |
| 18 | Chloroform | 2,500.1 | µg/mL | +/- | 14.5359 | µg/mL | Gravimetric |
| | CAS # 67-66-3.SEC (Lot 1297547) | | | +/- | 150.8436 | µg/mL | Unstressed |
| | Purity 99% | | | +/- | 151.2017 | µg/mL | Stressed |
| 19 | Bromochloromethane | 2,501.1 | µg/mL | +/- | 14.5418 | µg/mL | Gravimetric |
| | CAS # 74-97-5.SEC (Lot 5670200) | | | +/- | 150.9040 | µg/mL | Unstressed |
| | Purity 99% | | | +/- | 151.2622 | µg/mL | Stressed |
| 20 | Tetrahydrofuran | 5,002.3 | µg/mL | +/- | 29.0835 | µg/mL | Gravimetric |
| | CAS # 109-99-9.SEC (Lot K3V7J-SJ) | | | +/- | 301.8079 | µg/mL | Unstressed |
| | Purity 99% | | | +/- | 302.5245 | µg/mL | Stressed |
| 21 | 1,1,1-Trichloroethane | 2,500.7 | µg/mL | +/- | 14.5394 | µg/mL | Gravimetric |
| | CAS # 71-55-6.SEC (Lot CS160712) | | | +/- | 150.8792 | µg/mL | Unstressed |
| | Purity 98% | | | +/- | 151.2374 | µg/mL | Stressed |
| 22 | Cyclohexane | 2,501.0 | µg/mL | +/- | 14.5410 | µg/mL | Gravimetric |
| | CAS # 110-82-7.SEC (Lot YADRA) | | | +/- | 150.8964 | µg/mL | Unstressed |
| | Purity 99% | | | +/- | 151.2547 | µg/mL | Stressed |
| 23 | 1,1-Dichloropropene | 2,501.3 | µg/mL | +/- | 14.5427 | µg/mL | Gravimetric |
| | CAS # 563-58-6.SEC (Lot 5221100) | | | +/- | 150.9133 | µg/mL | Unstressed |
| | Purity 96% | | | +/- | 151.2716 | µg/mL | 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 | Trichloroethene 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 | Tetrachloroethene 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 150.8613 151.2194 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 41 | 1,2-Dibromoethane (EDB) CAS # 106-93-4.SEC Purity 99% | (Lot 3505900) | 2,500.5 | µg/mL | +/- +/- +/- | 14.5381 150.8662 151.2244 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 42 | Chlorobenzene CAS # 108-90-7.SEC Purity 99% | (Lot 1161936) | 2,501.0 | µg/mL | +/- +/- +/- | 14.5410 150.8964 151.2547 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 43 | m-Xylene CAS # 108-38-3.SEC Purity 99% | (Lot OUKMG-GB) | 1,250.9 | µg/mL | +/- +/- +/- | 7.2727 75.4708 75.6500 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 44 | p-Xylene CAS # 106-42-3.SEC Purity 99% | (Lot GM01) | 1,250.5 | µg/mL | +/- +/- +/- | 7.2705 75.4482 75.6273 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 45 | Ethylbenzene CAS # 100-41-4.SEC Purity 99% | (Lot PI4SE) | 2,500.9 | µg/mL | +/- +/- +/- | 14.5403 150.8889 151.2471 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 46 | 1,1,1,2-Tetrachloroethane CAS # 630-20-6.SEC Purity 99% | (Lot GC01) | 2,501.1 | µg/mL | +/- +/- +/- | 14.5418 150.9040 151.2622 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 47 | o-Xylene CAS # 95-47-6.SEC Purity 99% | (Lot FGL01-KTPK) | 2,500.9 | µg/mL | +/- +/- +/- | 14.5403 150.8889 151.2471 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 48 | Styrene CAS # 100-42-5.SEC Purity 99% | (Lot OFIOL-IA) | 2,500.4 | µg/mL | +/- +/- +/- | 14.5374 150.8587 151.2169 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 49 | Isopropylbenzene (cumene) CAS # 98-82-8.SEC Purity 99% | (Lot 2PHXG-IH) | 2,500.5 | µg/mL | +/- +/- +/- | 14.5381 150.8662 151.2244 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 50 | Bromoform CAS # 75-25-2.SEC Purity 99% | (Lot 5139000) | 2,502.3 | µg/mL | +/- +/- +/- | 14.5483 150.9718 151.3303 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 51 | Bromodichloromethane CAS # 75-27-4.SEC Purity 98% | (Lot 13780) | 2,500.1 | µg/mL | +/- +/- +/- | 14.5358 150.8423 151.2004 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 52 | 1,1,1,2-Tetrachloroethane CAS # 79-34-5.SEC Purity 99% | (Lot CFA4D-AQ) | 2,501.3 | µg/mL | +/- +/- +/- | 14.5425 150.9115 151.2698 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 53 | 1,2,3-Trichloropropane CAS # 96-18-4.SEC Purity 98% | (Lot OGI01) | 2,500.1 | µg/mL | +/- +/- +/- | 14.5358 150.8423 151.2004 | µg/mL µg/mL µg/mL | Gravimetric Unstressed 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 150.8940 151.2522 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |
| 55 | n-Propylbenzene CAS # 103-65-1.SEC Purity 99% | (Lot T2HFC-IT) | 2,500.0 | µg/mL | +/- +/- +/- | 14.5352 150.8361 151.1942 | µg/mL µg/mL µg/mL | Gravimetric Unstressed Stressed |

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

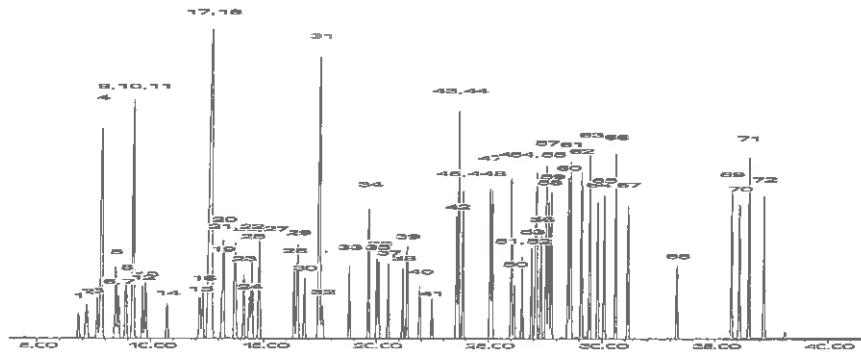
Carrier Gas:
helium-constant pressure 30 psi

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 Mays

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



CERTIFIED REFERENCE MATERIAL

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

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Certificate of Analysis



FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

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

Catalog No. : 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

CERTIFIED VALUES

| Elution Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L., K=2) | | |
|---------------|---|-----------------------------|--------------------------------------|-------|-------------|
| 1 | Dibromofluoromethane CAS # 1868-53-7 Purity 99% (Lot 022012) | 2,509.4 µg/mL | +/- 14.5899 | µg/mL | Gravimetric |
| | | | +/- 140.6996 | µg/mL | Unstressed |
| | | | +/- 143.9918 | µg/mL | Stressed |
| 2 | 1,2-Dichloroethane-d4 CAS # 17060-07-0 Purity 98% (Lot PR-25433) | 2,509.0 µg/mL | +/- 14.5875 | µg/mL | Gravimetric |
| | | | +/- 140.6769 | µg/mL | Unstressed |
| | | | +/- 143.9686 | µg/mL | Stressed |
| 3 | Toluene-d8 CAS # 2037-26-5 Purity 99% (Lot PR-26282) | 2,507.0 µg/mL | +/- 14.5759 | µg/mL | Gravimetric |
| | | | +/- 140.5650 | µg/mL | Unstressed |
| | | | +/- 143.8540 | µg/mL | Stressed |
| 4 | 1-Bromo-4-fluorobenzene (BFB) CAS # 460-00-4 Purity 99% (Lot 20401KOV) | 2,503.6 µg/mL | +/- 14.5561 | µg/mL | Gravimetric |
| | | | +/- 140.3744 | µg/mL | Unstressed |
| | | | +/- 143.6590 | µg/mL | Stressed |

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

Reagent

VOA8260SURRES_00122



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. : 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

CERTIFIED VALUES

| Elution Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L., K=2) | | | |
|---------------|-------------------------------|---------------------------------|--------------------------------------|----------|-------|-------------|
| 1 | Dibromofluoromethane | 2,509.4 µg/mL (Lot 022012) | +/- | 14.5899 | µg/mL | Gravimetric |
| | CAS # 1868-53-7 | | +/- | 140.6996 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 143.9918 | µg/mL | Stressed |
| 2 | 1,2-Dichloroethane-d4 | 2,509.0 µg/mL (Lot PR-25433) | +/- | 14.5875 | µg/mL | Gravimetric |
| | CAS # 17060-07-0 | | +/- | 140.6769 | µg/mL | Unstressed |
| | Purity 98% | | +/- | 143.9686 | µg/mL | Stressed |
| 3 | Toluene-d8 | 2,507.0 µg/mL (Lot PR-26282) | +/- | 14.5759 | µg/mL | Gravimetric |
| | CAS # 2037-26-5 | | +/- | 140.5650 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 143.8540 | µg/mL | Stressed |
| 4 | 1-Bromo-4-fluorobenzene (BFB) | 2,503.6 µg/mL (Lot 20401KOV) | +/- | 14.5561 | µg/mL | Gravimetric |
| | CAS # 460-00-4 | | +/- | 140.3744 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 143.6590 | µg/mL | Stressed |

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

Reagent

VOA8260VARES_00083



CERTIFIED REFERENCE MATERIAL

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Certificate of Analysis



FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

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

Catalog No. : 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.

CERTIFIED VALUES

| Elution Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L.; K=2) | | |
|---------------|--|-----------------------------|--------------------------------------|-------|-------------|
| 1 | Vinyl acetate CAS # 108-05-4 Purity 99% (Lot STBD7333V) | 5,027.0 µg/mL | +/- 29.5013 | µg/mL | Gravimetric |
| | | | +/- 303.3277 | µg/mL | Unstressed |
| | | | +/- 304.0477 | µg/mL | Stressed |

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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FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

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

Catalog No. : 568720 **Lot No.:** 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.

CERTIFIED VALUES

| Elution Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L.; K=2) |
|---------------|---|-----------------------------|--|
| 1 | Acrolein CAS # 107-02-8 Purity 99% (Lot 170123JLM) | 19,779.0 µg/mL | +/- 115.8104 µg/mL Gravimetric +/- 634.1769 µg/mL Unstressed +/- 737.1613 µg/mL Stressed |

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

Reagent

VOABFBRES_00056



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 2,500µg/mL, P&T Methanol, 1mL/ampul

Container Size : 2 mL **Pkg Amt:** > 1 mL

Expiration Date : November 30, 2021 **Storage:** 0°C or colder

CERTIFIED VALUES

| Elution Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L.; K=2) | | | |
|---------------|--|-----------------------------|--------------------------------------|-------|-------------|--|
| 1 | 1-Bromo-4-fluorobenzene (BFB) CAS # 460-00-4 (Lot 20401KOV) Purity 99% | 2,524.0 µg/mL | +/- 14.8122 | µg/mL | Gravimetric | |
| | | | +/- 141.5325 | µg/mL | Unstressed | |
| | | | +/- 144.8435 | µg/mL | 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 2,500µg/mL, P&T Methanol, 1mL/ampul

Container Size : 2 mL **Pkg Amt:** > 1 mL

Expiration Date : November 30, 2021 **Storage:** 0°C or colder

CERTIFIED VALUES

| Elution Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L.; K=2) | | | |
|---------------|--|-----------------------------|--------------------------------------|-------|-------------|--|
| 1 | 1-Bromo-4-fluorobenzene (BFB) CAS # 460-00-4 (Lot 20401KOV) Purity 99% | 2,524.0 µg/mL | +/- 14.8122 | µg/mL | Gravimetric | |
| | | | +/- 141.5325 | µg/mL | Unstressed | |
| | | | +/- 144.8435 | µg/mL | Stressed | |

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

Reagent

VOACEVERES_00127



CERTIFIED REFERENCE MATERIAL

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Certificate of Analysis



FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

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

Catalog No. : 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

2406027
28
29
30

CERTIFIED VALUES

| 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% (Lot MKBS6526V) | 2,503.5 µg/mL | +/- | 14.5556 | µg/mL | Gravimetric |
| | | | +/- | 53.6004 | µg/mL | Unstressed |
| | | | +/- | 55.1587 | µg/mL | Stressed |

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

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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.

2396751

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

Description : Custom EE Standard
Custom EE Standard 5,000µg/mL, P&T Methanol, 1mL/ampul

Container Size : 2 mL Pkg Amt: > 1 mL

Expiration Date : January 31, 2018 Storage: 0°C or colder

CERTIFIED VALUES

| Elution Order | Compound | Grav. Conc. (weight/volume) | Expanded Uncertainty (95% C.L.; K=2) | | | |
|---------------|------------------------------|----------------------------------|--------------------------------------|----------|-------|-------------|
| 1 | 3-Chlorobenzotrifluoride | 5,025.0 µg/mL (Lot 21324DO) | +/- | 29.4895 | µg/mL | Gravimetric |
| | CAS # 98-15-7 | | +/- | 281.7753 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 288.3671 | µg/mL | Stressed |
| 2 | 4-Chlorobenzotrifluoride | 5,031.0 µg/mL (Lot 08507BO) | +/- | 29.5247 | µg/mL | Gravimetric |
| | CAS # 98-56-6 | | +/- | 282.1117 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 288.7115 | µg/mL | Stressed |
| 3 | 2-Chlorobenzotrifluoride | 5,011.0 µg/mL (Lot I0316DQ) | +/- | 29.4074 | µg/mL | Gravimetric |
| | CAS # 88-16-4 | | +/- | 280.9902 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 287.5637 | µg/mL | Stressed |
| 4 | 3-Chlorotoluene | 5,046.0 µg/mL (Lot 13528LX) | +/- | 29.6128 | µg/mL | Gravimetric |
| | CAS # 108-41-8 | | +/- | 282.9528 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 289.5723 | µg/mL | Stressed |
| 5 | 2,4-Dichlorobenzotrifluoride | 5,018.0 µg/mL (Lot MKBL3552V) | +/- | 29.4484 | µg/mL | Gravimetric |
| | CAS # 320-60-5 | | +/- | 281.3828 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 287.9654 | µg/mL | Stressed |
| 6 | 3,4-Dichlorobenzotrifluoride | 5,031.0 µg/mL (Lot 11105EJV) | +/- | 29.5247 | µg/mL | Gravimetric |
| | CAS # 328-84-7 | | +/- | 282.1117 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 288.7115 | µg/mL | Stressed |
| 7 | 2,5-Dichlorobenzotrifluoride | 5,047.0 µg/mL (Lot 04415DSV) | +/- | 29.6186 | µg/mL | Gravimetric |
| | CAS # 320-50-3 | | +/- | 283.0089 | µg/mL | Unstressed |
| | Purity 99% | | +/- | 289.6296 | µg/mL | Stressed |

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.: 180-70873-1

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-22-0/1-0 | 180-70873-1 | 101 | 113 | 97 | 94 |
| HD-SPBA-CW-22-0/1-0 DL | 180-70873-1 DL | 109 | 117 | 99 | 94 |
| HD-QC4-0/1-2 | 180-70873-2 | 101 | 113 | 101 | 92 |
| | MB 180-224792/6 | 94 | 107 | 98 | 93 |
| | MB 180-224919/7 | 96 | 111 | 101 | 95 |
| | LCS 180-224792/4 | 98 | 106 | 113 | 107 |
| | LCS 180-224919/4 | 90 | 102 | 110 | 102 |

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

QC LIMITS
73-120
65-121
73-120
80-120

Column to be used to flag recovery values

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-70873-1

SDG No.: _____

Matrix: Water Level: Low

Lab File ID: 51003D04.D

Lab ID: LCS 180-224792/4

Client ID: _____

| COMPOUND | SPIKE ADDED (ug/L) | LCS CONCENTRATION (ug/L) | LCS % REC | QC LIMITS REC | # |
|-----------------------------|--------------------------|--------------------------------|-----------------|---------------------|---|
| Chloromethane | 10.0 | 11.6 | 116 | 49-135 | |
| Vinyl chloride | 10.0 | 11.8 | 118 | 52-136 | |
| Bromomethane | 10.0 | 10.1 | 101 | 37-150 | |
| Chloroethane | 10.0 | 13.5 | 135 | 44-139 | |
| 1,1-Dichloroethene | 10.0 | 10.3 | 103 | 64-131 | |
| Acetone | 20.0 | 22.4 | 112 | 24-150 | |
| Carbon disulfide | 10.0 | 9.37 | 94 | 20-150 | |
| Methylene Chloride | 10.0 | 9.77 | 98 | 66-123 | |
| trans-1,2-Dichloroethene | 10.0 | 10.0 | 100 | 70-123 | |
| Methyl tert-butyl ether | 10.0 | 9.62 | 96 | 66-130 | |
| 1,1-Dichloroethane | 10.0 | 9.85 | 98 | 66-122 | |
| cis-1,2-Dichloroethene | 10.0 | 9.81 | 98 | 73-120 | |
| Bromochloromethane | 10.0 | 9.53 | 95 | 73-122 | |
| 2-Butanone (MEK) | 20.0 | 21.0 | 105 | 37-150 | |
| Chloroform | 10.0 | 9.83 | 98 | 72-123 | |
| 1,1,1-Trichloroethane | 10.0 | 10.4 | 104 | 66-129 | |
| Carbon tetrachloride | 10.0 | 9.98 | 100 | 58-145 | |
| Benzene | 10.0 | 9.68 | 97 | 75-123 | |
| 1,2-Dichloroethane | 10.0 | 10.6 | 106 | 63-130 | |
| Trichloroethene | 10.0 | 9.28 | 93 | 74-121 | |
| 1,2-Dichloropropane | 10.0 | 9.20 | 92 | 67-119 | |
| Bromodichloromethane | 10.0 | 9.11 | 91 | 62-127 | |
| cis-1,3-Dichloropropene | 10.0 | 8.97 | 90 | 61-127 | |
| 4-Methyl-2-pentanone (MIBK) | 20.0 | 21.2 | 106 | 41-135 | |
| Toluene | 10.0 | 10.9 | 109 | 76-129 | |
| trans-1,3-Dichloropropene | 10.0 | 9.70 | 97 | 61-136 | |
| 1,1,2-Trichloroethane | 10.0 | 10.8 | 108 | 74-126 | |
| Tetrachloroethene | 10.0 | 10.5 | 105 | 76-128 | |
| 2-Hexanone | 20.0 | 21.2 | 106 | 37-150 | |
| Dibromochloromethane | 10.0 | 9.60 | 96 | 63-131 | |
| 1,2-Dibromoethane (EDB) | 10.0 | 9.86 | 99 | 76-128 | |
| Chlorobenzene | 10.0 | 10.3 | 103 | 79-124 | |
| 1,1,1,2-Tetrachloroethane | 10.0 | 10.6 | 106 | 70-130 | |
| Ethylbenzene | 10.0 | 9.83 | 98 | 77-124 | |
| Xylenes, Total | 20.0 | 20.2 | 101 | 76-124 | |
| Styrene | 10.0 | 10.0 | 100 | 80-125 | |
| Bromoform | 10.0 | 8.98 | 90 | 54-136 | |
| 1,1,2,2-Tetrachloroethane | 10.0 | 10.4 | 104 | 72-128 | |
| Acrylonitrile | 100 | 108 | 108 | 60-130 | |
| 1,4-Dioxane | 200 | 271 | 135 | 26-150 | |

Column to be used to flag recovery and RPD values

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Pittsburgh

Job No.: 180-70873-1

SDG No.: _____

Matrix: Water Level: Low

Lab File ID: 51004D04.D

Lab ID: LCS 180-224919/4

Client ID: _____

| COMPOUND | SPIKE ADDED (ug/L) | LCS CONCENTRATION (ug/L) | LCS % REC | QC LIMITS REC | # |
|-----------------------------|--------------------------|--------------------------------|-----------------|---------------------|---|
| Chloromethane | 10.0 | 12.0 | 120 | 49-135 | |
| Vinyl chloride | 10.0 | 12.6 | 126 | 52-136 | |
| Bromomethane | 10.0 | 12.4 | 124 | 37-150 | |
| Chloroethane | 10.0 | 12.5 | 125 | 44-139 | |
| 1,1-Dichloroethene | 10.0 | 9.95 | 100 | 64-131 | |
| Acetone | 20.0 | 21.7 | 109 | 24-150 | |
| Carbon disulfide | 10.0 | 9.04 | 90 | 20-150 | |
| Methylene Chloride | 10.0 | 8.90 | 89 | 66-123 | |
| trans-1,2-Dichloroethene | 10.0 | 9.52 | 95 | 70-123 | |
| Methyl tert-butyl ether | 10.0 | 9.23 | 92 | 66-130 | |
| 1,1-Dichloroethane | 10.0 | 9.51 | 95 | 66-122 | |
| cis-1,2-Dichloroethene | 10.0 | 8.65 | 86 | 73-120 | |
| Bromochloromethane | 10.0 | 8.75 | 88 | 73-122 | |
| 2-Butanone (MEK) | 20.0 | 18.3 | 91 | 37-150 | |
| Chloroform | 10.0 | 9.05 | 91 | 72-123 | |
| 1,1,1-Trichloroethane | 10.0 | 9.46 | 95 | 66-129 | |
| Carbon tetrachloride | 10.0 | 9.46 | 95 | 58-145 | |
| Benzene | 10.0 | 8.85 | 88 | 75-123 | |
| 1,2-Dichloroethane | 10.0 | 9.68 | 97 | 63-130 | |
| Trichloroethene | 10.0 | 8.20 | 82 | 74-121 | |
| 1,2-Dichloropropane | 10.0 | 8.44 | 84 | 67-119 | |
| Bromodichloromethane | 10.0 | 8.21 | 82 | 62-127 | |
| cis-1,3-Dichloropropene | 10.0 | 8.18 | 82 | 61-127 | |
| 4-Methyl-2-pentanone (MIBK) | 20.0 | 20.4 | 102 | 41-135 | |
| Toluene | 10.0 | 10.4 | 104 | 76-129 | |
| trans-1,3-Dichloropropene | 10.0 | 9.78 | 98 | 61-136 | |
| 1,1,2-Trichloroethane | 10.0 | 10.6 | 106 | 74-126 | |
| Tetrachloroethene | 10.0 | 9.77 | 98 | 76-128 | |
| 2-Hexanone | 20.0 | 19.4 | 97 | 37-150 | |
| Dibromochloromethane | 10.0 | 9.57 | 96 | 63-131 | |
| 1,2-Dibromoethane (EDB) | 10.0 | 9.26 | 93 | 76-128 | |
| Chlorobenzene | 10.0 | 9.72 | 97 | 79-124 | |
| 1,1,1,2-Tetrachloroethane | 10.0 | 9.84 | 98 | 70-130 | |
| Ethylbenzene | 10.0 | 9.31 | 93 | 77-124 | |
| Xylenes, Total | 20.0 | 18.9 | 95 | 76-124 | |
| Styrene | 10.0 | 9.11 | 91 | 80-125 | |
| Bromoform | 10.0 | 8.75 | 88 | 54-136 | |
| 1,1,2,2-Tetrachloroethane | 10.0 | 9.67 | 97 | 72-128 | |
| Acrylonitrile | 100 | 102 | 102 | 60-130 | |
| 1,4-Dioxane | 200 | 189 J | 94 | 26-150 | |

Column to be used to flag recovery and RPD values

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
SDG No.: _____
Lab File ID: 51003D06.D Lab Sample ID: MB 180-224792/6
Matrix: Water Heated Purge: (Y/N) N
Instrument ID: CHHP5 Date Analyzed: 10/04/2017 02:21
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 180-224792/4 | 51003D04.D | 10/04/2017 01:24 |
| HD-QC4-0/1-2 | 180-70873-2 | 51003D12.D | 10/04/2017 04:59 |
| HD-SPBA-CW-22-0/1-0 | 180-70873-1 | 51003D13.D | 10/04/2017 05:23 |

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Lab File ID: 51004D07.D Lab Sample ID: MB 180-224919/7
 Matrix: Water Heated Purge: (Y/N) N
 Instrument ID: CHHP5 Date Analyzed: 10/05/2017 02:41
 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 180-224919/4 | 51004D04.D | 10/05/2017 01:09 |
| HD-SPBA-CW-22-0/1-0 DL | 180-70873-1 DL | 51004D16.D | 10/05/2017 06:31 |

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

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Lab File ID: 50727D01.D BFB Injection Date: 07/27/2017
 Instrument ID: CHHP5 BFB Injection Time: 00:22
 Analysis Batch No.: 218218

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 50 | 15.0 - 40.0 % of mass 95 | 16.0 |
| 75 | 30.0 - 60.0 % of mass 95 | 47.0 |
| 95 | Base Peak, 100% relative abundance | 100.0 |
| 96 | 5.0 - 9.0 % of mass 95 | 7.9 |
| 173 | Less than 2.0 % of mass 174 | 0.4 (0.5) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 75.4 |
| 175 | 5.0 - 9.0 % of mass 174 | 5.4 (7.2) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 74.0 (98.2) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.8 (6.5) 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 |
|------------------|-------------------|-------------|---------------|---------------|
| | IC 180-218218/2 | 50727D02.D | 07/27/2017 | 00:51 |
| | IC 180-218218/3 | 50727D03.D | 07/27/2017 | 01:15 |
| | ICIS 180-218218/4 | 50727D04.D | 07/27/2017 | 01:39 |
| | IC 180-218218/5 | 50727D05.D | 07/27/2017 | 02:02 |
| | IC 180-218218/6 | 50727D06.D | 07/27/2017 | 02:26 |
| | IC 180-218218/8 | 50727D08.D | 07/27/2017 | 03:13 |
| | IC 180-218218/10 | 50727D10.D | 07/27/2017 | 04:00 |
| | IC 180-218218/11 | 50727D11.D | 07/27/2017 | 04:24 |

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

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Lab File ID: 51003D01.D BFB Injection Date: 10/03/2017
 Instrument ID: CHHP5 BFB Injection Time: 23:49
 Analysis Batch No.: 224792

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 50 | 15.0 - 40.0 % of mass 95 | 17.2 |
| 75 | 30.0 - 60.0 % of mass 95 | 45.1 |
| 95 | Base Peak, 100% relative abundance | 100.0 |
| 96 | 5.0 - 9.0 % of mass 95 | 5.5 |
| 173 | Less than 2.0 % of mass 174 | 0.0 (0.0) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 63.6 |
| 175 | 5.0 - 9.0 % of mass 174 | 5.0 (7.9) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 63.8 (100.3) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 3.3 (5.2) 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 180-224792/2 | 51003D02.D | 10/04/2017 | 00:22 |
| | LCS 180-224792/4 | 51003D04.D | 10/04/2017 | 01:24 |
| | MB 180-224792/6 | 51003D06.D | 10/04/2017 | 02:21 |
| HD-QC4-0/1-2 | 180-70873-2 | 51003D12.D | 10/04/2017 | 04:59 |
| HD-SPBA-CW-22-0/1-0 | 180-70873-1 | 51003D13.D | 10/04/2017 | 05:23 |

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

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Lab File ID: 51004D01.D BFB Injection Date: 10/04/2017
 Instrument ID: CHHP5 BFB Injection Time: 22:24
 Analysis Batch No.: 224919

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|-----|------------------------------------|----------------------|----------|
| 50 | 15.0 - 40.0 % of mass 95 | 17.7 | |
| 75 | 30.0 - 60.0 % of mass 95 | 44.6 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 7.8 | |
| 173 | Less than 2.0 % of mass 174 | 0.4 | (0.6) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 68.2 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.4 | (7.9) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 67.8 | (99.4) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.1 | (6.1) 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 |
|------------------------|--------------------|-------------|---------------|---------------|
| | CCVIS 180-224919/2 | 51004D02.D | 10/04/2017 | 23:29 |
| | LCS 180-224919/4 | 51004D04.D | 10/05/2017 | 01:09 |
| | MB 180-224919/7 | 51004D07.D | 10/05/2017 | 02:41 |
| HD-SPBA-CW-22-0/1-0 DL | 180-70873-1 DL | 51004D16.D | 10/05/2017 | 06:31 |

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

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Sample No.: CCVIS 180-224792/2 Date Analyzed: 10/04/2017 00:22
 Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm)
 Lab File ID (Standard): 51003D02.D Heated Purge: (Y/N) N
 Calibration ID: 35038

| | TBA _d 9 | | FB | | CBN _{Zd} 5 | | |
|------------------|---------------------|--------|--------|--------|---------------------|-------|-------|
| | AREA # | RT # | AREA # | RT # | AREA # | RT # | |
| 12/24 HOUR STD | 166571 | 4.38 | 398262 | 7.33 | 81838 | 10.43 | |
| UPPER LIMIT | 333142 | 4.88 | 796524 | 7.83 | 163676 | 10.93 | |
| LOWER LIMIT | 83286 | 3.88 | 199131 | 6.83 | 40919 | 9.93 | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| LCS 180-224792/4 | 183806 | 4.36 | 350927 | 7.34 | 75706 | 10.43 | |
| MB 180-224792/6 | 193789 | 4.35 | 384752 | 7.34 | 86410 | 10.43 | |
| 180-70873-2 | HD-QC4-0/1-2 | 177857 | 4.36 | 356184 | 7.34 | 75850 | 10.43 |
| 180-70873-1 | HD-SPBA-CW-22-0/1-0 | 187446 | 4.35 | 362965 | 7.34 | 85357 | 10.43 |

TBA_d9 = TBA-d9 (IS)

FB = Fluorobenzene (IS)

CBN_{Zd}5 = 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.: 180-70873-1
 SDG No.: _____
 Sample No.: CCVIS 180-224792/2 Date Analyzed: 10/04/2017 00:22
 Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm)
 Lab File ID (Standard): 51003D02.D Heated Purge: (Y/N) N
 Calibration ID: 35038

| | | DCBd4 | | | | | |
|------------------|---------------------|--------|-------|--------|------|--------|------|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 115198 | 12.77 | | | | |
| UPPER LIMIT | | 230396 | 13.27 | | | | |
| LOWER LIMIT | | 57599 | 12.27 | | | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| LCS 180-224792/4 | | 111504 | 12.77 | | | | |
| MB 180-224792/6 | | 122049 | 12.77 | | | | |
| 180-70873-2 | HD-QC4-0/1-2 | 109308 | 12.77 | | | | |
| 180-70873-1 | HD-SPBA-CW-22-0/1-0 | 114932 | 12.77 | | | | |

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.: 180-70873-1
 SDG No.: _____
 Sample No.: CCVIS 180-224919/2 Date Analyzed: 10/04/2017 23:29
 Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm)
 Lab File ID (Standard): 51004D02.D Heated Purge: (Y/N) N
 Calibration ID: 35038

| | TBA _d 9 | | FB | | CBN _{Zd} 5 | | |
|------------------|---------------------------|--------|--------|--------|---------------------|-------|-------|
| | AREA # | RT # | AREA # | RT # | AREA # | RT # | |
| 12/24 HOUR STD | 150069 | 4.37 | 334551 | 7.34 | 71810 | 10.43 | |
| UPPER LIMIT | 300138 | 4.87 | 669102 | 7.84 | 143620 | 10.93 | |
| LOWER LIMIT | 75035 | 3.87 | 167276 | 6.84 | 35905 | 9.93 | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| LCS 180-224919/4 | | 181022 | 4.37 | 356874 | 7.34 | 73297 | 10.43 |
| MB 180-224919/7 | | 189568 | 4.36 | 369267 | 7.34 | 81221 | 10.43 |
| 180-70873-1 DL | HD-SPBA-CW-22-0/1-0 DL | 149753 | 4.36 | 323921 | 7.34 | 71240 | 10.43 |

TBA_d9 = TBA-d9 (IS)

FB = Fluorobenzene (IS)

CBN_{Zd}5 = 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.: 180-70873-1
 SDG No.: _____
 Sample No.: CCVIS 180-224919/2 Date Analyzed: 10/04/2017 23:29
 Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm)
 Lab File ID (Standard): 51004D02.D Heated Purge: (Y/N) N
 Calibration ID: 35038

| | | DCBd4 | | | | | |
|------------------|---------------------------|--------|-------|--------|------|--------|------|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 99164 | 12.77 | | | | |
| UPPER LIMIT | | 198328 | 13.27 | | | | |
| LOWER LIMIT | | 49582 | 12.27 | | | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| LCS 180-224919/4 | | 101073 | 12.77 | | | | |
| MB 180-224919/7 | | 115186 | 12.77 | | | | |
| 180-70873-1 DL | HD-SPBA-CW-22-0/1-0 DL | 100827 | 12.77 | | | | |

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.: 180-70873-1
 SDG No.: _____
 Client Sample ID: HD-SPBA-CW-22-0/1-0 Lab Sample ID: 180-70873-1
 Matrix: Water Lab File ID: 51003D13.D
 Analysis Method: 8260C Date Collected: 09/29/2017 09:40
 Sample wt/vol: 5 (mL) Date Analyzed: 10/04/2017 05:23
 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.: 224792 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|------------|-----------------------------|--------|------|-----|------|
| 74-87-3 | Chloromethane | 1.0 | U ^c | 1.0 | 0.38 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 1.0 | 0.17 |
| 74-83-9 | Bromomethane | 1.0 | U | 1.0 | 0.59 |
| 75-00-3 | Chloroethane | 1.0 | U | 1.0 | 0.58 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.32 |
| 67-64-1 | Acetone | 5.0 | U | 5.0 | 3.1 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 1.0 | 0.53 |
| 75-09-2 | Methylene Chloride | 1.0 | U | 1.0 | 0.94 |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.20 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.34 |
| 156-59-2 | cis-1,2-Dichloroethene | 1.7 | | 1.0 | 0.30 |
| 74-97-5 | Bromochloromethane | 1.0 | U | 1.0 | 0.36 |
| 78-93-3 | 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 |
| 67-66-3 | Chloroform | 1.0 | U | 1.0 | 0.27 |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.27 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 1.0 | 0.56 |
| 71-43-2 | Benzene | 1.0 | U | 1.0 | 0.18 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.24 |
| 79-01-6 | Trichloroethene | 220 | E | 1.0 | 0.20 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.35 |
| 75-27-4 | Bromodichloromethane | 1.0 | U ^c | 1.0 | 0.57 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U ^c | 1.0 | 0.32 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 5.0 | U | 5.0 | 2.2 |
| 108-88-3 | Toluene | 1.0 | U | 1.0 | 0.16 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.22 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.31 |
| 127-18-4 | Tetrachloroethene | 320 | E | 1.0 | 0.24 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 5.0 | 2.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 1.0 | 0.44 |
| 106-93-4 | 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.51 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 1.0 | 0.15 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.49 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 1.0 | 0.25 |
| 1330-20-7 | Xylenes, Total | 2.0 | U | 2.0 | 0.27 |
| 100-42-5 | Styrene | 1.0 | U | 1.0 | 0.22 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Client Sample ID: HD-SPBA-CW-22-0/1-0 Lab Sample ID: 180-70873-1
 Matrix: Water Lab File ID: 51003D13.D
 Analysis Method: 8260C Date Collected: 09/29/2017 09:40
 Sample wt/vol: 5 (mL) Date Analyzed: 10/04/2017 05:23
 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.: 224792 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|----------|---------------------------|--------|------|-----|------|
| 75-25-2 | Bromoform | 1.0 | U ^c | 1.0 | 0.76 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.37 |
| 107-13-1 | Acrylonitrile | 20 | U | 20 | 3.3 |
| 123-91-1 | 1,4-Dioxane | 200 | U | 200 | 16 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|------------|------------------------------|------|---|--------|
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 113 | | 65-121 |
| 2037-26-5 | Toluene-d8 (Surr) | 97 | | 73-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 94 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 73-120 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D13.D
 Lims ID: 180-70873-A-1
 Client ID: HD-SPBA-CW-22-0/1-0
 Sample Type: Client
 Inject. Date: 04-Oct-2017 05:23:30 ALS Bottle#: 13 Worklist Smp#: 13
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018710-013
 Misc. Info.: 180-70873-A-1
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 04-Oct-2017 21:10:21 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK005

First Level Reviewer: bungardf

Date: 04-Oct-2017 20:55:35

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|----|----------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.353 | 4.347 | 0.006 | 0 | 187446 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.340 | 7.334 | 0.006 | 99 | 362965 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.430 | 10.431 | -0.001 | 86 | 85357 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.772 | 12.773 | -0.001 | 97 | 114932 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.616 | 6.616 | 0.000 | 92 | 88499 | 50.7 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.987 | 6.981 | 0.006 | 0 | 120691 | 56.7 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.982 | 8.977 | 0.005 | 93 | 330827 | 48.7 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.610 | 11.611 | -0.001 | 84 | 115743 | 47.2 | |
| 12 Chloromethane | 50 | | 1.823 | | | | ND | |
| 13 Vinyl chloride | 62 | | 1.963 | | | | ND | |
| 15 Bromomethane | 94 | | 2.291 | | | | ND | |
| 16 Chloroethane | 64 | | 2.461 | | | | ND | |
| 22 1,1-Dichloroethene | 96 | | 3.423 | | | | ND | |
| 24 Acetone | 43 | 3.537 | 3.526 | 0.011 | 72 | 6332 | 6.67 | |
| 26 Carbon disulfide | 76 | | 3.708 | | | | ND | |
| 31 Methylene Chloride | 84 | | 4.232 | | | | ND | |
| 33 Acrylonitrile | 53 | | 4.609 | | | | ND | |
| 34 trans-1,2-Dichloroethene | 96 | | 4.633 | | | | ND | |
| 35 Methyl tert-butyl ether | 73 | 4.651 | 4.657 | -0.006 | 18 | 1691 | 0.3115 | |
| 37 1,1-Dichloroethane | 63 | | 5.272 | | | | ND | |
| 45 cis-1,2-Dichloroethene | 96 | 6.007 | 6.008 | -0.001 | 79 | 19530 | 8.43 | |
| 46 2-Butanone (MEK) | 43 | | 6.026 | | | | ND | |
| 49 Chlorobromomethane | 128 | | 6.294 | | | | ND | |
| 52 Chloroform | 83 | 6.427 | 6.440 | -0.013 | 37 | 4434 | 1.26 | |
| 53 1,1,1-Trichloroethane | 97 | | 6.592 | | | | ND | |
| 56 Carbon tetrachloride | 117 | | 6.762 | | | | ND | |
| 58 Benzene | 78 | | 6.994 | | | | ND | |
| 59 1,2-Dichloroethane | 62 | | 7.067 | | | | ND | |
| 64 Trichloroethene | 130 | 7.723 | 7.724 | -0.001 | 97 | 2397052 | 1079.3 | E |
| 67 1,2-Dichloropropane | 63 | | 7.997 | | | | ND | |
| 70 1,4-Dioxane | 88 | | 8.082 | | | | ND | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|---------------|----|----------|--------------|-------|
| 71 Dichlorobromomethane | 83 | | 8.277 | | | | ND | |
| 74 cis-1,3-Dichloropropene | 75 | | 8.721 | | | | ND | |
| 75 4-Methyl-2-pentanone (MIBK) | 43 | | 8.873 | | | | ND | |
| 76 Toluene | 91 | | 9.044 | | | | ND | |
| 77 trans-1,3-Dichloropropene | 75 | | 9.293 | | | | ND | |
| 79 1,1,2-Trichloroethane | 97 | | 9.488 | | | | ND | |
| 80 Tetrachloroethene | 164 | 9.560 | 9.561 | -0.001 | 91 | 2568655 | 1582.5 | E |
| 82 2-Hexanone | 43 | | 9.707 | | | | ND | |
| 84 Chlorodibromomethane | 129 | | 9.853 | | | | ND | |
| 85 Ethylene Dibromide | 107 | | 9.974 | | | | ND | |
| 87 Chlorobenzene | 112 | | 10.455 | | | | ND | |
| 89 1,1,1,2-Tetrachloroethane | 131 | | 10.552 | | | | ND | |
| 90 Ethylbenzene | 106 | | 10.558 | | | | ND | |
| 91 m-Xylene & p-Xylene | 106 | | 10.692 | | | | ND | |
| 92 o-Xylene | 106 | | 11.069 | | | | ND | |
| 93 Styrene | 104 | | 11.094 | | | | ND | |
| 94 Bromoform | 173 | | 11.270 | | | | ND | |
| 99 1,1,2,2-Tetrachloroethane | 83 | | 11.751 | | | | ND | |
| S 133 Xylenes, Total | 106 | | 1.000 | | | | ND | |

QC Flag Legend

Processing Flags

E - Exceeded Maximum Amount

Reagents:

VOA8260INT_00074

Amount Added: 2.00

Units: uL

Run Reagent

VOA8260SURR_00073

Amount Added: 2.00

Units: uL

Run Reagent

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D13.D

Injection Date: 04-Oct-2017 05:23:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: 180-70873-A-1

Lab Sample ID: 180-70873-1

Worklist Smp#: 13

Client ID: HD-SPBA-CW-22-0/1-0

Purge Vol: 5.000 mL

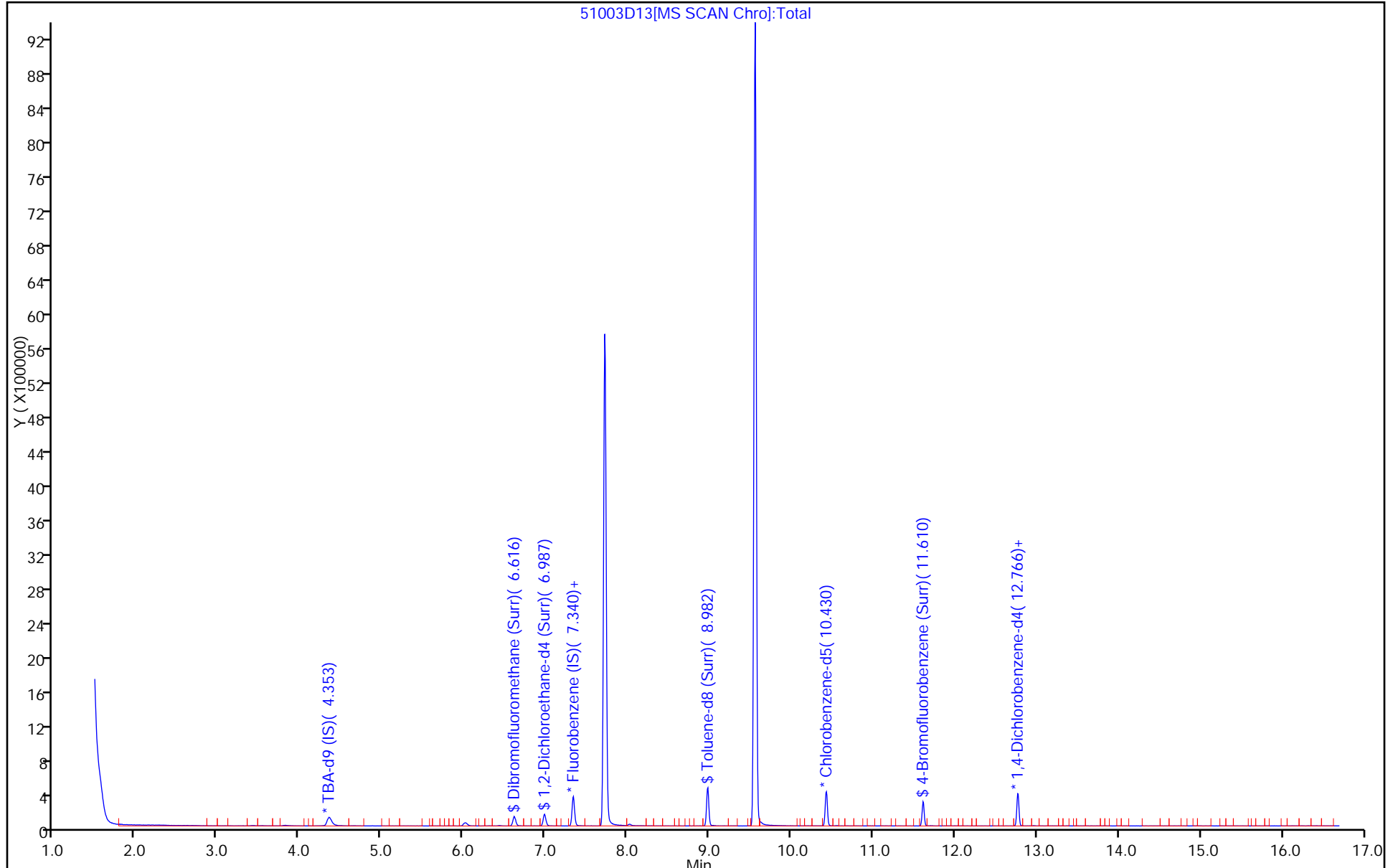
Dil. Factor: 1.0000

ALS Bottle#: 13

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Recovery Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D13.D
 Lims ID: 180-70873-A-1
 Client ID: HD-SPBA-CW-22-0/1-0
 Sample Type: Client
 Inject. Date: 04-Oct-2017 05:23:30 ALS Bottle#: 13 Worklist Smp#: 13
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018710-013
 Misc. Info.: 180-70873-A-1
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 04-Oct-2017 21:10:21 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK005

First Level Reviewer: bungardf

Date: 04-Oct-2017 20:55:35

| Compound | Amount Added | Amount Recovered | % Rec. |
|-----------------------------------|--------------|------------------|--------|
| \$ 5 Dibromofluoromethane (Surr) | 50.0 | 50.7 | 101.35 |
| \$ 6 1,2-Dichloroethane-d4 (Surr) | 50.0 | 56.7 | 113.32 |
| \$ 7 Toluene-d8 (Surr) | 50.0 | 48.7 | 97.40 |
| \$ 8 4-Bromofluorobenzene (Surr) | 50.0 | 47.2 | 94.35 |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D13.D

Injection Date: 04-Oct-2017 05:23:30

Instrument ID: CHHP5

Lims ID: 180-70873-A-1

Lab Sample ID: 180-70873-1

Client ID: HD-SPBA-CW-22-0/1-0

Operator ID: 034635

ALS Bottle#: 13

Worklist Smp#: 13

Purge Vol: 5.000 mL

Dil. Factor: 1.0000

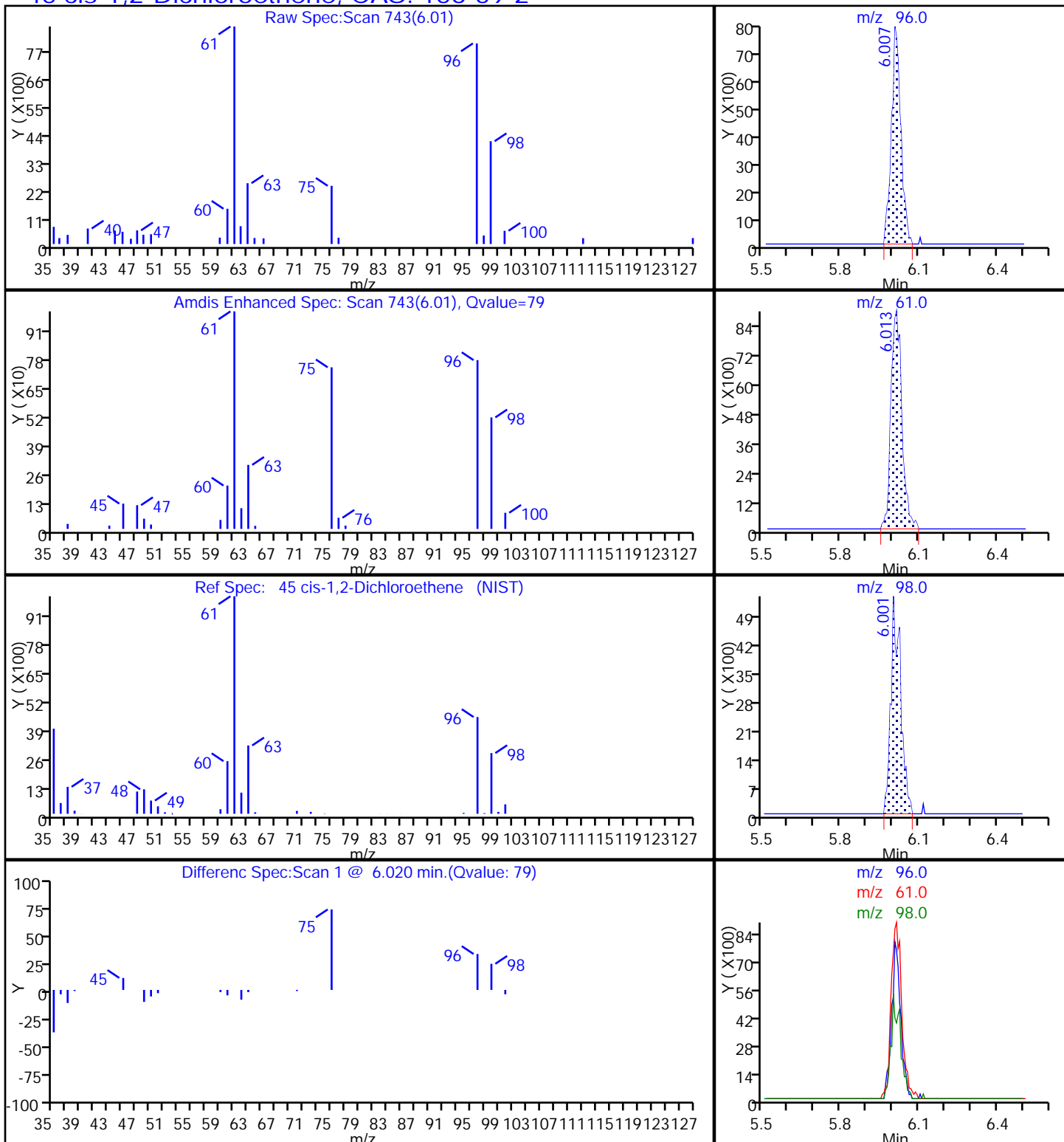
Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)

Detector: MS SCAN

45 cis-1,2-Dichloroethene, CAS: 156-59-2



TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D13.D

Injection Date: 04-Oct-2017 05:23:30

Instrument ID: CHHP5

Lims ID: 180-70873-A-1

Lab Sample ID: 180-70873-1

Client ID: HD-SPBA-CW-22-0/1-0

Operator ID: 034635

ALS Bottle#: 13

Worklist Smp#: 13

Purge Vol: 5.000 mL

Dil. Factor: 1.0000

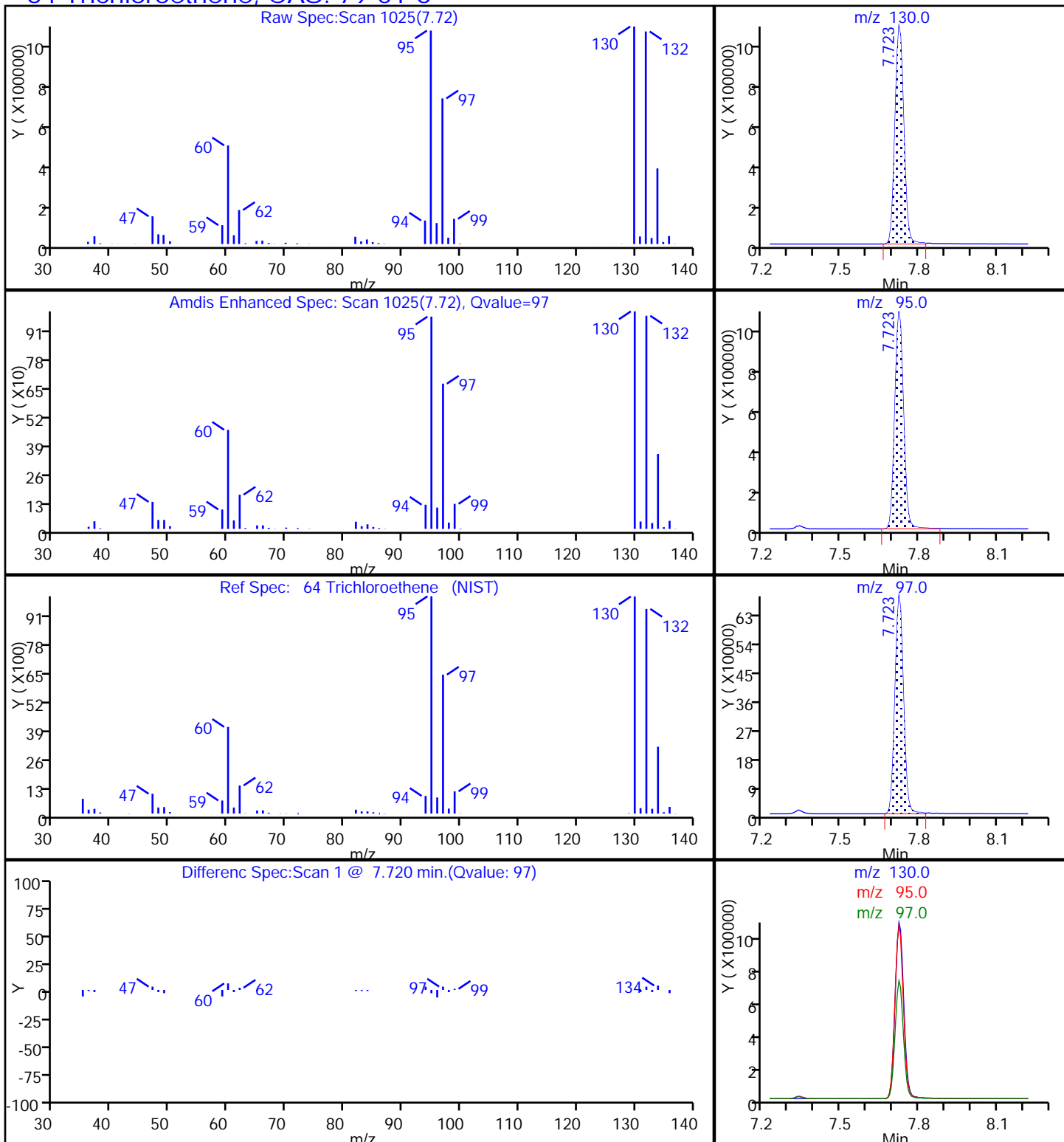
Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)

Detector: MS SCAN

64 Trichloroethene, CAS: 79-01-6



TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D13.D

Injection Date: 04-Oct-2017 05:23:30

Instrument ID: CHHP5

Lims ID: 180-70873-A-1

Lab Sample ID: 180-70873-1

Client ID: HD-SPBA-CW-22-0/1-0

Operator ID: 034635

ALS Bottle#: 13

Worklist Smp#: 13

Purge Vol: 5.000 mL

Dil. Factor: 1.0000

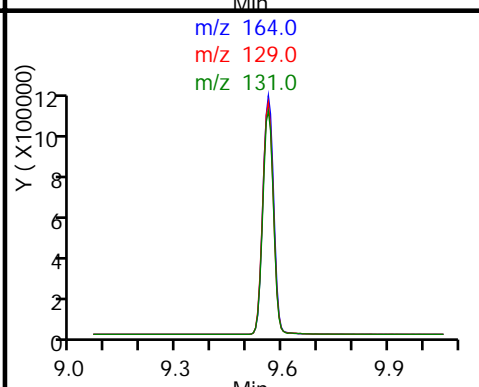
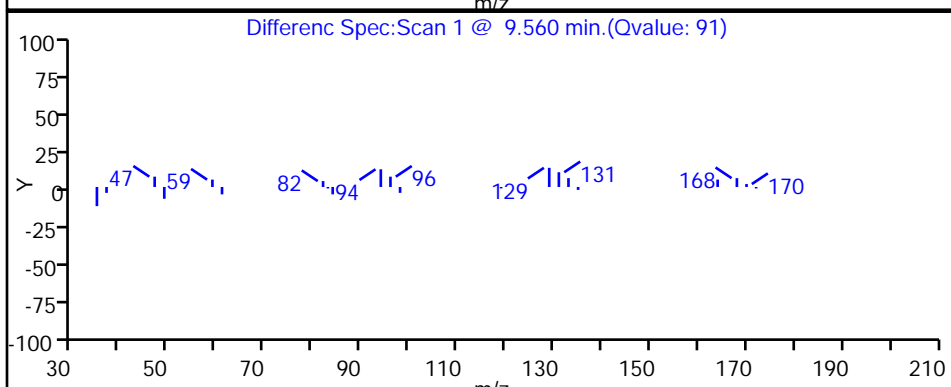
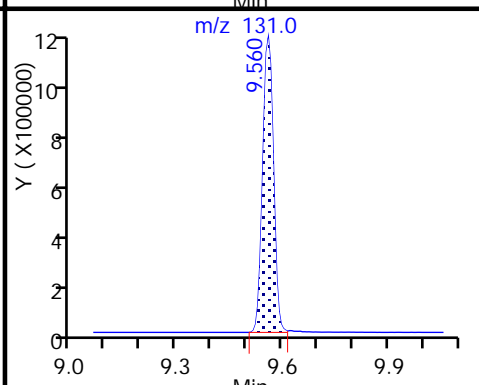
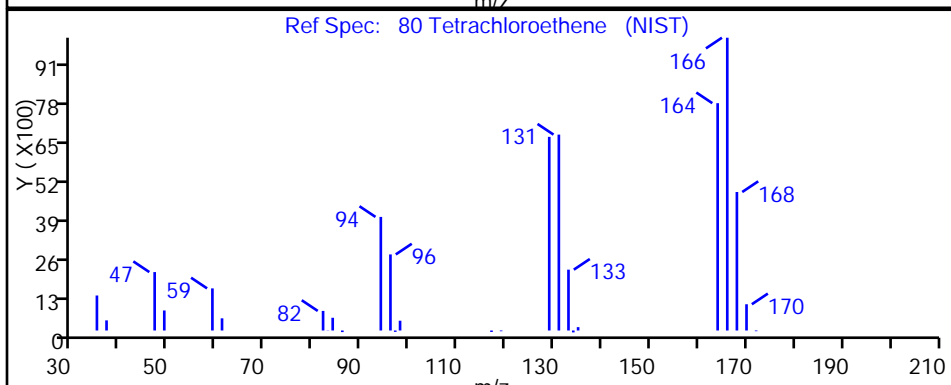
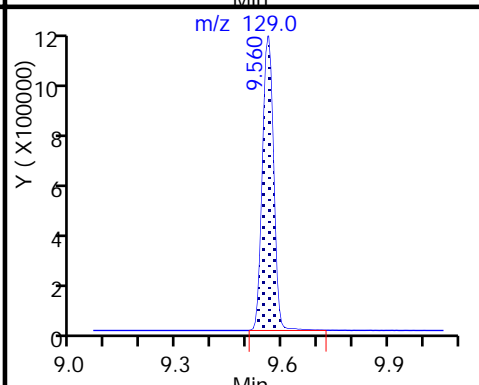
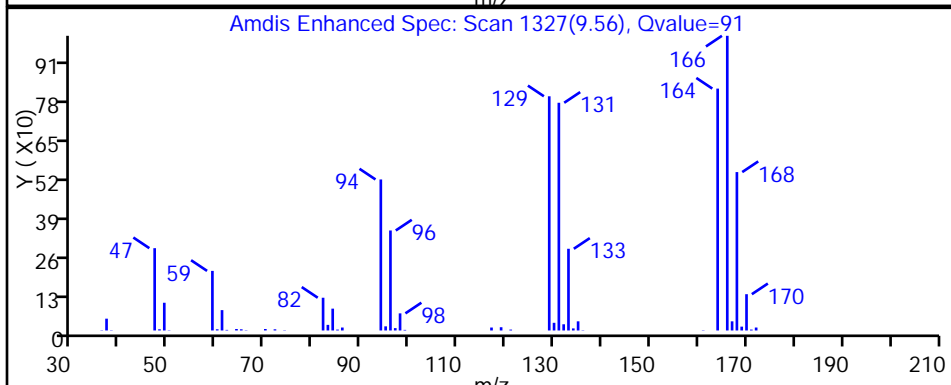
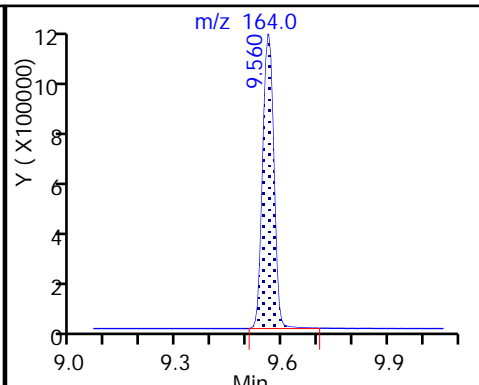
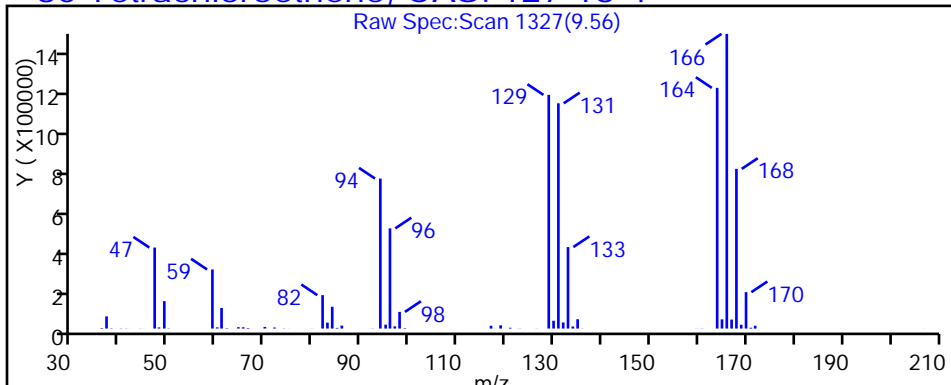
Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)

Detector: MS SCAN

80 Tetrachloroethene, CAS: 127-18-4



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Client Sample ID: HD-SPBA-CW-22-0/1-0 DL Lab Sample ID: 180-70873-1 DL
 Matrix: Water Lab File ID: 51004D16.D
 Analysis Method: 8260C Date Collected: 09/29/2017 09:40
 Sample wt/vol: 5 (mL) Date Analyzed: 10/05/2017 06:31
 Soil Aliquot Vol: _____ Dilution Factor: 12.5
 Soil Extract Vol.: _____ GC Column: DB-624 ID: 0.18 (mm)
 % Moisture: _____ Level: (low/med) Low
 Analysis Batch No.: 224919 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|------------|-----------------------------|--------|------|----|-----|
| 74-87-3 | Chloromethane | 13 | U ^c | 13 | 4.8 |
| 75-01-4 | Vinyl chloride | 13 | U ^c | 13 | 2.1 |
| 74-83-9 | Bromomethane | 13 | U | 13 | 7.3 |
| 75-00-3 | Chloroethane | 13 | U ^c | 13 | 7.2 |
| 75-35-4 | 1,1-Dichloroethene | 13 | U | 13 | 4.0 |
| 67-64-1 | Acetone | 63 | U | 63 | 39 |
| 75-15-0 | Carbon disulfide | 13 | U | 13 | 6.6 |
| 75-09-2 | Methylene Chloride | 13 | U | 13 | 12 |
| 156-60-5 | trans-1,2-Dichloroethene | 13 | U | 13 | 2.5 |
| 1634-04-4 | Methyl tert-butyl ether | 13 | U | 13 | 2.4 |
| 75-34-3 | 1,1-Dichloroethane | 13 | U | 13 | 4.2 |
| 156-59-2 | cis-1,2-Dichloroethene | 13 | U | 13 | 3.8 |
| 74-97-5 | Bromochloromethane | 13 | U | 13 | 4.5 |
| 78-93-3 | 2-Butanone (MEK) | 63 | U | 63 | 32 |
| 67-66-3 | Chloroform | 13 | U | 13 | 3.3 |
| 71-55-6 | 1,1,1-Trichloroethane | 13 | U | 13 | 3.4 |
| 56-23-5 | Carbon tetrachloride | 13 | U | 13 | 7.0 |
| 71-43-2 | Benzene | 13 | U | 13 | 2.3 |
| 107-06-2 | 1,2-Dichloroethane | 13 | U | 13 | 3.0 |
| 79-01-6 | Trichloroethene | 190 | | 13 | 2.5 |
| 78-87-5 | 1,2-Dichloropropane | 13 | U | 13 | 4.3 |
| 75-27-4 | Bromodichloromethane | 13 | U | 13 | 7.1 |
| 10061-01-5 | cis-1,3-Dichloropropene | 13 | U | 13 | 4.0 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 63 | U | 63 | 27 |
| 108-88-3 | Toluene | 13 | U | 13 | 2.0 |
| 10061-02-6 | trans-1,3-Dichloropropene | 13 | U | 13 | 2.8 |
| 79-00-5 | 1,1,2-Trichloroethane | 13 | U | 13 | 3.8 |
| 127-18-4 | Tetrachloroethene | 290 | | 13 | 3.1 |
| 591-78-6 | 2-Hexanone | 63 | U | 63 | 25 |
| 124-48-1 | Dibromochloromethane | 13 | U | 13 | 5.5 |
| 106-93-4 | 1,2-Dibromoethane (EDB) | 13 | U | 13 | 6.4 |
| 108-90-7 | Chlorobenzene | 13 | U | 13 | 1.8 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 13 | U | 13 | 6.2 |
| 100-41-4 | Ethylbenzene | 13 | U | 13 | 3.2 |
| 1330-20-7 | Xylenes, Total | 25 | U | 25 | 3.4 |
| 100-42-5 | Styrene | 13 | U | 13 | 2.7 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Client Sample ID: HD-SPBA-CW-22-0/1-0 DL Lab Sample ID: 180-70873-1 DL
 Matrix: Water Lab File ID: 51004D16.D
 Analysis Method: 8260C Date Collected: 09/29/2017 09:40
 Sample wt/vol: 5 (mL) Date Analyzed: 10/05/2017 06:31
 Soil Aliquot Vol: _____ Dilution Factor: 12.5
 Soil Extract Vol.: _____ GC Column: DB-624 ID: 0.18 (mm)
 % Moisture: _____ Level: (low/med) Low
 Analysis Batch No.: 224919 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|----------|----------------------------------|--------|---|------|-----|
| 75-25-2 | <i>Bromoform</i> | 13 | U | 13 | 9.5 |
| 79-34-5 | <i>1,1,2,2-Tetrachloroethane</i> | 13 | U | 13 | 4.6 |
| 107-13-1 | <i>Acrylonitrile</i> | 250 | U | 250 | 42 |
| 123-91-1 | <i>1,4-Dioxane</i> | 2500 | U | 2500 | 200 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|------------|------------------------------|------|---|--------|
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 117 | | 65-121 |
| 2037-26-5 | Toluene-d8 (Surr) | 99 | | 73-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 94 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 109 | | 73-120 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D16.D
 Lims ID: 180-70873-C-1
 Client ID: HD-SPBA-CW-22-0/1-0
 Sample Type: Client
 Inject. Date: 05-Oct-2017 06:31:30 ALS Bottle#: 16 Worklist Smp#: 16
 Purge Vol: 5.000 mL Dil. Factor: 12.5000
 Sample Info: 180-0018725-016
 Misc. Info.: 180-70873-C-1
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 08-Oct-2017 20:41:27 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK026

First Level Reviewer: bungardf

Date: 08-Oct-2017 20:36:16

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|----|----------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.361 | 4.352 | 0.009 | 0 | 149753 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.342 | 7.339 | 0.003 | 98 | 323921 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.426 | 10.429 | -0.003 | 86 | 71240 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.768 | 12.771 | -0.003 | 96 | 100827 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.618 | 6.613 | 0.005 | 92 | 85145 | 54.6 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.983 | 6.990 | -0.007 | 0 | 110735 | 58.3 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.979 | 8.979 | -0.001 | 93 | 280607 | 49.5 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.613 | 11.613 | 0.000 | 84 | 95981 | 46.9 | |
| 12 Chloromethane | 50 | | 1.825 | | | | ND | |
| 13 Vinyl chloride | 62 | | 1.959 | | | | ND | |
| 15 Bromomethane | 94 | | 2.300 | | | | ND | |
| 16 Chloroethane | 64 | | 2.470 | | | | ND | |
| 22 1,1-Dichloroethene | 96 | | 3.413 | | | | ND | |
| 24 Acetone | 43 | | 3.529 | | | | ND | |
| 26 Carbon disulfide | 76 | | 3.699 | | | | ND | |
| 31 Methylene Chloride | 84 | | 4.222 | | | | ND | |
| 33 Acrylonitrile | 53 | | 4.605 | | | | ND | |
| 34 trans-1,2-Dichloroethene | 96 | | 4.642 | | | | ND | |
| 35 Methyl tert-butyl ether | 73 | | 4.660 | | | | ND | |
| 37 1,1-Dichloroethane | 63 | | 5.268 | | | | ND | |
| 45 cis-1,2-Dichloroethene | 96 | | 6.011 | | | | ND | |
| 46 2-Butanone (MEK) | 43 | | 6.023 | | | | ND | |
| 49 Chlorobromomethane | 128 | | 6.290 | | | | ND | |
| 52 Chloroform | 83 | 6.436 | 6.436 | 0.000 | 1 | 667 | 0.2126 | |
| 53 1,1,1-Trichloroethane | 97 | | 6.595 | | | | ND | |
| 56 Carbon tetrachloride | 117 | | 6.759 | | | | ND | |
| 58 Benzene | 78 | | 6.996 | | | | ND | |
| 59 1,2-Dichloroethane | 62 | | 7.069 | | | | ND | |
| 64 Trichloroethene | 130 | 7.725 | 7.720 | 0.005 | 99 | 147589 | 74.5 | |
| 67 1,2-Dichloropropane | 63 | | 7.994 | | | | ND | |
| 70 1,4-Dioxane | 88 | | 8.085 | | | | ND | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Diff RT (min.) | Q | Response | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|----------------|----|----------|--------------|-------|
| 71 Dichlorobromomethane | 83 | | 8.280 | | | | ND | |
| 74 cis-1,3-Dichloropropene | 75 | | 8.718 | | | | ND | |
| 75 4-Methyl-2-pentanone (MIBK) | 43 | | 8.876 | | | | ND | |
| 76 Toluene | 91 | | 9.046 | | | | ND | |
| 77 trans-1,3-Dichloropropene | 75 | | 9.296 | | | | ND | |
| 79 1,1,2-Trichloroethane | 97 | | 9.490 | | | | ND | |
| 80 Tetrachloroethene | 164 | 9.563 | 9.563 | -0.001 | 94 | 158129 | 116.7 | |
| 82 2-Hexanone | 43 | | 9.703 | | | | ND | |
| 84 Chlorodibromomethane | 129 | | 9.855 | | | | ND | |
| 85 Ethylene Dibromide | 107 | | 9.971 | | | | ND | |
| 87 Chlorobenzene | 112 | | 10.458 | | | | ND | |
| 89 1,1,1,2-Tetrachloroethane | 131 | | 10.549 | | | | ND | |
| 90 Ethylbenzene | 106 | | 10.555 | | | | ND | |
| 91 m-Xylene & p-Xylene | 106 | | 10.689 | | | | ND | |
| 92 o-Xylene | 106 | | 11.072 | | | | ND | |
| 93 Styrene | 104 | | 11.090 | | | | ND | |
| 94 Bromoform | 173 | | 11.273 | | | | ND | |
| 99 1,1,2,2-Tetrachloroethane | 83 | | 11.747 | | | | ND | |
| S 133 Xylenes, Total | 106 | | 1.000 | | | | ND | |

Reagents:

VOA8260INT_00074

Amount Added: 2.00

Units: uL

Run Reagent

VOA8260SURR_00073

Amount Added: 2.00

Units: uL

Run Reagent

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D16.D

Injection Date: 05-Oct-2017 06:31:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: 180-70873-C-1

Lab Sample ID: 180-70873-1

Worklist Smp#: 16

Client ID: HD-SPBA-CW-22-0/1-0

Purge Vol: 5.000 mL

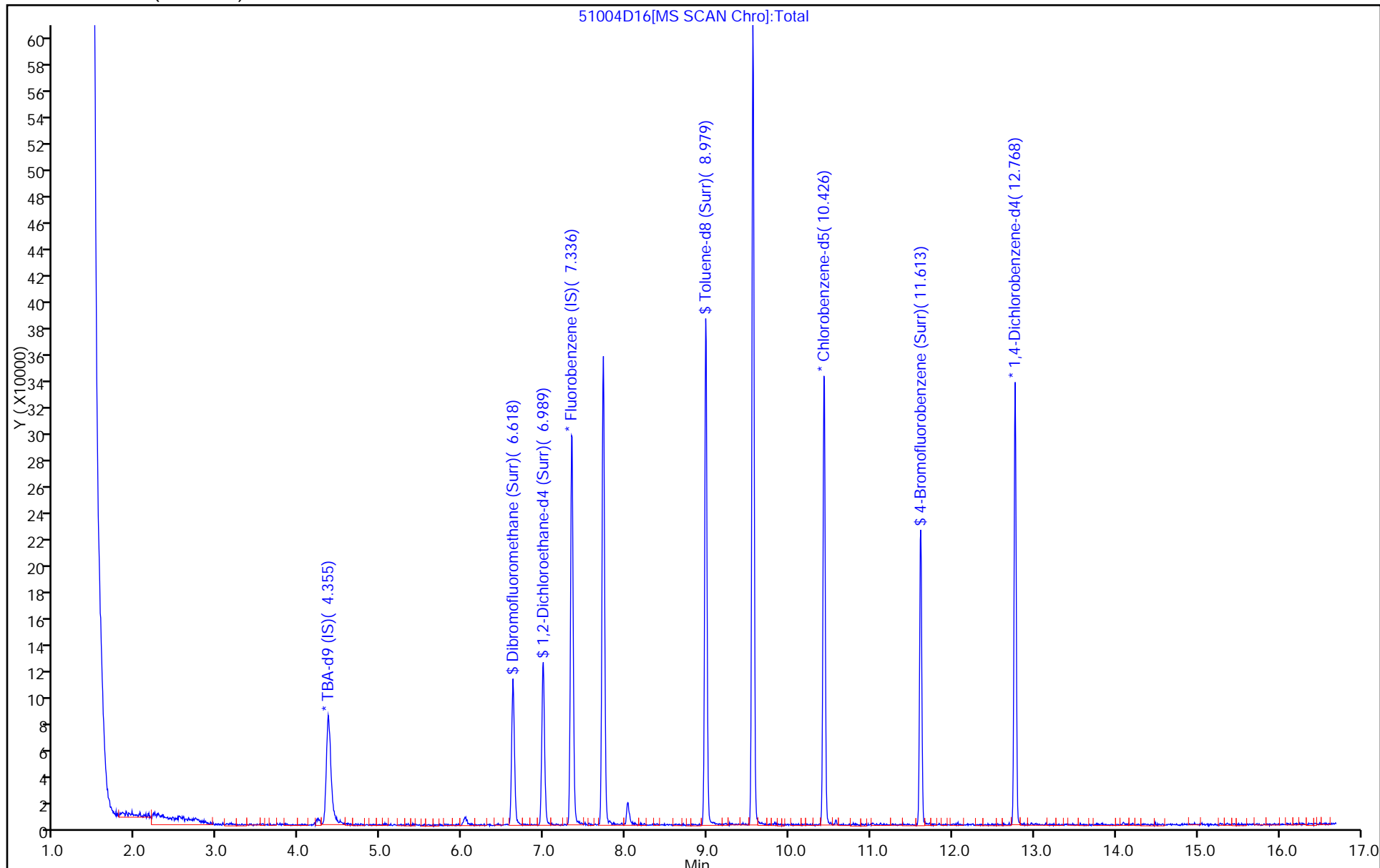
Dil. Factor: 12.5000

ALS Bottle#: 16

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Recovery Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D16.D
 Lims ID: 180-70873-C-1
 Client ID: HD-SPBA-CW-22-0/1-0
 Sample Type: Client
 Inject. Date: 05-Oct-2017 06:31:30 ALS Bottle#: 16 Worklist Smp#: 16
 Purge Vol: 5.000 mL Dil. Factor: 12.5000
 Sample Info: 180-0018725-016
 Misc. Info.: 180-70873-C-1
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 08-Oct-2017 20:41:27 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK026

First Level Reviewer: bungardf

Date: 08-Oct-2017 20:36:16

| Compound | Amount Added | Amount Recovered | % Rec. |
|-----------------------------------|--------------|------------------|--------|
| \$ 5 Dibromofluoromethane (Surr) | 50.0 | 54.6 | 109.26 |
| \$ 6 1,2-Dichloroethane-d4 (Surr) | 50.0 | 58.3 | 116.51 |
| \$ 7 Toluene-d8 (Surr) | 50.0 | 49.5 | 98.98 |
| \$ 8 4-Bromofluorobenzene (Surr) | 50.0 | 46.9 | 93.74 |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D16.D

Injection Date: 05-Oct-2017 06:31:30

Instrument ID: CHHP5

Lims ID: 180-70873-C-1

Lab Sample ID: 180-70873-1

Client ID: HD-SPBA-CW-22-0/1-0

Operator ID: 034635

ALS Bottle#: 16

Worklist Smp#: 16

Purge Vol: 5.000 mL

Dil. Factor: 12.5000

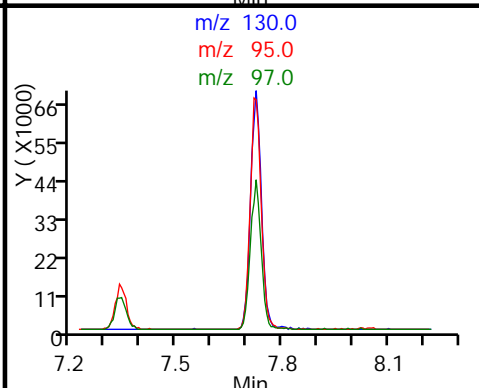
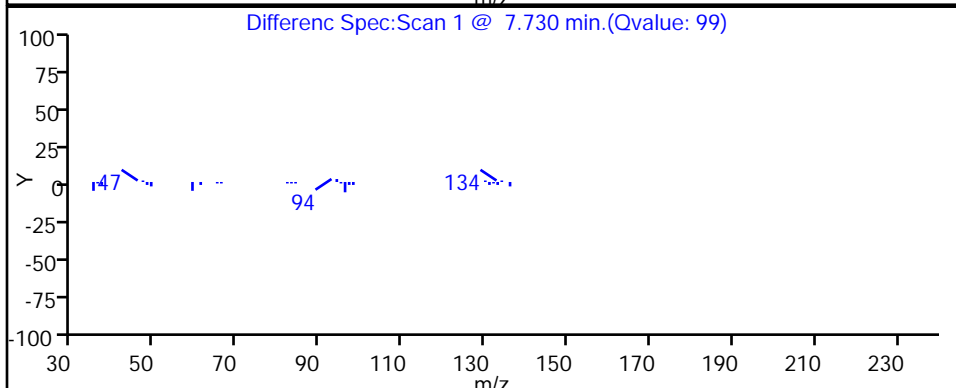
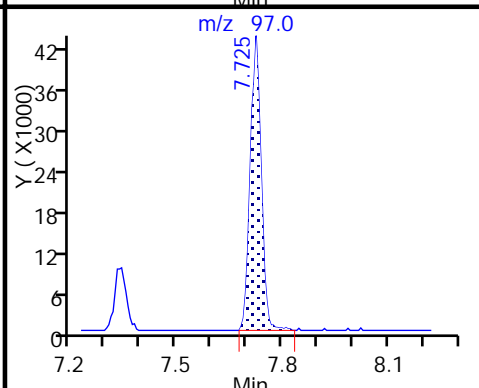
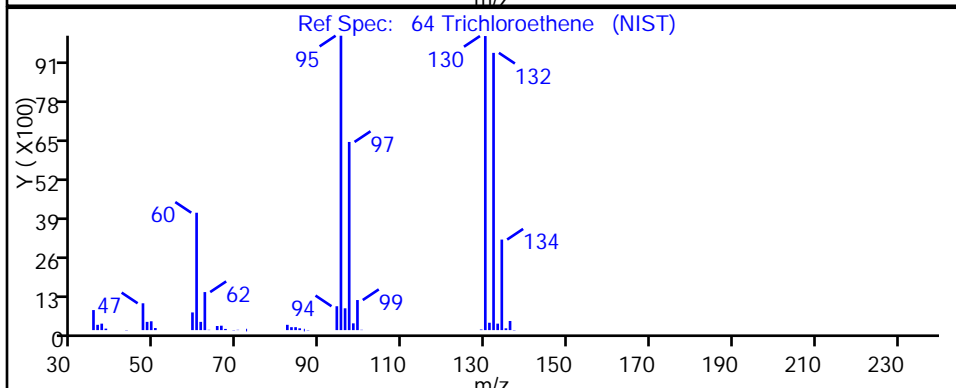
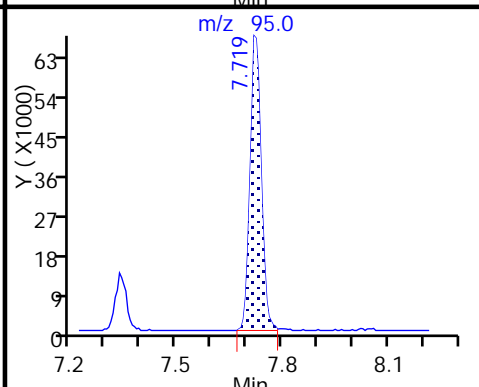
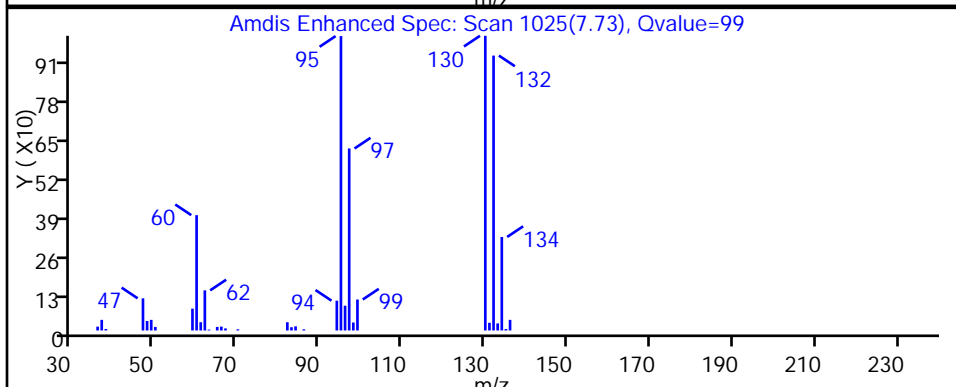
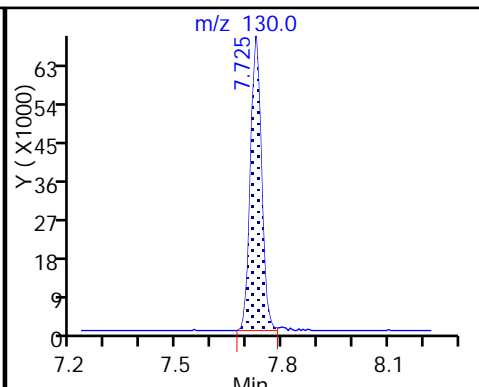
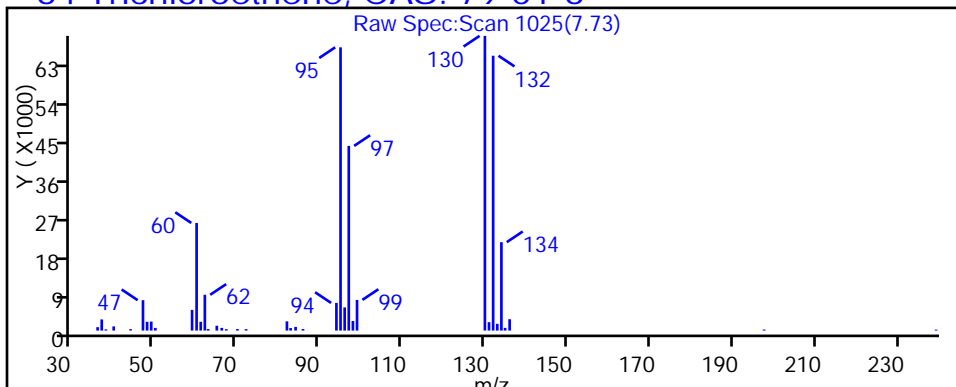
Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)

Detector: MS SCAN

64 Trichloroethene, CAS: 79-01-6



TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D16.D

Injection Date: 05-Oct-2017 06:31:30

Instrument ID: CHHP5

Lims ID: 180-70873-C-1

Lab Sample ID: 180-70873-1

Client ID: HD-SPBA-CW-22-0/1-0

Operator ID: 034635

ALS Bottle#: 16

Worklist Smp#: 16

Purge Vol: 5.000 mL

Dil. Factor: 12.5000

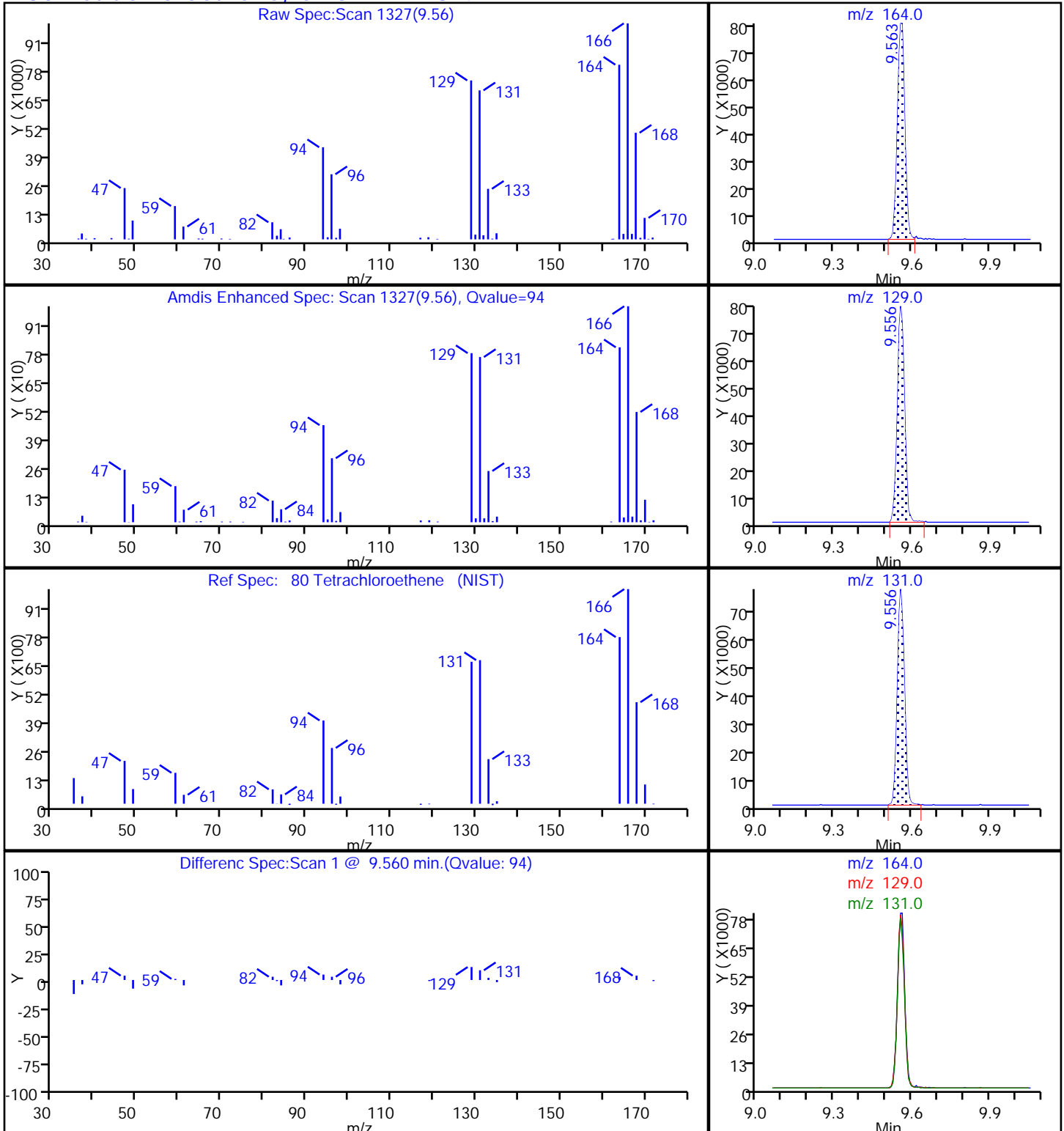
Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)

Detector: MS SCAN

80 Tetrachloroethene, CAS: 127-18-4



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Client Sample ID: HD-QC4-0/1-2 Lab Sample ID: 180-70873-2
 Matrix: Water Lab File ID: 51003D12.D
 Analysis Method: 8260C Date Collected: 09/29/2017 12:00
 Sample wt/vol: 5 (mL) Date Analyzed: 10/04/2017 04:59
 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.: 224792 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|------------|-----------------------------|--------|------|-----|------|
| 74-87-3 | Chloromethane | 1.0 | U ^c | 1.0 | 0.38 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 1.0 | 0.17 |
| 74-83-9 | Bromomethane | 1.0 | U | 1.0 | 0.59 |
| 75-00-3 | Chloroethane | 1.0 | U | 1.0 | 0.58 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.32 |
| 67-64-1 | Acetone | 5.0 | U | 5.0 | 3.1 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 1.0 | 0.53 |
| 75-09-2 | Methylene Chloride | 1.0 | U | 1.0 | 0.94 |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.20 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.34 |
| 156-59-2 | cis-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.30 |
| 74-97-5 | Bromochloromethane | 1.0 | U | 1.0 | 0.36 |
| 78-93-3 | 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 |
| 67-66-3 | Chloroform | 1.0 | U | 1.0 | 0.27 |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.27 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 1.0 | 0.56 |
| 71-43-2 | Benzene | 1.0 | U | 1.0 | 0.18 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.24 |
| 79-01-6 | Trichloroethene | 1.0 | U | 1.0 | 0.20 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.35 |
| 75-27-4 | Bromodichloromethane | 1.0 | U ^c | 1.0 | 0.57 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U ^c | 1.0 | 0.32 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 5.0 | U | 5.0 | 2.2 |
| 108-88-3 | Toluene | 1.0 | U | 1.0 | 0.16 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.22 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.31 |
| 127-18-4 | Tetrachloroethene | 1.0 | U | 1.0 | 0.24 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 5.0 | 2.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 1.0 | 0.44 |
| 106-93-4 | 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.51 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 1.0 | 0.15 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.49 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 1.0 | 0.25 |
| 1330-20-7 | Xylenes, Total | 2.0 | U | 2.0 | 0.27 |
| 100-42-5 | Styrene | 1.0 | U | 1.0 | 0.22 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Client Sample ID: HD-QC4-0/1-2 Lab Sample ID: 180-70873-2
 Matrix: Water Lab File ID: 51003D12.D
 Analysis Method: 8260C Date Collected: 09/29/2017 12:00
 Sample wt/vol: 5 (mL) Date Analyzed: 10/04/2017 04:59
 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.: 224792 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|----------|---------------------------|--------|------|-----|------|
| 75-25-2 | Bromoform | 1.0 | U ^c | 1.0 | 0.76 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.37 |
| 107-13-1 | Acrylonitrile | 20 | U | 20 | 3.3 |
| 123-91-1 | 1,4-Dioxane | 200 | U | 200 | 16 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|------------|------------------------------|------|---|--------|
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 113 | | 65-121 |
| 2037-26-5 | Toluene-d8 (Surr) | 101 | | 73-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 92 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 73-120 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D12.D
 Lims ID: 180-70873-A-2
 Client ID: HD-QC4-0/1-2
 Sample Type: Client
 Inject. Date: 04-Oct-2017 04:59:30 ALS Bottle#: 12 Worklist Smp#: 12
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018710-012
 Misc. Info.: 180-70873-A-2
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 04-Oct-2017 21:10:21 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK005

First Level Reviewer: bungardf

Date: 04-Oct-2017 20:54:30

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|----|----------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.358 | 4.347 | 0.011 | 0 | 177857 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.339 | 7.334 | 0.005 | 99 | 356184 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.430 | 10.431 | -0.001 | 86 | 75850 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.772 | 12.773 | -0.001 | 96 | 109308 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.622 | 6.616 | 0.006 | 94 | 86590 | 50.5 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.993 | 6.981 | 0.012 | 0 | 118160 | 56.5 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.982 | 8.977 | 0.005 | 93 | 304177 | 50.4 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.610 | 11.611 | -0.001 | 86 | 100564 | 46.1 | |
| 12 Chloromethane | 50 | | 1.823 | | | | ND | |
| 13 Vinyl chloride | 62 | | 1.963 | | | | ND | |
| 15 Bromomethane | 94 | | 2.291 | | | | ND | |
| 16 Chloroethane | 64 | | 2.461 | | | | ND | |
| 22 1,1-Dichloroethene | 96 | | 3.423 | | | | ND | |
| 24 Acetone | 43 | 3.543 | 3.526 | 0.017 | 93 | 11198 | 12.0 | |
| 26 Carbon disulfide | 76 | | 3.708 | | | | ND | |
| 31 Methylene Chloride | 84 | | 4.232 | | | | ND | |
| 33 Acrylonitrile | 53 | | 4.609 | | | | ND | |
| 34 trans-1,2-Dichloroethene | 96 | | 4.633 | | | | ND | |
| 35 Methyl tert-butyl ether | 73 | | 4.657 | | | | ND | |
| 37 1,1-Dichloroethane | 63 | | 5.272 | | | | ND | |
| 45 cis-1,2-Dichloroethene | 96 | | 6.008 | | | | ND | |
| 46 2-Butanone (MEK) | 43 | | 6.026 | | | | ND | |
| 49 Chlorobromomethane | 128 | | 6.294 | | | | ND | |
| 52 Chloroform | 83 | | 6.440 | | | | ND | |
| 53 1,1,1-Trichloroethane | 97 | | 6.592 | | | | ND | |
| 56 Carbon tetrachloride | 117 | | 6.762 | | | | ND | |
| 58 Benzene | 78 | | 6.994 | | | | ND | |
| 59 1,2-Dichloroethane | 62 | | 7.067 | | | | ND | |
| 64 Trichloroethene | 130 | | 7.724 | | | | ND | |
| 67 1,2-Dichloropropane | 63 | | 7.997 | | | | ND | |
| 70 1,4-Dioxane | 88 | | 8.082 | | | | ND | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|---------------|---|----------|--------------|-------|
| 71 Dichlorobromomethane | 83 | | 8.277 | | | | ND | |
| 74 cis-1,3-Dichloropropene | 75 | | 8.721 | | | | ND | |
| 75 4-Methyl-2-pentanone (MIBK) | 43 | | 8.873 | | | | ND | |
| 76 Toluene | 91 | | 9.044 | | | | ND | |
| 77 trans-1,3-Dichloropropene | 75 | | 9.293 | | | | ND | |
| 79 1,1,2-Trichloroethane | 97 | | 9.488 | | | | ND | |
| 80 Tetrachloroethene | 164 | | 9.561 | | | | ND | |
| 82 2-Hexanone | 43 | | 9.707 | | | | ND | |
| 84 Chlorodibromomethane | 129 | | 9.853 | | | | ND | |
| 85 Ethylene Dibromide | 107 | | 9.974 | | | | ND | |
| 87 Chlorobenzene | 112 | | 10.455 | | | | ND | |
| 89 1,1,1,2-Tetrachloroethane | 131 | | 10.552 | | | | ND | |
| 90 Ethylbenzene | 106 | | 10.558 | | | | ND | |
| 91 m-Xylene & p-Xylene | 106 | | 10.692 | | | | ND | |
| 92 o-Xylene | 106 | | 11.069 | | | | ND | |
| 93 Styrene | 104 | | 11.094 | | | | ND | |
| 94 Bromoform | 173 | | 11.270 | | | | ND | |
| 99 1,1,2,2-Tetrachloroethane | 83 | | 11.751 | | | | ND | |
| S 133 Xylenes, Total | 106 | | 1.000 | | | | ND | |

Reagents:

VOA8260INT_00074

Amount Added: 2.00

Units: uL

Run Reagent

VOA8260SURR_00073

Amount Added: 2.00

Units: uL

Run Reagent

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D12.D

Injection Date: 04-Oct-2017 04:59:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: 180-70873-A-2

Lab Sample ID: 180-70873-2

Worklist Smp#: 12

Client ID: HD-QC4-0/1-2

Purge Vol: 5.000 mL

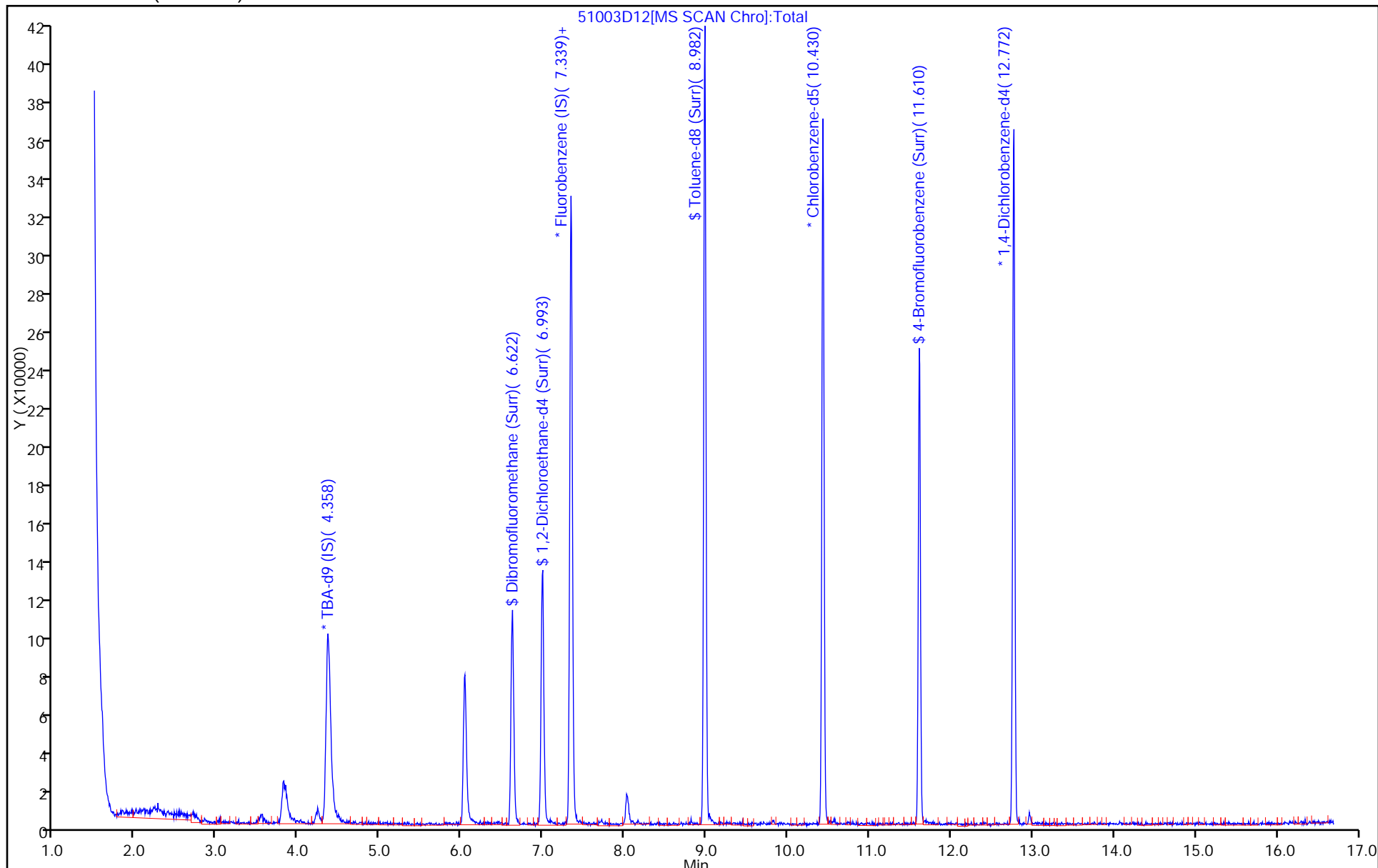
Dil. Factor: 1.0000

ALS Bottle#: 12

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Recovery Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D12.D
 Lims ID: 180-70873-A-2
 Client ID: HD-QC4-0/1-2
 Sample Type: Client
 Inject. Date: 04-Oct-2017 04:59:30 ALS Bottle#: 12 Worklist Smp#: 12
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018710-012
 Misc. Info.: 180-70873-A-2
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 04-Oct-2017 21:10:21 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK005

First Level Reviewer: bungardf Date: 04-Oct-2017 20:54:30

| Compound | Amount Added | Amount Recovered | % Rec. |
|-----------------------------------|--------------|------------------|--------|
| \$ 5 Dibromofluoromethane (Surr) | 50.0 | 50.5 | 101.05 |
| \$ 6 1,2-Dichloroethane-d4 (Surr) | 50.0 | 56.5 | 113.06 |
| \$ 7 Toluene-d8 (Surr) | 50.0 | 50.4 | 100.78 |
| \$ 8 4-Bromofluorobenzene (Surr) | 50.0 | 46.1 | 92.25 |

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

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1 Analy Batch No.: 218218

SDG No.: _____

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

Calibration Start Date: 07/27/2017 00:51 Calibration End Date: 07/27/2017 04:24 Calibration ID: 35038

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---------|-------------------|--------------|
| Level 1 | IC 180-218218/2 | 50727D02.D |
| Level 2 | IC 180-218218/3 | 50727D03.D |
| Level 3 | ICIS 180-218218/4 | 50727D04.D |
| Level 4 | IC 180-218218/5 | 50727D05.D |
| Level 5 | IC 180-218218/6 | 50727D06.D |
| Level 6 | IC 180-218218/10 | 50727D10.D |
| Level 7 | IC 180-218218/8 | 50727D08.D |
| Level 8 | IC 180-218218/11 | 50727D11.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---------------------------------------|------------------|------------------|------------------|--------|--------|------------|-------------|--------|----|--------|---------|------|------|----------|------------|---|----------------|
| | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| | LVL 6 | LVL 7 | LVL 8 | | | | | | | | | | | | | | |
| Dichlorodifluoromethane | 0.3099 0.3034 | 0.3143 0.2538 | 0.2964 0.2820 | 0.2910 | 0.2753 | Ave | | 0.2907 | | 0.1000 | 6.9 | | 20.0 | | | | |
| Chloromethane | 0.3638 0.2790 | 0.2935 0.2586 | 0.2871 0.2672 | 0.2979 | 0.2905 | Ave | | 0.2922 | | 0.1000 | 10.9 | | 20.0 | | | | |
| Vinyl chloride | 0.3612 0.2960 | 0.3073 0.2570 | 0.3014 0.2855 | 0.2838 | 0.2802 | Ave | | 0.2965 | | 0.1000 | 10.2 | | 20.0 | | | | |
| 1,3-Butadiene | 0.3317 0.2714 | 0.2771 0.2281 | 0.2660 0.2684 | 0.2619 | 0.2505 | Ave | | 0.2694 | | 0.0100 | 10.9 | | 20.0 | | | | |
| Bromomethane | 0.1274 0.1338 | 0.1569 0.1290 | 0.1507 0.1244 | 0.1438 | 0.1556 | Ave | | 0.1402 | | 0.0500 | 9.4 | | 20.0 | | | | |
| Chloroethane | 0.1972 0.1593 | 0.1757 0.1437 | 0.1605 0.1363 | 0.1653 | 0.1659 | Ave | | 0.1630 | | 0.0500 | 11.5 | | 20.0 | | | | |
| Trichlorofluoromethane | 0.4130 0.3605 | 0.3896 0.3164 | 0.3801 0.3348 | 0.3631 | 0.3573 | Ave | | 0.3643 | | 0.1000 | 8.4 | | 20.0 | | | | |
| Ethyl ether | 0.2690 0.2226 | 0.2473 0.2272 | 0.2344 0.2016 | 0.2419 | 0.2520 | Ave | | 0.2370 | | 0.0100 | 8.6 | | 20.0 | | | | |
| Acrolein | 0.0588 0.0564 | 0.0546 0.0639 | 0.0629 0.0550 | 0.0633 | 0.0629 | Ave | | 0.0597 | | 0.0100 | 6.7 | | 20.0 | | | | |
| 1,1-Dichloroethene | 0.2633 0.2529 | 0.2525 0.2180 | 0.2438 0.2452 | 0.2449 | 0.2377 | Ave | | 0.2448 | | 0.1000 | 5.4 | | 20.0 | | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.3346 0.2678 | 0.2745 0.2382 | 0.2615 0.2547 | 0.2644 | 0.2534 | Ave | | 0.2686 | | 0.1000 | 10.7 | | 20.0 | | | | |
| Acetone | 0.1396 0.1048 | 0.1447 0.1163 | 0.1388 0.1038 | 0.1460 | 0.1519 | Ave | | 0.1308 | | 0.0500 | 14.8 | | 20.0 | | | | |
| Iodomethane | 0.4213 0.3803 | 0.3860 0.3716 | 0.3712 0.3619 | 0.3906 | 0.3928 | Ave | | 0.3845 | | 0.0100 | 4.8 | | 20.0 | | | | |
| Carbon disulfide | 0.5698 ++++ | 0.4896 0.5397 | 0.4946 0.6108 | 0.5168 | 0.5392 | Ave | | 0.5372 | | 0.1000 | 8.0 | | 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.: 180-70873-1 Analy Batch No.: 218218

SDG No.: _____

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

Calibration Start Date: 07/27/2017 00:51 Calibration End Date: 07/27/2017 04:24 Calibration ID: 35038

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|--------------------------|------------------|------------------|------------------|--------|--------|------------|-------------|--------|----|--------|---------|------|------|----------|------------|--------|----------------|
| | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| | LVL 6 | LVL 7 | LVL 8 | LVL 5 | | | | | | | | | | | | | |
| Allyl chloride | 0.1501 0.1710 | 0.1485 0.1632 | 0.1541 0.1645 | 0.1561 | 0.1579 | Ave | | 0.1582 | | 0.0100 | 4.8 | | 20.0 | | | | |
| Methyl acetate | 0.2888 0.2364 | 0.2463 0.2614 | 0.2631 0.2382 | 0.2688 | 0.2686 | Ave | | 0.2589 | | 0.1000 | 6.8 | | 20.0 | | | | |
| Methylene Chloride | 0.4748 0.2821 | 0.3152 0.2910 | 0.3044 0.2676 | 0.3112 | 0.3108 | Lin2 | 0.9532 | 0.2841 | | 0.1000 | | | | 0.9980 | | 0.9900 | |
| tert-Butyl alcohol | 1.3346 1.2872 | 1.1570 1.0277 | 1.1638 1.2343 | 1.1314 | 1.1253 | Ave | | 1.1826 | | 0.0100 | 8.3 | | 20.0 | | | | |
| Acrylonitrile | 0.1353 0.1106 | 0.1251 0.1245 | 0.1313 0.1150 | 0.1320 | 0.1333 | Ave | | 0.1259 | | 0.0100 | 7.1 | | 20.0 | | | | |
| trans-1,2-Dichloroethene | 0.3167 0.2789 | 0.2730 0.2547 | 0.2727 0.2653 | 0.2850 | 0.2851 | Ave | | 0.2789 | | 0.1000 | 6.6 | | 20.0 | | | | |
| Methyl tert-butyl ether | 0.7081 0.7482 | 0.7314 0.7800 | 0.7230 0.7142 | 0.7872 | 0.7909 | Ave | | 0.7479 | | 0.1000 | 4.5 | | 20.0 | | | | |
| Hexane | 0.4597 0.3561 | 0.3588 0.3156 | 0.3449 0.3625 | 0.3424 | 0.3242 | Ave | | 0.3580 | | 0.0100 | 12.4 | | 20.0 | | | | |
| 1,1-Dichloroethane | 0.5228 0.4797 | 0.4979 0.4638 | 0.4852 0.4528 | 0.4864 | 0.4910 | Ave | | 0.4850 | | 0.2000 | 4.4 | | 20.0 | | | | |
| Vinyl acetate | 0.5018 0.5003 | 0.4274 0.5345 | 0.4556 0.5012 | 0.5130 | 0.5116 | Ave | | 0.4932 | | 0.0100 | 7.0 | | 20.0 | | | | |
| 2,2-Dichloropropane | 0.0696 0.0640 | 0.0591 0.0559 | 0.0577 0.0619 | 0.0627 | 0.0632 | Ave | | 0.0617 | | 0.0100 | 6.9 | | 20.0 | | | | |
| cis-1,2-Dichloroethene | 0.3297 0.3143 | 0.3194 0.3060 | 0.3200 0.2963 | 0.3326 | 0.3338 | Ave | | 0.3190 | | 0.1000 | 4.1 | | 20.0 | | | | |
| 2-Butanone (MEK) | 0.1854 0.1607 | 0.1969 0.1772 | 0.1989 0.1584 | 0.2064 | 0.2051 | Ave | | 0.1861 | | 0.0500 | 10.2 | | 20.0 | | | | |
| Bromochloromethane | 0.1517 0.1366 | 0.1414 0.1398 | 0.1402 0.1299 | 0.1453 | 0.1494 | Ave | | 0.1418 | | 0.0100 | 4.9 | | 20.0 | | | | |
| Tetrahydrofuran | 0.1371 0.0928 | 0.0982 0.1088 | 0.1088 0.1003 | 0.1130 | 0.1079 | Ave | | 0.1084 | | 0.0100 | 12.4 | | 20.0 | | | | |
| Chloroform | 0.5466 0.4636 | 0.4996 0.4621 | 0.4713 0.4342 | 0.4992 | 0.4977 | Ave | | 0.4843 | | 0.2000 | 7.0 | | 20.0 | | | | |
| 1,1,1-Trichloroethane | 0.3786 0.3800 | 0.3677 0.3465 | 0.3637 0.3610 | 0.3661 | 0.3690 | Ave | | 0.3666 | | 0.1000 | 2.9 | | 20.0 | | | | |
| Cyclohexane | 0.4979 0.4744 | 0.4616 0.4108 | 0.4435 0.4590 | 0.4424 | 0.4292 | Ave | | 0.4524 | | 0.1000 | 6.0 | | 20.0 | | | | |
| Carbon tetrachloride | 0.3181 0.3198 | 0.2990 0.2880 | 0.3018 0.3038 | 0.3054 | 0.3047 | Ave | | 0.3051 | | 0.1000 | 3.3 | | 20.0 | | | | |
| 1,1-Dichloropropene | 0.4064 0.4059 | 0.4083 0.3679 | 0.3990 0.3876 | 0.4006 | 0.3928 | Ave | | 0.3961 | | 0.0100 | 3.4 | | 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.: 180-70873-1 Analy Batch No.: 218218

SDG No.: _____

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

Calibration Start Date: 07/27/2017 00:51 Calibration End Date: 07/27/2017 04:24 Calibration ID: 35038

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R ² OR COD | # | MIN R ² OR COD |
|-----------------------------|------------------|------------------|------------------|--------|--------|------------|-------------|--------|----|---|---------|------|------|----------|-----------------------|---|---------------------------|
| | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| | LVL 6 | LVL 7 | LVL 8 | | | | | | | | | | | | | | |
| Isobutyl alcohol | 0.0097 0.0085 | 0.0091 0.0105 | 0.0102 0.0094 | 0.0111 | 0.0112 | Ave | | 0.0099 | | * | 0.0100 | 9.6 | 20.0 | | | | |
| Benzene | 1.3787 1.1520 | 1.2628 1.1081 | 1.2398 1.0692 | 1.2590 | 1.2563 | Ave | | 1.2157 | | | 0.5000 | 8.2 | 20.0 | | | | |
| 1,2-Dichloroethane | 0.3884 0.3320 | 0.3554 0.3421 | 0.3528 0.3189 | 0.3753 | 0.3703 | Ave | | 0.3544 | | | 0.1000 | 6.5 | 20.0 | | | | |
| n-Heptane | 0.3037 0.2967 | 0.3011 0.2552 | 0.2860 0.3036 | 0.2755 | 0.2684 | Ave | | 0.2863 | | | 0.0100 | 6.4 | 20.0 | | | | |
| Trichloroethene | 0.3229 0.3036 | 0.3087 0.2884 | 0.3052 0.2920 | 0.3101 | 0.3167 | Ave | | 0.3059 | | | 0.2000 | 3.8 | 20.0 | | | | |
| Methylcyclohexane | 0.4727 0.4875 | 0.4672 0.4232 | 0.4697 0.4715 | 0.4601 | 0.4491 | Ave | | 0.4626 | | | 0.1000 | 4.2 | 20.0 | | | | |
| 1,2-Dichloropropane | 0.3012 0.2794 | 0.2779 0.2782 | 0.2782 0.2612 | 0.2913 | 0.2975 | Ave | | 0.2831 | | | 0.1000 | 4.6 | 20.0 | | | | |
| 1,4-Dioxane | 0.0022 0.0027 | 0.0028 0.0030 | 0.0031 0.0031 | 0.0030 | 0.0032 | Ave | | 0.0029 | | * | 0.0100 | 11.4 | 20.0 | | | | |
| Dibromomethane | 0.1595 0.1606 | 0.1708 0.1667 | 0.1638 0.1549 | 0.1734 | 0.1774 | Ave | | 0.1659 | | | 0.0100 | 4.6 | 20.0 | | | | |
| Bromodichloromethane | 0.3001 0.3336 | 0.3125 0.3351 | 0.3169 0.3110 | 0.3438 | 0.3519 | Ave | | 0.3256 | | | 0.2000 | 5.6 | 20.0 | | | | |
| 2-Chloroethyl vinyl ether | 0.1669 0.2025 | 0.1917 0.2176 | 0.2032 0.2031 | 0.2200 | 0.2248 | Ave | | 0.2037 | | | 0.0100 | 9.1 | 20.0 | | | | |
| cis-1,3-Dichloropropene | 0.3596 0.4128 | 0.3596 0.4158 | 0.3786 0.3959 | 0.4116 | 0.4298 | Ave | | 0.3955 | | | 0.2000 | 6.8 | 20.0 | | | | |
| 4-Methyl-2-pentanone (MIBK) | 1.3560 1.1652 | 1.2491 1.2232 | 1.3592 1.1532 | 1.3610 | 1.3926 | Ave | | 1.2824 | | | 0.1000 | 7.5 | 20.0 | | | | |
| Toluene | 6.1005 4.5990 | 5.6903 4.2081 | 5.2159 4.0277 | 5.0185 | 5.0243 | Ave | | 4.9855 | | | 0.4000 | 14.1 | 20.0 | | | | |
| trans-1,3-Dichloropropene | 1.2257 1.4397 | 1.2796 1.4086 | 1.2851 1.3247 | 1.3956 | 1.4937 | Ave | | 1.3566 | | | 0.1000 | 6.8 | 20.0 | | | | |
| Ethyl methacrylate | 1.3604 1.6673 | 1.5623 1.6591 | 1.6724 1.5738 | 1.7698 | 1.8222 | Ave | | 1.6359 | | | 0.0100 | 8.7 | 20.0 | | | | |
| 1,1,2-Trichloroethane | 1.2522 0.9633 | 1.0992 0.9427 | 1.0403 0.8887 | 1.0530 | 1.0694 | Ave | | 1.0386 | | | 0.1000 | 10.8 | 20.0 | | | | |
| Tetrachloroethene | 1.1481 0.9182 | 1.0929 0.8058 | 0.9505 0.8459 | 0.9238 | 0.9211 | Ave | | 0.9508 | | | 0.2000 | 12.2 | 20.0 | | | | |
| 1,3-Dichloropropane | 2.2370 1.7852 | 2.0694 1.7532 | 1.9307 1.6348 | 1.9958 | 1.9532 | Ave | | 1.9199 | | | 0.0100 | 10.0 | 20.0 | | | | |
| 2-Hexanone | 0.9818 0.8998 | 0.9941 0.9190 | 1.0485 0.8780 | 1.0518 | 1.0958 | Ave | | 0.9836 | | | 0.1000 | 8.1 | 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.: 180-70873-1

Analy Batch No.: 218218

SDG No.: _____

Instrument ID: CHHP5

GC Column: DB-624

ID: 0.18 (mm)

Heated Purge: (Y/N) N

Calibration Start Date: 07/27/2017 00:51

Calibration End Date: 07/27/2017 04:24

Calibration ID: 35038

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R ² OR COD | # | MIN R ² OR COD |
|-----------------------------|------------------|------------------|------------------|--------|--------|------------|-------------|--------|----|---|---------|------|------|----------|-----------------------|---|---------------------------|
| | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| | LVL 6 | LVL 7 | LVL 8 | | | | | | | | | | | | | | |
| Dibromochloromethane | 0.7989 0.9016 | 0.8620 0.8947 | 0.8650 0.8322 | 0.9093 | 0.9598 | Ave | | 0.8779 | | | 0.1000 | 5.7 | 20.0 | | | | |
| 1,2-Dibromoethane (EDB) | 1.1425 1.0146 | 1.0956 1.0059 | 1.0726 0.9575 | 1.1227 | 1.1100 | Ave | | 1.0652 | | | 0.1000 | 6.1 | 20.0 | | | | |
| 3-Chlorobenzotrifluoride | 2.1508 1.6103 | 1.7646 1.4397 | 1.6777 1.5967 | 1.7670 | 1.7382 | Ave | | 1.7181 | | | 0.0100 | 12.0 | 20.0 | | | | |
| Chlorobenzene | 4.0368 3.0317 | 3.5186 2.8231 | 3.2468 2.6869 | 3.3119 | 3.3091 | Ave | | 3.2456 | | | 0.5000 | 13.0 | 20.0 | | | | |
| 4-Chlorobenzotrifluoride | 1.8614 1.5230 | 1.6468 1.3432 | 1.5641 1.5178 | 1.6419 | 1.5859 | Ave | | 1.5855 | | | 0.0100 | 9.3 | 20.0 | | | | |
| 1,1,1,2-Tetrachloroethane | 1.0682 1.0211 | 1.0658 0.9781 | 1.0366 0.9303 | 1.0666 | 1.0896 | Ave | | 1.0321 | | | 0.0100 | 5.2 | 20.0 | | | | |
| Ethylbenzene | 1.9199 1.7723 | 1.9530 1.6113 | 1.8804 1.6150 | 1.8616 | 1.8815 | Ave | | 1.8119 | | | 0.1000 | 7.3 | 20.0 | | | | |
| m-Xylene & p-Xylene | 2.1686 2.2054 | 2.4439 2.0173 | 2.3106 1.9980 | 2.2675 | 2.3006 | Ave | | 2.2140 | | | 0.1000 | 6.8 | 20.0 | | | | |
| o-Xylene | 2.1421 2.0826 | 2.2379 1.9206 | 2.1746 1.8793 | 2.2085 | 2.2321 | Ave | | 2.1097 | | | 0.3000 | 6.6 | 20.0 | | | | |
| Styrene | 3.6332 3.4371 | 3.9143 3.2595 | 3.7554 3.0478 | 3.7413 | 3.7778 | Ave | | 3.5708 | | | 0.3000 | 8.3 | 20.0 | | | | |
| Bromoform | 0.5105 0.5727 | 0.4852 0.5813 | 0.5106 0.5484 | 0.5622 | 0.5938 | Ave | | 0.5456 | | | 0.1000 | 7.2 | 20.0 | | | | |
| 2-Chlorobenzotrifluoride | 1.7885 1.5489 | 1.7322 1.4506 | 1.6281 1.5406 | 1.7502 | 1.7146 | Ave | | 1.6442 | | | 0.0100 | 7.4 | 20.0 | | | | |
| Isopropylbenzene | 5.5110 4.9386 | 5.7732 4.4163 | 5.4683 4.3345 | 5.4199 | 5.3367 | Ave | | 5.1498 | | | 0.1000 | 10.3 | 20.0 | | | | |
| Bromobenzene | 0.9987 0.9743 | 0.9872 0.9390 | 0.9377 0.9146 | 0.9980 | 1.0140 | Ave | | 0.9704 | | | 0.0100 | 3.7 | 20.0 | | | | |
| 1,1,2,2-Tetrachloroethane | 1.7609 1.4046 | 1.6228 1.4415 | 1.5952 1.3351 | 1.5862 | 1.5551 | Ave | | 1.5377 | | | 0.3000 | 8.9 | 20.0 | | | | |
| trans-1,4-Dichloro-2-butene | 0.2598 0.2949 | 0.2743 0.2979 | 0.2825 0.3083 | 0.3195 | 0.3037 | Ave | | 0.2926 | | | 0.0100 | 6.6 | 20.0 | | | | |
| 1,2,3-Trichloropropane | 0.4104 0.3768 | 0.3859 0.3949 | 0.4160 0.3815 | 0.4181 | 0.4204 | Ave | | 0.4005 | | | 0.0100 | 4.4 | 20.0 | | | | |
| N-Propylbenzene | 1.0871 1.1604 | 1.1279 1.0214 | 1.1341 1.0987 | 1.1152 | 1.1268 | Ave | | 1.1089 | | | 0.0100 | 3.8 | 20.0 | | | | |
| 2-Chlorotoluene | 0.9007 0.9835 | 0.9855 0.9238 | 0.9604 0.9321 | 0.9790 | 1.0033 | Ave | | 0.9585 | | | 0.0100 | 3.7 | 20.0 | | | | |
| 3-Chlorotoluene | 1.0064 1.0049 | 1.0309 0.9798 | 1.0614 1.0388 | 1.1086 | 1.1105 | Ave | | 1.0427 | | | 0.0100 | 4.6 | 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.: 180-70873-1 Analy Batch No.: 218218

SDG No.: _____

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

Calibration Start Date: 07/27/2017 00:51 Calibration End Date: 07/27/2017 04:24 Calibration ID: 35038

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|------------------------------------|------------------|------------------|------------------|--------|--------|------------|-------------|--------|----|---|---------|------|------|----------|------------|---|----------------|
| | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| | LVL 6 | LVL 7 | LVL 8 | | | | | | | | | | | | | | |
| 1,3,5-Trimethylbenzene | 3.0303 3.1789 | 3.4364 2.8871 | 3.3130 2.9071 | 3.3121 | 3.3198 | Ave | | 3.1731 | | | 0.0100 | 6.6 | 20.0 | | | | |
| 4-Chlorotoluene | 1.0553 1.0614 | 1.0524 0.9741 | 1.0341 0.9970 | 1.0305 | 1.0761 | Ave | | 1.0351 | | | 0.0100 | 3.3 | 20.0 | | | | |
| tert-Butylbenzene | 2.5746 2.7227 | 2.8017 2.3880 | 2.7530 2.5138 | 2.7587 | 2.7116 | Ave | | 2.6530 | | | 0.0100 | 5.5 | 20.0 | | | | |
| 1,2,4-Trimethylbenzene | 3.1254 3.2212 | 3.4166 2.9826 | 3.3711 2.9395 | 3.3815 | 3.3664 | Ave | | 3.2255 | | | 0.0100 | 5.9 | 20.0 | | | | |
| 3,4-Dichlorobenzotrifluoride | 0.9400 0.7764 | 0.7679 0.7160 | 0.7941 0.8232 | 0.8410 | 0.8065 | Ave | | 0.8081 | | | 0.0100 | 8.1 | 20.0 | | | | |
| sec-Butylbenzene | 3.7533 3.7112 | 3.9865 3.2645 | 3.8932 3.4225 | 3.8001 | 3.7790 | Ave | | 3.7013 | | | 0.0100 | 6.5 | 20.0 | | | | |
| 1,3-Dichlorobenzene | 1.8909 1.6927 | 1.7949 1.6042 | 1.7488 1.5884 | 1.7678 | 1.7840 | Ave | | 1.7340 | | | 0.6000 | 5.8 | 20.0 | | | | |
| 4-Isopropyltoluene | 2.9547 3.1220 | 3.2883 2.7812 | 3.2665 2.8873 | 3.2019 | 3.1605 | Ave | | 3.0828 | | | 0.0100 | 6.0 | 20.0 | | | | |
| 1,4-Dichlorobenzene | 1.9782 1.7336 | 1.8319 1.6481 | 1.8074 1.6177 | 1.8136 | 1.8124 | Ave | | 1.7804 | | | 0.5000 | 6.4 | 20.0 | | | | |
| 2,4-Dichlorobenzotrifluoride | 0.7762 0.7410 | 0.7684 0.6560 | 0.7174 0.7931 | 0.7890 | 0.7781 | Ave | | 0.7524 | | | 0.0100 | 6.2 | 20.0 | | | | |
| 2,5-Dichlorobenzotrifluoride | 0.8709 0.7991 | 0.7991 0.7661 | 0.8033 0.8193 | 0.8304 | 0.8133 | Ave | | 0.8127 | | | 0.0100 | 3.7 | 20.0 | | | | |
| n-Butylbenzene | 2.4429 2.5807 | 2.6260 2.2815 | 2.6042 2.4382 | 2.5661 | 2.5760 | Ave | | 2.5144 | | | 0.0100 | 4.7 | 20.0 | | | | |
| 1,2-Dichlorobenzene | 1.8724 1.5966 | 1.7261 1.5319 | 1.6636 1.4748 | 1.6744 | 1.6818 | Ave | | 1.6527 | | | 0.4000 | 7.4 | 20.0 | | | | |
| 1,2-Dibromo-3-Chloropropane | 0.1676 0.1857 | 0.1676 0.2001 | 0.1774 0.1873 | 0.1829 | 0.1992 | Ave | | 0.1835 | | | 0.0500 | 6.8 | 20.0 | | | | |
| 2,4- & 2,5- & 2,6- Dichlorotoluene | 0.9836 1.0182 | 1.0277 0.9802 | 1.0819 1.0447 | 1.1339 | 1.1166 | Ave | | 1.0483 | | | 0.0100 | 5.5 | 20.0 | | | | |
| 2,3- & 3,4- Dichlorotoluene | 0.9469 1.0658 | 1.0253 1.0486 | 1.0886 1.1261 | 1.1868 | 1.1843 | Ave | | 1.0841 | | | 0.0100 | 7.5 | 20.0 | | | | |
| 1,2,4-Trichlorobenzene | 0.7563 0.7556 | 0.7184 0.7286 | 0.7717 0.7766 | 0.7671 | 0.7765 | Ave | | 0.7563 | | | 0.2000 | 2.9 | 20.0 | | | | |
| Hexachlorobutadiene | 0.2941 0.2697 | 0.2848 0.2377 | 0.2809 0.2898 | 0.2829 | 0.2739 | Ave | | 0.2767 | | | 0.0100 | 6.4 | 20.0 | | | | |
| Naphthalene | 2.0979 2.6004 | 2.2731 2.6494 | 2.6660 2.6327 | 2.8062 | 2.8819 | Ave | | 2.5759 | | | 0.0100 | 10.2 | 20.0 | | | | |
| 1,2,3-Trichlorobenzene | 0.7106 0.6701 | 0.6788 0.6564 | 0.6707 0.7130 | 0.7070 | 0.7206 | Ave | | 0.6909 | | | 0.0100 | 3.5 | 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.: 180-70873-1 Analy Batch No.: 218218
 SDG No.: _____
 Instrument ID: CHHP5 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N
 Calibration Start Date: 07/27/2017 00:51 Calibration End Date: 07/27/2017 04:24 Calibration ID: 35038

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|------------------------------|------------------|------------------|------------------|--------|--------|---------------|-------------|--------|----|---|---------|------|------|-------------|---------------|---|-------------------|
| | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| | LVL 6 | LVL 7 | LVL 8 | | | | | | | | | | | | | | |
| 2,4,5-Trichlorotoluene | 0.3224 0.3475 | 0.2818 0.3346 | 0.3064 ++++ | 0.3498 | 0.3564 | Ave | | 0.3284 | | | 0.0100 | 8.2 | 20.0 | | | | |
| 2,3,6-Trichlorotoluene | 0.2545 0.3128 | 0.2731 0.3131 | 0.3085 ++++ | 0.3418 | 0.3347 | Ave | | 0.3055 | | | 0.0100 | 10.3 | 20.0 | | | | |
| Dibromofluoromethane (Surr) | 0.2565 0.2365 | 0.2433 0.2326 | 0.2366 0.2242 | 0.2475 | 0.2474 | Ave | | 0.2406 | | | | 4.2 | 20.0 | | | | |
| 1,2-Dichloroethane-d4 (Surr) | 0.3401 0.2693 | 0.3050 0.2801 | 0.2948 0.2619 | 0.3004 | 0.2957 | Ave | | 0.2934 | | | | 8.3 | 20.0 | | | | |
| Toluene-d8 (Surr) | 5.1161 3.6702 | 4.5030 3.3148 | 4.0781 3.3147 | 3.9154 | 3.9228 | Ave | | 3.9794 | | | | 15.2 | 20.0 | | | | |
| 4-Bromofluorobenzene (Surr) | 1.6317 1.3781 | 1.5302 1.3139 | 1.4390 1.2793 | 1.4518 | 1.4735 | Ave | | 1.4372 | | | | 8.0 | 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
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1 Analy Batch No.: 218218

SDG No.: _____

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

Calibration Start Date: 07/27/2017 00:51 Calibration End Date: 07/27/2017 04:24 Calibration ID: 35038

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---------|-------------------|--------------|
| Level 1 | IC 180-218218/2 | 50727D02.D |
| Level 2 | IC 180-218218/3 | 50727D03.D |
| Level 3 | ICIS 180-218218/4 | 50727D04.D |
| Level 4 | IC 180-218218/5 | 50727D05.D |
| Level 5 | IC 180-218218/6 | 50727D06.D |
| Level 6 | IC 180-218218/10 | 50727D10.D |
| Level 7 | IC 180-218218/8 | 50727D08.D |
| Level 8 | IC 180-218218/11 | 50727D11.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 | 16788 647803 | 84559 569791 | 159957 857078 | 226899 | 286388 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Chloromethane | FB | Ave | 19706 595751 | 78965 580608 | 154943 811941 | 232300 | 302276 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Vinyl chloride | FB | Ave | 19568 632153 | 82670 577090 | 162634 867536 | 221295 | 291558 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,3-Butadiene | FB | Ave | 17968 579584 | 74553 512032 | 143576 815610 | 204212 | 260580 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Bromomethane | FB | Ave | 6901 285707 | 42224 289712 | 81346 377950 | 112119 | 161865 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Chloroethane | FB | Ave | 10685 340168 | 47273 322589 | 86601 414342 | 128899 | 172552 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Trichlorofluoromethane | FB | Ave | 22371 769762 | 104824 710415 | 205127 1017488 | 283194 | 371684 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Ethyl ether | FB | Ave | 14571 475422 | 66542 510033 | 126496 612640 | 188662 | 262150 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Acrolein | FB | Ave | 63695 154738 | 73476 179414 | 101829 183852 | 115103 | 130923 | 100 225 | 125 250 | 150 275 | 175 | 200 |
| 1,1-Dichloroethene | FB | Ave | 14263 540044 | 67928 489503 | 131576 745282 | 190985 | 247279 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | FB | Ave | 18126 571742 | 73846 534815 | 141127 774058 | 206212 | 263603 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Acetone | FB | Ave | 37823 447756 | 77890 522287 | 149782 630881 | 227784 | 316026 | 25.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| Iodomethane | FB | Ave | 22822 811997 | 103869 834240 | 200342 1099819 | 304618 | 408622 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Carbon disulfide | FB | Ave | 30868 +++++ | 131730 1211678 | 266935 1856339 | 403056 | 561008 | 5.00 +++++ | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Allyl chloride | FB | Ave | 8133 365237 | 39946 366340 | 83167 500032 | 121734 | 164305 | 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.: 180-70873-1

Analy Batch No.: 218218

SDG No.: _____

Instrument ID: CHHP5

GC Column: DB-624

ID: 0.18 (mm)

Heated Purge: (Y/N) N

Calibration Start Date: 07/27/2017 00:51

Calibration End Date: 07/27/2017 04:24

Calibration ID: 35038

| 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 | 31286 1009713 | 132543 1173609 | 283974 1447736 | 419273 | 558912 | 10.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| Methylene Chloride | FB | Lin2 | 25720 602402 | 84822 653341 | 164284 813282 | 242665 | 323324 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| tert-Butyl alcohol | TBAd 9 | Ave | 16447 524619 | 64738 519054 | 139891 568135 | 204334 | 283777 | 50.0 1750 | 250 2000 | 500 2500 | 750 | 1000 |
| Acrylonitrile | FB | Ave | 73302 2362587 | 336508 2794353 | 708552 3495451 | 1029651 | 1387354 | 50.0 1750 | 250 2000 | 500 2500 | 750 | 1000 |
| trans-1,2-Dichloroethene | FB | Ave | 17158 595572 | 73445 571864 | 147191 806194 | 222245 | 296608 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Methyl tert-butyl ether | FB | Ave | 38357 1597553 | 196780 1751345 | 390184 2170401 | 613933 | 822838 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Hexane | FB | Ave | 24902 760411 | 96542 708650 | 186124 1101558 | 266987 | 337300 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,1-Dichloroethane | FB | Ave | 28319 1024340 | 133976 1041269 | 261874 1376176 | 379320 | 510811 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Vinyl acetate | FB | Ave | 27185 1068205 | 115000 1200052 | 245879 1523056 | 400099 | 532250 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2,2-Dichloropropane | FB | Ave | 3769 136605 | 15889 125406 | 31118 188250 | 48893 | 65750 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| cis-1,2-Dichloroethene | FB | Ave | 17858 671208 | 85931 687049 | 172690 900432 | 259385 | 347303 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2-Butanone (MEK) | FB | Ave | 50216 686266 | 105960 795793 | 214731 962704 | 321867 | 426755 | 25.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| Bromochloromethane | FB | Ave | 8216 291754 | 38047 313977 | 75687 394763 | 113290 | 155416 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Tetrahydrofuran | FB | Ave | 14858 396477 | 52866 488432 | 117485 609910 | 176266 | 224432 | 10.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| Chloroform | FB | Ave | 29608 989929 | 134431 1037446 | 254354 1319564 | 389323 | 517765 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,1,1-Trichloroethane | FB | Ave | 20508 811476 | 98927 777880 | 196286 1097196 | 285488 | 383868 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Cyclohexane | FB | Ave | 26974 1012965 | 124196 922281 | 239333 1394833 | 345041 | 446560 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Carbon tetrachloride | FB | Ave | 17231 682784 | 80446 646700 | 162849 923177 | 238173 | 317033 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,1-Dichloropropene | FB | Ave | 22014 866715 | 109851 825970 | 215336 1178056 | 312373 | 408627 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Isobutyl alcohol | FB | Ave | 13122 452876 | 61305 587752 | 136973 715201 | 216532 | 290317 | 125 4375 | 625 5000 | 1250 6250 | 1875 | 2500 |
| Benzene | FB | Ave | 74686 2459963 | 339765 2487856 | 669098 3249284 | 981851 | 1307056 | 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.: 180-70873-1 Analy Batch No.: 218218

SDG No.: _____

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

Calibration Start Date: 07/27/2017 00:51 Calibration End Date: 07/27/2017 04:24 Calibration ID: 35038

| 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 | 21038 708898 | 95627 767974 | 190422 969148 | 292683 | 385206 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| n-Heptane | FB | Ave | 16453 633483 | 81002 573064 | 154370 922592 | 214813 | 279216 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Trichloroethene | FB | Ave | 17490 648262 | 83072 647404 | 164695 887332 | 241861 | 329499 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Methylcyclohexane | FB | Ave | 25605 1041060 | 125697 950167 | 253511 1432791 | 358781 | 467268 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2-Dichloropropane | FB | Ave | 16316 596512 | 74777 624637 | 150135 793667 | 227133 | 309491 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,4-Dioxane | FB | Ave | 2333 115916 | 15162 135844 | 33209 187034 | 46920 | 65688 | 100 3500 | 500 4000 | 1000 5000 | 1500 | 2000 |
| Dibromomethane | FB | Ave | 8641 342853 | 45949 374289 | 88395 470836 | 135198 | 184529 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Bromodichloromethane | FB | Ave | 16257 712434 | 84070 752352 | 171049 945026 | 268080 | 366097 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2-Chloroethyl vinyl ether | FB | Ave | 18086 864836 | 103158 977190 | 219328 1234429 | 343066 | 467677 | 10.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| cis-1,3-Dichloropropene | FB | Ave | 19479 881560 | 96744 933591 | 204344 1203144 | 320956 | 447138 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 4-Methyl-2-pentanone (MIBK) | CBNZ d5 | Ave | 79892 1265241 | 154465 1476808 | 361112 1863520 | 542662 | 738839 | 25.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| Toluene | CBNZ d5 | Ave | 71883 2496911 | 351840 2540251 | 692901 3254284 | 1000479 | 1332783 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| trans-1,3-Dichloropropene | CBNZ d5 | Ave | 14443 781619 | 79122 850338 | 170710 1070347 | 278226 | 396221 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Ethyl methacrylate | CBNZ d5 | Ave | 16030 905216 | 96602 1001550 | 222171 1271580 | 352819 | 483364 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,1,2-Trichloroethane | CBNZ d5 | Ave | 14755 523017 | 67966 569083 | 138196 718069 | 209928 | 283688 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Tetrachloroethene | CBNZ d5 | Ave | 13528 498519 | 67579 486427 | 126273 683462 | 184171 | 244346 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,3-Dichloropropane | CBNZ d5 | Ave | 26359 969241 | 127957 1058308 | 256477 1320887 | 397870 | 518120 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2-Hexanone | CBNZ d5 | Ave | 57842 977068 | 122936 1109580 | 278579 1418811 | 419354 | 581383 | 25.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| Dibromochloromethane | CBNZ d5 | Ave | 9414 489506 | 53302 540065 | 114911 672369 | 181267 | 254603 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2-Dibromoethane (EDB) | CBNZ d5 | Ave | 13462 550826 | 67745 607203 | 142489 773664 | 223815 | 294438 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 3-Chlorobenzotrifluoride | CBNZ d5 | Ave | 25343 874266 | 109109 869071 | 222871 1290067 | 352260 | 461082 | 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.: 180-70873-1

Analy Batch No.: 218218

SDG No.: _____

Instrument ID: CHHP5

GC Column: DB-624

ID: 0.18 (mm)

Heated Purge: (Y/N) N

Calibration Start Date: 07/27/2017 00:51

Calibration End Date: 07/27/2017 04:24

Calibration ID: 35038

| 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 | 47566 1645967 | 217561 1704167 | 431311 2170926 | 660247 | 877804 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 4-Chlorobenzotrifluoride | CBNZ d5 | Ave | 21933 826850 | 101825 810848 | 207774 1226371 | 327327 | 420704 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,1,1,2-Tetrachloroethane | CBNZ d5 | Ave | 12587 554351 | 65901 590452 | 137710 751692 | 212641 | 289044 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Ethylbenzene | CBNZ d5 | Ave | 22622 962208 | 120759 972676 | 249792 1304914 | 371119 | 499116 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| m-Xylene & p-Xylene | CBNZ d5 | Ave | 25553 1197380 | 151114 1217768 | 306948 1614353 | 452043 | 610286 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| o-Xylene | CBNZ d5 | Ave | 25240 1130677 | 138375 1159372 | 288885 1518391 | 440285 | 592117 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Styrene | CBNZ d5 | Ave | 42810 1866053 | 242031 1967591 | 498873 2462559 | 745860 | 1002147 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Bromoform | CBNZ d5 | Ave | 6015 310948 | 30000 350923 | 67829 443094 | 112077 | 157509 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2-Chlorobenzotrifluoride | CBNZ d5 | Ave | 21074 840920 | 107103 875687 | 216286 1244752 | 348911 | 454842 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Isopropylbenzene | CBNZ d5 | Ave | 64937 2681266 | 356966 2665903 | 726432 3502176 | 1080505 | 1415676 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Bromobenzene | DCBd 4 | Ave | 16032 659984 | 83376 711710 | 163748 889999 | 261052 | 348475 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,1,2,2-Tetrachloroethane | CBNZ d5 | Ave | 20749 762601 | 100341 870164 | 211912 1078742 | 316221 | 412534 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| trans-1,4-Dichloro-2-butene | DCBd 4 | Ave | 4170 199800 | 23168 225821 | 49334 299994 | 83561 | 104361 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2,3-Trichloropropane | DCBd 4 | Ave | 6588 255265 | 32588 299299 | 72643 371250 | 109372 | 144469 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| N-Propylbenzene | DCBd 4 | Ave | 17451 786064 | 95261 774184 | 198029 1069171 | 291693 | 387234 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2-Chlorotoluene | DCBd 4 | Ave | 14458 666236 | 83234 700158 | 167713 907016 | 256066 | 344800 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 3-Chlorotoluene | DCBd 4 | Ave | 16155 680717 | 87067 742625 | 185343 1010916 | 289960 | 381649 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,3,5-Trimethylbenzene | DCBd 4 | Ave | 48645 2153457 | 290219 2188229 | 578518 2828999 | 866332 | 1140888 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 4-Chlorotoluene | DCBd 4 | Ave | 16940 719035 | 88877 738280 | 180584 970169 | 269544 | 369832 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| tert-Butylbenzene | DCBd 4 | Ave | 41329 1844417 | 236619 1809964 | 480729 2446270 | 721573 | 931884 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2,4-Trimethylbenzene | DCBd 4 | Ave | 50171 2182090 | 288545 2260604 | 588662 2860516 | 884487 | 1156912 | 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.: 180-70873-1 Analy Batch No.: 218218

SDG No.: _____

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

Calibration Start Date: 07/27/2017 00:51 Calibration End Date: 07/27/2017 04:24 Calibration ID: 35038

| 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 | 15090 525922 | 64854 542681 | 138659 801099 | 219982 | 277157 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| sec-Butylbenzene | DCBd 4 | Ave | 60251 2514051 | 336681 2474312 | 679839 3330508 | 993968 | 1298722 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,3-Dichlorobenzene | DCBd 4 | Ave | 30355 1146674 | 151590 1215884 | 305374 1545747 | 462404 | 613101 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 4-Isopropyltoluene | DCBd 4 | Ave | 47431 2114911 | 277710 2107989 | 570403 2809716 | 837492 | 1086140 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,4-Dichlorobenzene | DCBd 4 | Ave | 31756 1174377 | 154714 1249173 | 315614 1574222 | 474362 | 622850 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2,4-Dichlorobenzotrifluoride | DCBd 4 | Ave | 12460 501975 | 64892 497225 | 125268 771761 | 206368 | 267418 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2,5-Dichlorobenzotrifluoride | DCBd 4 | Ave | 13980 541324 | 67486 580659 | 140272 797256 | 217211 | 279514 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| n-Butylbenzene | DCBd 4 | Ave | 39215 1748217 | 221777 1729209 | 454742 2372703 | 671190 | 885288 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2-Dichlorobenzene | DCBd 4 | Ave | 30057 1081541 | 145778 1161072 | 290492 1435184 | 437966 | 577962 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2-Dibromo-3-Chloropropane | DCBd 4 | Ave | 2690 125814 | 14158 151695 | 30986 182290 | 47827 | 68470 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2,4- & 2,5- & 2,6- Dichlorotoluene | DCBd 4 | Ave | 47367 2069215 | 260387 2228710 | 566788 3049908 | 889724 | 1151252 | 15.0 525 | 75.0 600 | 150 750 | 225 | 300 |
| 2,3- & 3,4- Dichlorotoluene | DCBd 4 | Ave | 30402 1443949 | 173187 1589536 | 380181 2191624 | 620870 | 814032 | 10.0 350 | 50.0 400 | 100 500 | 150 | 200 |
| 1,2,4-Trichlorobenzene | DCBd 4 | Ave | 12140 511830 | 60672 552245 | 134753 755690 | 200638 | 266863 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Hexachlorobutadiene | DCBd 4 | Ave | 4721 182711 | 24054 180140 | 49048 282046 | 73984 | 94134 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Naphthalene | DCBd 4 | Ave | 33677 1761559 | 191971 2008065 | 465533 2561966 | 733996 | 990398 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2,3-Trichlorobenzene | DCBd 4 | Ave | 11407 453926 | 57325 497473 | 117120 693791 | 184932 | 247660 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 2,4,5-Trichlorotoluene | DCBd 4 | Ave | 5175 235417 | 23799 253594 | 53498 ++++ | 91488 | 122498 | 5.00 175 | 25.0 200 | 50.0 ++++ | 75.0 | 100 |
| 2,3,6-Trichlorotoluene | DCBd 4 | Ave | 4086 211883 | 23065 237299 | 53869 ++++ | 89402 | 115009 | 5.00 175 | 25.0 200 | 50.0 ++++ | 75.0 | 100 |
| Dibromofluoromethane (Surr) | FB | Ave | 13893 505019 | 65453 522323 | 127700 681339 | 193042 | 257355 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| 1,2-Dichloroethane-d4 (Surr) | FB | Ave | 18421 575099 | 82071 628942 | 159071 795993 | 234269 | 307676 | 5.00 175 | 25.0 200 | 50.0 250 | 75.0 | 100 |
| Toluene-d8 (Surr) | CBNZ d5 | Ave | 60283 1992609 | 278432 2000995 | 541748 2678162 | 780569 | 1040595 | 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.: 180-70873-1 Analy Batch No.: 218218

SDG No.: _____

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

Calibration Start Date: 07/27/2017 00:51 Calibration End Date: 07/27/2017 04:24 Calibration ID: 35038

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG) | | | | |
|-----------------------------|------------|------------|-----------------|-----------------|-------------------|--------|--------|--------------------|-------------|-------------|-------|-------|
| | | | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| | | | LVL 6 | LVL 7 | LVL 8 | | | LVL 6 | LVL 7 | LVL 8 | | |
| 4-Bromofluorobenzene (Surr) | CBNZ d5 | Ave | 19227 748217 | 94618 793129 | 191158 1033645 | 289432 | 390879 | 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.: 180-70873-1 Analy Batch No.: 218218

SDG No.: _____

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

Calibration Start Date: 07/27/2017 00:51 Calibration End Date: 07/27/2017 04:24 Calibration ID: 35038

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---------|-------------------|--------------|
| Level 1 | IC 180-218218/2 | 50727D02.D |
| Level 2 | IC 180-218218/3 | 50727D03.D |
| Level 3 | ICIS 180-218218/4 | 50727D04.D |
| Level 4 | IC 180-218218/5 | 50727D05.D |
| Level 5 | IC 180-218218/6 | 50727D06.D |
| Level 6 | IC 180-218218/10 | 50727D10.D |
| Level 7 | IC 180-218218/8 | 50727D08.D |
| Level 8 | IC 180-218218/11 | 50727D11.D |

| 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 |
| | LVL 7 # | LVL 8 # | | | | | LVL 7 | LVL 8 | | | | |
| Dichlorodifluoromethane | 6.6 -12.7 | 8.1 -3.0 | 1.9 | 0.1 | -5.3 | 4.3 | 50 30 | 30 | 30 | 30 | 30 | 30 |
| Chloromethane | 24.5 -11.5 | 0.4 -8.6 | -1.7 | 1.9 | -0.6 | -4.5 | 50 30 | 30 | 30 | 30 | 30 | 30 |
| Vinyl chloride | 21.8 -13.3 | 3.6 -3.7 | 1.6 | -4.3 | -5.5 | -0.2 | 50 30 | 30 | 30 | 30 | 30 | 30 |
| 1,3-Butadiene | 23.1 -15.3 | 2.9 -0.4 | -1.2 | -2.8 | -7.0 | 0.8 | 50 30 | 30 | 30 | 30 | 30 | 30 |
| Bromomethane | -9.1 -8.0 | 11.9 -11.3 | 7.5 | 2.5 | 11.0 | -4.6 | 50 30 | 30 | 30 | 30 | 30 | 30 |
| Chloroethane | 21.0 -11.8 | 7.8 -16.3 | -1.5 | 1.4 | 1.8 | -2.3 | 50 30 | 30 | 30 | 30 | 30 | 30 |
| Trichlorofluoromethane | 13.3 -13.2 | 6.9 -8.1 | 4.3 | -0.3 | -1.9 | -1.1 | 50 30 | 30 | 30 | 30 | 30 | 30 |
| Ethyl ether | 13.5 -4.1 | 4.4 -14.9 | -1.1 | 2.1 | 6.3 | -6.1 | 50 30 | 30 | 30 | 30 | 30 | 30 |
| Acrolein | -1.6 7.0 | -8.5 -7.9 | 5.3 | 5.9 | 5.4 | -5.6 | 50 30 | 30 | 30 | 30 | 30 | 30 |
| 1,1-Dichloroethene | 7.6 -10.9 | 3.1 0.2 | -0.4 | 0.0 | -2.9 | 3.3 | 50 30 | 30 | 30 | 30 | 30 | 30 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 24.6 -11.3 | 2.2 -5.2 | -2.7 | -1.6 | -5.7 | -0.3 | 50 30 | 30 | 30 | 30 | 30 | 30 |
| Acetone | 6.8 -11.0 | 10.7 -20.6 | 6.1 | 11.7 | 16.2 | -19.8 | 50 30 | 30 | 30 | 30 | 30 | 30 |
| Iodomethane | 9.6 -3.4 | 0.4 -5.9 | -3.4 | 1.6 | 2.2 | -1.1 | 50 30 | 30 | 30 | 30 | 30 | 30 |
| Carbon disulfide | 6.1 0.5 | -8.9 13.7 | -7.9 | -3.8 | 0.4 | +++++ | 50 30 | 30 | 30 | 30 | 30 | 30 |
| Allyl chloride | -5.1 3.1 | -6.1 4.0 | -2.6 | -1.3 | -0.2 | 8.1 | 50 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.: 180-70873-1 Analy Batch No.: 218218

SDG No.: _____

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

Calibration Start Date: 07/27/2017 00:51 Calibration End Date: 07/27/2017 04:24 Calibration ID: 35038

| 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 |
| | LVL 7 # | LVL 8 # | | | | | LVL 7 | LVL 8 | | | | |
| Methyl acetate | 11.5 | -4.9 | 1.6 | 3.8 | 3.7 | -8.7 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 0.9 | -8.0 | | | | | 30 | 30 | | | | |
| Methylene Chloride | 0.0 | -2.5 | 0.4 | 5.0 | 6.0 | -2.6 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 0.7 | -7.2 | | | | | 30 | 30 | | | | |
| tert-Butyl alcohol | 12.8 | -2.2 | -1.6 | -4.3 | -4.9 | 8.8 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -13.1 | 4.4 | | | | | 30 | 30 | | | | |
| Acrylonitrile | 7.5 | -0.7 | 4.3 | 4.9 | 5.9 | -12.1 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -1.1 | -8.6 | | | | | 30 | 30 | | | | |
| trans-1,2-Dichloroethene | 13.6 | -2.1 | -2.2 | 2.2 | 2.2 | 0.0 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -8.7 | -4.9 | | | | | 30 | 30 | | | | |
| Methyl tert-butyl ether | -5.3 | -2.2 | -3.3 | 5.3 | 5.8 | 0.0 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 4.3 | -4.5 | | | | | 30 | 30 | | | | |
| Hexane | 28.4 | 0.2 | -3.7 | -4.4 | -9.4 | -0.5 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -11.8 | 1.2 | | | | | 30 | 30 | | | | |
| 1,1-Dichloroethane | 7.8 | 2.7 | 0.1 | 0.3 | 1.2 | -1.1 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -4.4 | -6.6 | | | | | 30 | 30 | | | | |
| Vinyl acetate | 1.8 | -13.3 | -7.6 | 4.0 | 3.7 | 1.4 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 8.4 | 1.6 | | | | | 30 | 30 | | | | |
| 2,2-Dichloropropane | 12.7 | -4.4 | -6.6 | 1.5 | 2.4 | 3.6 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -9.5 | 0.3 | | | | | 30 | 30 | | | | |
| cis-1,2-Dichloroethene | 3.3 | 0.1 | 0.3 | 4.3 | 4.6 | -1.5 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -4.1 | -7.1 | | | | | 30 | 30 | | | | |
| 2-Butanone (MEK) | -0.4 | 5.8 | 6.9 | 10.9 | 10.2 | -13.7 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -4.8 | -14.9 | | | | | 30 | 30 | | | | |
| Bromochloromethane | 7.0 | -0.3 | -1.1 | 2.5 | 5.4 | -3.6 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -1.4 | -8.4 | | | | | 30 | 30 | | | | |
| Tetrahydrofuran | 26.5 | -9.4 | 0.4 | 4.3 | -0.5 | -14.3 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 0.4 | -7.4 | | | | | 30 | 30 | | | | |
| Chloroform | 12.9 | 3.2 | -2.7 | 3.1 | 2.8 | -4.3 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -4.6 | -10.3 | | | | | 30 | 30 | | | | |
| 1,1,1-Trichloroethane | 3.3 | 0.3 | -0.8 | -0.1 | 0.7 | 3.7 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -5.5 | -1.5 | | | | | 30 | 30 | | | | |
| Cyclohexane | 10.1 | 2.0 | -2.0 | -2.2 | -5.1 | 4.9 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -9.2 | 1.5 | | | | | 30 | 30 | | | | |
| Carbon tetrachloride | 4.3 | -2.0 | -1.1 | 0.1 | -0.1 | 4.8 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -5.6 | -0.4 | | | | | 30 | 30 | | | | |
| 1,1-Dichloropropene | 2.6 | 3.1 | 0.7 | 1.1 | -0.8 | 2.5 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -7.1 | -2.1 | | | | | 30 | 30 | | | | |
| Isobutyl alcohol | -2.6 | -8.4 | 2.0 | 11.6 | 12.2 | -14.7 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 5.2 | -5.4 | | | | | 30 | 30 | | | | |
| Benzene | 13.4 | 3.9 | 2.0 | 3.6 | 3.3 | -5.2 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -8.9 | -12.1 | | | | | 30 | 30 | | | | |

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

Lab Name: TestAmerica Pittsburgh

Job No.: 180-70873-1

Analy Batch No.: 218218

SDG No.: _____

Instrument ID: CHHP5

GC Column: DB-624

ID: 0.18 (mm)

Heated Purge: (Y/N) N

Calibration Start Date: 07/27/2017 00:51

Calibration End Date: 07/27/2017 04:24

Calibration ID: 35038

| 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 |
| | LVL 7 # | LVL 8 # | | | | | LVL 7 | LVL 8 | | | | |
| 1,2-Dichloroethane | 9.6 | 0.3 | -0.4 | 5.9 | 4.5 | -6.3 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -3.5 | -10.0 | | | | | 30 | 30 | | | | |
| n-Heptane | 6.1 | 5.2 | -0.1 | -3.8 | -6.2 | 3.6 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -10.8 | 6.0 | | | | | 30 | 30 | | | | |
| Trichloroethene | 5.5 | 0.9 | -0.3 | 1.4 | 3.5 | -0.8 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -5.8 | -4.6 | | | | | 30 | 30 | | | | |
| Methylcyclohexane | 2.2 | 1.0 | 1.5 | -0.6 | -2.9 | 5.4 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -8.5 | 1.9 | | | | | 30 | 30 | | | | |
| 1,2-Dichloropropane | 6.4 | -1.8 | -1.7 | 2.9 | 5.1 | -1.3 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -1.7 | -7.7 | | | | | 30 | 30 | | | | |
| 1,4-Dioxane | -25.2 | -2.1 | 6.9 | 4.5 | 9.7 | -5.7 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 5.1 | 6.9 | | | | | 30 | 30 | | | | |
| Dibromomethane | -3.8 | 3.0 | -1.3 | 4.5 | 6.9 | -3.2 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 0.5 | -6.6 | | | | | 30 | 30 | | | | |
| Bromodichloromethane | -7.8 | -4.0 | -2.7 | 5.6 | 8.1 | 2.5 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 2.9 | -4.5 | | | | | 30 | 30 | | | | |
| 2-Chloroethyl vinyl ether | -18.1 | -5.9 | -0.3 | 8.0 | 10.3 | -0.6 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 6.8 | -0.3 | | | | | 30 | 30 | | | | |
| cis-1,3-Dichloropropene | -9.1 | -9.1 | -4.3 | 4.1 | 8.7 | 4.4 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 5.1 | 0.1 | | | | | 30 | 30 | | | | |
| 4-Methyl-2-pentanone (MIBK) | 5.7 | -2.6 | 6.0 | 6.1 | 8.6 | -9.1 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -4.6 | -10.1 | | | | | 30 | 30 | | | | |
| Toluene | 22.4 | 14.1 | 4.6 | 0.7 | 0.8 | -7.8 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -15.6 | -19.2 | | | | | 30 | 30 | | | | |
| trans-1,3-Dichloropropene | -9.6 | -5.7 | -5.3 | 2.9 | 10.1 | 6.1 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 3.8 | -2.3 | | | | | 30 | 30 | | | | |
| Ethyl methacrylate | -16.8 | -4.5 | 2.2 | 8.2 | 11.4 | 1.9 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 1.4 | -3.8 | | | | | 30 | 30 | | | | |
| 1,1,2-Trichloroethane | 20.6 | 5.8 | 0.2 | 1.4 | 3.0 | -7.2 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -9.2 | -14.4 | | | | | 30 | 30 | | | | |
| Tetrachloroethene | 20.7 | 14.9 | 0.0 | -2.8 | -3.1 | -3.4 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -15.3 | -11.0 | | | | | 30 | 30 | | | | |
| 1,3-Dichloropropane | 16.5 | 7.8 | 0.6 | 4.0 | 1.7 | -7.0 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -8.7 | -14.8 | | | | | 30 | 30 | | | | |
| 2-Hexanone | -0.2 | 1.1 | 6.6 | 6.9 | 11.4 | -8.5 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -6.6 | -10.7 | | | | | 30 | 30 | | | | |
| Dibromochloromethane | -9.0 | -1.8 | -1.5 | 3.6 | 9.3 | 2.7 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 1.9 | -5.2 | | | | | 30 | 30 | | | | |
| 1,2-Dibromoethane (EDB) | 7.3 | 2.9 | 0.7 | 5.4 | 4.2 | -4.8 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -5.6 | -10.1 | | | | | 30 | 30 | | | | |
| 3-Chlorobenzotrifluoride | 25.2 | 2.7 | -2.4 | 2.8 | 1.2 | -6.3 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -16.2 | -7.1 | | | | | 30 | 30 | | | | |

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

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1 Analy Batch No.: 218218

SDG No.: _____

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

Calibration Start Date: 07/27/2017 00:51 Calibration End Date: 07/27/2017 04:24 Calibration ID: 35038

| 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 |
| | LVL 7 # | LVL 8 # | | | | | LVL 7 | LVL 8 | | | | |
| Chlorobenzene | 24.4 | 8.4 | 0.0 | 2.0 | 2.0 | -6.6 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -13.0 | -17.2 | | | | | 30 | 30 | | | | |
| 4-Chlorobenzotrifluoride | 17.4 | 3.9 | -1.4 | 3.6 | 0.0 | -3.9 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -15.3 | -4.3 | | | | | 30 | 30 | | | | |
| 1,1,1,2-Tetrachloroethane | 3.5 | 3.3 | 0.4 | 3.4 | 5.6 | -1.1 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -5.2 | -9.9 | | | | | 30 | 30 | | | | |
| Ethylbenzene | 6.0 | 7.8 | 3.8 | 2.7 | 3.8 | -2.2 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -11.1 | -10.9 | | | | | 30 | 30 | | | | |
| m-Xylene & p-Xylene | -2.1 | 10.4 | 4.4 | 2.4 | 3.9 | -0.4 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -8.9 | -9.8 | | | | | 30 | 30 | | | | |
| o-Xylene | 1.5 | 6.1 | 3.1 | 4.7 | 5.8 | -1.3 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -9.0 | -10.9 | | | | | 30 | 30 | | | | |
| Styrene | 1.7 | 9.6 | 5.2 | 4.8 | 5.8 | -3.7 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -8.7 | -14.6 | | | | | 30 | 30 | | | | |
| Bromoform | -6.4 | -11.1 | -6.4 | 3.0 | 8.8 | 5.0 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 6.6 | 0.5 | | | | | 30 | 30 | | | | |
| 2-Chlorobenzotrifluoride | 8.8 | 5.3 | -1.0 | 6.4 | 4.3 | -5.8 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -11.8 | -6.3 | | | | | 30 | 30 | | | | |
| Isopropylbenzene | 7.0 | 12.1 | 6.2 | 5.2 | 3.6 | -4.1 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -14.2 | -15.8 | | | | | 30 | 30 | | | | |
| Bromobenzene | 2.9 | 1.7 | -3.4 | 2.8 | 4.5 | 0.4 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -3.2 | -5.8 | | | | | 30 | 30 | | | | |
| 1,1,2,2-Tetrachloroethane | 14.5 | 5.5 | 3.7 | 3.2 | 1.1 | -8.7 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -6.3 | -13.2 | | | | | 30 | 30 | | | | |
| trans-1,4-Dichloro-2-butene | -11.2 | -6.3 | -3.4 | 9.2 | 3.8 | 0.8 | 50 | 30 | 30 | 30 | 30 | 30 |
| | 1.8 | 5.4 | | | | | 30 | 30 | | | | |
| 1,2,3-Trichloropropane | 2.5 | -3.7 | 3.9 | 4.4 | 5.0 | -5.9 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -1.4 | -4.7 | | | | | 30 | 30 | | | | |
| N-Propylbenzene | -2.0 | 1.7 | 2.3 | 0.6 | 1.6 | 4.6 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -7.9 | -0.9 | | | | | 30 | 30 | | | | |
| 2-Chlorotoluene | -6.0 | 2.8 | 0.2 | 2.1 | 4.7 | 2.6 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -3.6 | -2.8 | | | | | 30 | 30 | | | | |
| 3-Chlorotoluene | -3.5 | -1.1 | 1.8 | 6.3 | 6.5 | -3.6 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -6.0 | -0.4 | | | | | 30 | 30 | | | | |
| 1,3,5-Trimethylbenzene | -4.5 | 8.3 | 4.4 | 4.4 | 4.6 | 0.2 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -9.0 | -8.4 | | | | | 30 | 30 | | | | |
| 4-Chlorotoluene | 1.9 | 1.7 | -0.1 | -0.4 | 4.0 | 2.5 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -5.9 | -3.7 | | | | | 30 | 30 | | | | |
| tert-Butylbenzene | -3.0 | 5.6 | 3.8 | 4.0 | 2.2 | 2.6 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -10.0 | -5.2 | | | | | 30 | 30 | | | | |
| 1,2,4-Trimethylbenzene | -3.1 | 5.9 | 4.5 | 4.8 | 4.4 | -0.1 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -7.5 | -8.9 | | | | | 30 | 30 | | | | |

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

Lab Name: TestAmerica Pittsburgh

Job No.: 180-70873-1

Analy Batch No.: 218218

SDG No.: _____

Instrument ID: CHHP5

GC Column: DB-624

ID: 0.18 (mm)

Heated Purge: (Y/N) N

Calibration Start Date: 07/27/2017 00:51

Calibration End Date: 07/27/2017 04:24

Calibration ID: 35038

| 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 |
| | LVL 7 # | LVL 8 # | | | | | LVL 7 | LVL 8 | | | | |
| 3,4-Dichlorobenzotrifluoride | 16.3 -11.4 | -5.0 1.9 | -1.7 | 4.1 | -0.2 | -3.9 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| sec-Butylbenzene | 1.4 -11.8 | 7.7 -7.5 | 5.2 | 2.7 | 2.1 | 0.3 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,3-Dichlorobenzene | 9.1 -7.5 | 3.5 -8.4 | 0.9 | 2.0 | 2.9 | -2.4 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 4-Isopropyltoluene | -4.2 -9.8 | 6.7 -6.3 | 6.0 | 3.9 | 2.5 | 1.3 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,4-Dichlorobenzene | 11.1 -7.4 | 2.9 -9.1 | 1.5 | 1.9 | 1.8 | -2.6 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2,4-Dichlorobenzotrifluoride | 3.2 -12.8 | 2.1 5.4 | -4.7 | 4.9 | 3.4 | -1.5 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2,5-Dichlorobenzotrifluoride | 7.2 -5.7 | -1.7 0.8 | -1.2 | 2.2 | 0.1 | -1.7 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| n-Butylbenzene | -2.8 -9.3 | 4.4 -3.0 | 3.6 | 2.1 | 2.4 | 2.6 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,2-Dichlorobenzene | 13.3 -7.3 | 4.4 -10.8 | 0.7 | 1.3 | 1.8 | -3.4 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,2-Dibromo-3-Chloropropane | -8.7 9.1 | -8.6 2.1 | -3.3 | -0.3 | 8.6 | 1.2 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2,4- & 2,5- & 2,6- Dichlorotoluene | -6.2 -6.5 | -2.0 -0.3 | 3.2 | 8.2 | 6.5 | -2.9 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2,3- & 3,4- Dichlorotoluene | -12.6 -3.3 | -5.4 3.9 | 0.4 | 9.5 | 9.3 | -1.7 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,2,4-Trichlorobenzene | 0.0 -3.7 | -5.0 2.7 | 2.0 | 1.4 | 2.7 | -0.1 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Hexachlorobutadiene | 6.3 -14.1 | 2.9 4.7 | 1.5 | 2.2 | -1.0 | -2.5 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Naphthalene | -18.6 2.9 | -11.8 2.2 | 3.5 | 8.9 | 11.9 | 1.0 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,2,3-Trichlorobenzene | 2.9 -5.0 | -1.8 3.2 | -2.9 | 2.3 | 4.3 | -3.0 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2,4,5-Trichlorotoluene | -1.8 1.9 | -14.2 ++++ | -6.7 | 6.5 | 8.5 | 5.8 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 2,3,6-Trichlorotoluene | -16.7 2.5 | -10.6 ++++ | 1.0 | 11.9 | 9.5 | 2.4 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Dibromofluoromethane (Surr) | 6.6 -3.3 | 1.1 -6.8 | -1.6 | 2.9 | 2.8 | -1.7 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| 1,2-Dichloroethane-d4 (Surr) | 15.9 -4.5 | 4.0 -10.7 | 0.5 | 2.4 | 0.8 | -8.2 | 50 30 | 30 30 | 30 | 30 | 30 | 30 |
| Toluene-d8 (Surr) | 28.6 -16.7 | 13.2 -16.7 | 2.5 | -1.6 | -1.4 | -7.8 | 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.: 180-70873-1 Analy Batch No.: 218218

SDG No.: _____

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

Calibration Start Date: 07/27/2017 00:51 Calibration End Date: 07/27/2017 04:24 Calibration ID: 35038

| 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 |
| | LVL 7 # | LVL 8 # | | | | | LVL 7 | LVL 8 | | | | |
| 4-Bromofluorobenzene (Surr) | 13.5 | 6.5 | 0.1 | 1.0 | 2.5 | -4.1 | 50 | 30 | 30 | 30 | 30 | 30 |
| | -8.6 | -11.0 | | | | | 30 | 30 | | | | |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D02.D
 Lims ID: IC VSTD1
 Client ID:
 Sample Type: IC Calib Level: 1
 Inject. Date: 27-Jul-2017 00:51:30 ALS Bottle#: 2 Worklist Smp#: 2
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017756-002
 Misc. Info.: IC VSTD1
 Operator ID: 034635 Instrument ID: CHHP5
 Sublist: chrom-MSVOA_LL_CHHP5*sub12
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 28-Jul-2017 01:04:45 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK029

First Level Reviewer: bungardf

Date: 27-Jul-2017 03:08:26

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.317 | 4.323 | -0.006 | 0 | 246479 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.298 | 7.298 | 0.000 | 99 | 541701 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.406 | 10.406 | 0.000 | 85 | 117831 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.773 | 12.773 | 0.000 | 96 | 160528 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.574 | 6.574 | 0.000 | 90 | 13893 | 5.00 | 5.33 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.951 | 6.945 | 0.006 | 0 | 18421 | 5.00 | 5.79 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.946 | 8.946 | 0.000 | 92 | 60283 | 5.00 | 6.43 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.599 | 11.599 | 0.000 | 87 | 19227 | 5.00 | 5.68 | |
| 11 Dichlorodifluoromethane | 85 | 1.665 | 1.646 | 0.018 | 68 | 16788 | 5.00 | 5.33 | |
| 12 Chloromethane | 50 | 1.804 | 1.804 | 0.000 | 97 | 19706 | 5.00 | 6.22 | |
| 13 Vinyl chloride | 62 | 1.932 | 1.944 | -0.012 | 95 | 19568 | 5.00 | 6.09 | |
| 14 Butadiene | 39 | 1.963 | 1.969 | -0.005 | 95 | 17968 | 5.00 | 6.16 | |
| 15 Bromomethane | 94 | 2.273 | 2.254 | 0.019 | 90 | 6901 | 5.00 | 4.54 | |
| 16 Chloroethane | 64 | 2.419 | 2.419 | 0.000 | 89 | 10685 | 5.00 | 6.05 | |
| 17 Dichlorofluoromethane | 67 | 2.699 | 2.699 | 0.000 | 97 | 26531 | 5.00 | 5.94 | |
| 18 Trichlorofluoromethane | 101 | 2.760 | 2.741 | 0.019 | 45 | 22371 | 5.00 | 5.67 | M |
| 20 Ethyl ether | 59 | 3.076 | 3.076 | 0.000 | 88 | 14571 | 5.00 | 5.67 | |
| 21 Acrolein | 56 | 3.252 | 3.252 | 0.000 | 99 | 63695 | 100.0 | 98.4 | |
| 22 1,1-Dichloroethene | 96 | 3.368 | 3.368 | 0.000 | 77 | 14263 | 5.00 | 5.38 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.441 | 3.441 | 0.000 | 74 | 18126 | 5.00 | 6.23 | |
| 24 Acetone | 43 | 3.483 | 3.477 | 0.006 | 99 | 37823 | 25.0 | 26.7 | |
| 25 Iodomethane | 142 | 3.569 | 3.562 | 0.007 | 95 | 22822 | 5.00 | 5.48 | |
| 26 Carbon disulfide | 76 | 3.654 | 3.648 | 0.006 | 98 | 30868 | 5.00 | 5.30 | |
| 28 3-Chloro-1-propene | 76 | 3.940 | 3.946 | -0.006 | 90 | 8133 | 5.00 | 4.75 | |
| 30 Methyl acetate | 43 | 3.970 | 3.976 | -0.006 | 95 | 31286 | 10.0 | 11.2 | |
| 31 Methylene Chloride | 84 | 4.177 | 4.165 | 0.012 | 84 | 25720 | 5.00 | 5.00 | |
| 32 2-Methyl-2-propanol | 59 | 4.432 | 4.451 | -0.019 | 92 | 16447 | 50.0 | 56.4 | |
| 33 Acrylonitrile | 53 | 4.554 | 4.554 | 0.000 | 98 | 73302 | 50.0 | 53.7 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.591 | 4.584 | 0.007 | 74 | 17158 | 5.00 | 5.68 | |
| 35 Methyl tert-butyl ether | 73 | 4.603 | 4.603 | 0.000 | 84 | 38357 | 5.00 | 4.73 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 36 Hexane | 57 | 5.004 | 4.998 | 0.006 | 89 | 24902 | 5.00 | 6.42 | |
| 37 1,1-Dichloroethane | 63 | 5.211 | 5.217 | -0.006 | 96 | 28319 | 5.00 | 5.39 | |
| 38 Vinyl acetate | 43 | 5.272 | 5.272 | 0.000 | 97 | 27185 | 5.00 | 5.09 | |
| 44 2,2-Dichloropropane | 97 | 5.947 | 5.959 | -0.012 | 46 | 3769 | 5.00 | 5.63 | |
| 45 cis-1,2-Dichloroethene | 96 | 5.953 | 5.965 | -0.012 | 79 | 17858 | 5.00 | 5.17 | |
| 46 2-Butanone (MEK) | 43 | 5.984 | 5.978 | 0.006 | 98 | 50216 | 25.0 | 24.9 | |
| 49 Chlorobromomethane | 128 | 6.245 | 6.245 | 0.000 | 93 | 8216 | 5.00 | 5.35 | |
| 51 Tetrahydrofuran | 42 | 6.264 | 6.263 | 0.001 | 93 | 14858 | 10.0 | 12.7 | |
| 52 Chloroform | 83 | 6.391 | 6.391 | 0.000 | 91 | 29608 | 5.00 | 5.64 | |
| 53 1,1,1-Trichloroethane | 97 | 6.556 | 6.549 | 0.007 | 97 | 20508 | 5.00 | 5.16 | |
| 54 Cyclohexane | 56 | 6.616 | 6.622 | -0.006 | 87 | 26974 | 5.00 | 5.50 | |
| 56 Carbon tetrachloride | 117 | 6.726 | 6.726 | 0.000 | 88 | 17231 | 5.00 | 5.21 | |
| 55 1,1-Dichloropropene | 75 | 6.738 | 6.738 | 0.000 | 96 | 22014 | 5.00 | 5.13 | |
| 57 Isobutyl alcohol | 41 | 6.951 | 6.945 | 0.006 | 43 | 13122 | 125.0 | 121.7 | |
| 58 Benzene | 78 | 6.951 | 6.951 | 0.000 | 96 | 74686 | 5.00 | 5.67 | |
| 59 1,2-Dichloroethane | 62 | 7.030 | 7.030 | 0.000 | 97 | 21038 | 5.00 | 5.48 | |
| 62 n-Heptane | 43 | 7.316 | 7.316 | 0.000 | 56 | 16453 | 5.00 | 5.30 | |
| 64 Trichloroethene | 130 | 7.681 | 7.687 | -0.006 | 95 | 17490 | 5.00 | 5.28 | |
| 66 Methylcyclohexane | 83 | 7.918 | 7.918 | 0.000 | 86 | 25605 | 5.00 | 5.11 | |
| 67 1,2-Dichloropropane | 63 | 7.955 | 7.961 | -0.006 | 93 | 16316 | 5.00 | 5.32 | |
| 68 Dibromomethane | 93 | 8.046 | 8.046 | 0.000 | 90 | 8641 | 5.00 | 4.81 | |
| 70 1,4-Dioxane | 88 | 8.040 | 8.052 | -0.012 | 5 | 2333 | 100.0 | 74.8 | |
| 71 Dichlorobromomethane | 83 | 8.241 | 8.241 | 0.000 | 99 | 16257 | 5.00 | 4.61 | |
| 73 2-Chloroethyl vinyl ether | 63 | 8.551 | 8.545 | 0.006 | 92 | 18086 | 10.0 | 8.19 | |
| 74 cis-1,3-Dichloropropene | 75 | 8.691 | 8.685 | 0.006 | 95 | 19479 | 5.00 | 4.55 | |
| 75 4-Methyl-2-pentanone (MIBK) | 43 | 8.843 | 8.843 | 0.000 | 96 | 79892 | 25.0 | 26.4 | |
| 76 Toluene | 91 | 9.019 | 9.019 | 0.000 | 98 | 71883 | 5.00 | 6.12 | |
| 77 trans-1,3-Dichloropropene | 75 | 9.263 | 9.269 | -0.006 | 92 | 14443 | 5.00 | 4.52 | |
| 78 Ethyl methacrylate | 69 | 9.330 | 9.330 | 0.000 | 90 | 16030 | 5.00 | 4.16 | |
| 79 1,1,2-Trichloroethane | 97 | 9.457 | 9.457 | 0.000 | 89 | 14755 | 5.00 | 6.03 | |
| 80 Tetrachloroethene | 164 | 9.530 | 9.530 | 0.000 | 95 | 13528 | 5.00 | 6.04 | |
| 81 1,3-Dichloropropane | 76 | 9.616 | 9.615 | 0.001 | 90 | 26359 | 5.00 | 5.83 | |
| 82 2-Hexanone | 43 | 9.683 | 9.682 | 0.000 | 98 | 57842 | 25.0 | 25.0 | |
| 84 Chlorodibromomethane | 129 | 9.835 | 9.834 | 0.001 | 92 | 9414 | 5.00 | 4.55 | |
| 85 Ethylene Dibromide | 107 | 9.944 | 9.944 | 0.000 | 98 | 13462 | 5.00 | 5.36 | |
| 86 3-Chlorobenzotrifluoride | 180 | 10.413 | 10.412 | 0.001 | 90 | 25343 | 5.00 | 6.26 | |
| 87 Chlorobenzene | 112 | 10.437 | 10.437 | 0.000 | 94 | 47566 | 5.00 | 6.22 | |
| 88 4-Chlorobenzotrifluoride | 180 | 10.498 | 10.498 | 0.000 | 96 | 21933 | 5.00 | 5.87 | |
| 89 1,1,1,2-Tetrachloroethane | 131 | 10.528 | 10.528 | 0.000 | 88 | 12587 | 5.00 | 5.18 | |
| 90 Ethylbenzene | 106 | 10.534 | 10.534 | 0.000 | 98 | 22622 | 5.00 | 5.30 | |
| 91 m-Xylene & p-Xylene | 106 | 10.668 | 10.668 | 0.000 | 0 | 25553 | 5.00 | 4.90 | |
| 92 o-Xylene | 106 | 11.051 | 11.051 | 0.000 | 95 | 25240 | 5.00 | 5.08 | |
| 93 Styrene | 104 | 11.076 | 11.069 | 0.007 | 93 | 42810 | 5.00 | 5.09 | |
| 94 Bromoform | 173 | 11.252 | 11.252 | 0.000 | 92 | 6015 | 5.00 | 4.68 | |
| 96 2-Chlorobenzotrifluoride | 180 | 11.325 | 11.325 | 0.000 | 96 | 21074 | 5.00 | 5.44 | |
| 97 Isopropylbenzene | 105 | 11.422 | 11.422 | 0.000 | 96 | 64937 | 5.00 | 5.35 | |
| 100 Bromobenzene | 156 | 11.739 | 11.739 | 0.000 | 93 | 16032 | 5.00 | 5.15 | |
| 99 1,1,2,2-Tetrachloroethane | 83 | 11.739 | 11.745 | -0.006 | 77 | 20749 | 5.00 | 5.73 | |
| 102 trans-1,4-Dichloro-2-buten | 53 | 11.787 | 11.775 | 0.012 | 75 | 4170 | 5.00 | 4.44 | |
| 101 1,2,3-Trichloropropane | 110 | 11.800 | 11.793 | 0.007 | 85 | 6588 | 5.00 | 5.12 | |
| 103 N-Propylbenzene | 120 | 11.842 | 11.842 | 0.000 | 99 | 17451 | 5.00 | 4.90 | |
| 104 2-Chlorotoluene | 126 | 11.927 | 11.927 | 0.000 | 96 | 14458 | 5.00 | 4.70 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 105 3-Chlorotoluene | 126 | 11.994 | 11.994 | 0.000 | 96 | 16155 | 5.00 | 4.83 | |
| 106 1,3,5-Trimethylbenzene | 105 | 12.031 | 12.031 | 0.000 | 95 | 48645 | 5.00 | 4.78 | |
| 107 4-Chlorotoluene | 126 | 12.061 | 12.055 | 0.006 | 96 | 16940 | 5.00 | 5.10 | |
| 108 tert-Butylbenzene | 119 | 12.347 | 12.347 | 0.000 | 93 | 41329 | 5.00 | 4.85 | |
| 110 1,2,4-Trimethylbenzene | 105 | 12.408 | 12.408 | 0.000 | 97 | 50171 | 5.00 | 4.84 | |
| 111 1,2-dichloro-4-(trifluorom | 214 | 12.457 | 12.456 | 0.001 | 95 | 15090 | 5.00 | 5.82 | |
| 112 sec-Butylbenzene | 105 | 12.572 | 12.572 | 0.000 | 94 | 60251 | 5.00 | 5.07 | |
| 113 1,3-Dichlorobenzene | 146 | 12.694 | 12.688 | 0.006 | 96 | 30355 | 5.00 | 5.45 | |
| 114 4-Isopropyltoluene | 119 | 12.736 | 12.730 | 0.006 | 97 | 47431 | 5.00 | 4.79 | |
| 115 1,4-Dichlorobenzene | 146 | 12.797 | 12.797 | 0.000 | 95 | 31756 | 5.00 | 5.56 | |
| 116 2,4-Dichloro-1-(trifluorom | 214 | 12.840 | 12.828 | 0.012 | 94 | 12460 | 5.00 | 5.16 | |
| 118 2,5-Dichlorobenzotrifluori | 214 | 12.882 | 12.870 | 0.012 | 0 | 13980 | 5.00 | 5.36 | |
| 120 n-Butylbenzene | 91 | 13.156 | 13.150 | 0.006 | 96 | 39215 | 5.00 | 4.86 | |
| 121 1,2-Dichlorobenzene | 146 | 13.162 | 13.156 | 0.006 | 85 | 30057 | 5.00 | 5.66 | |
| 122 1,2-Dibromo-3-Chloropropan | 75 | 13.977 | 13.971 | 0.006 | 81 | 2690 | 5.00 | 4.57 | |
| 123 2,4- & 2,5- & 2,6- Dichlor | 125 | 14.130 | 14.117 | 0.013 | 0 | 47367 | 15.0 | 14.1 | |
| 125 2,3- & 3,4- Dichlorotoluen | 125 | 14.561 | 14.555 | 0.006 | 0 | 30402 | 10.0 | 8.74 | |
| 126 1,2,4-Trichlorobenzene | 180 | 14.847 | 14.829 | 0.018 | 92 | 12140 | 5.00 | 5.00 | |
| 127 Hexachlorobutadiene | 225 | 15.012 | 14.993 | 0.019 | 91 | 4721 | 5.00 | 5.31 | |
| 128 Naphthalene | 128 | 15.127 | 15.103 | 0.024 | 96 | 33677 | 5.00 | 4.07 | |
| 129 1,2,3-Trichlorobenzene | 180 | 15.371 | 15.346 | 0.025 | 95 | 11407 | 5.00 | 5.14 | |
| 131 2,4,5-Trichlorotoluene | 159 | 16.240 | 16.198 | 0.042 | 0 | 5175 | 5.00 | 4.91 | |
| 130 2,3,6-Trichlorotoluene | 159 | 16.338 | 16.307 | 0.031 | 88 | 4086 | 5.00 | 4.17 | |
| 149 3,4-Dichlorotoluene | 1 | | 0.000 | | | | ND | ND | |
| S 133 Xylenes, Total | 106 | | | | 0 | | 10.0 | 9.97 | |
| S 134 1,2-Dichloroethene, Total | 96 | | | | 0 | | 10.0 | 10.8 | |
| S 135 1,3-Dichloropropene, Total | 1 | | | | 0 | | 10.0 | 9.06 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

Reagents:

| | | |
|---------------------|--------------------|-----------|
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 0.20 | Units: uL |
| VOA8260VOAPRI_00263 | Amount Added: 0.20 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 4.00 | Units: uL |
| voaWVA1stRest_00017 | Amount Added: 0.20 | Units: uL |
| voaWEEmix1stR_00009 | Amount Added: 0.20 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 0.20 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 0.80 | Units: uL |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D02.D

Injection Date: 27-Jul-2017 00:51:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: IC VSTD1

Worklist Smp#: 2

Client ID:

Purge Vol: 5.000 mL

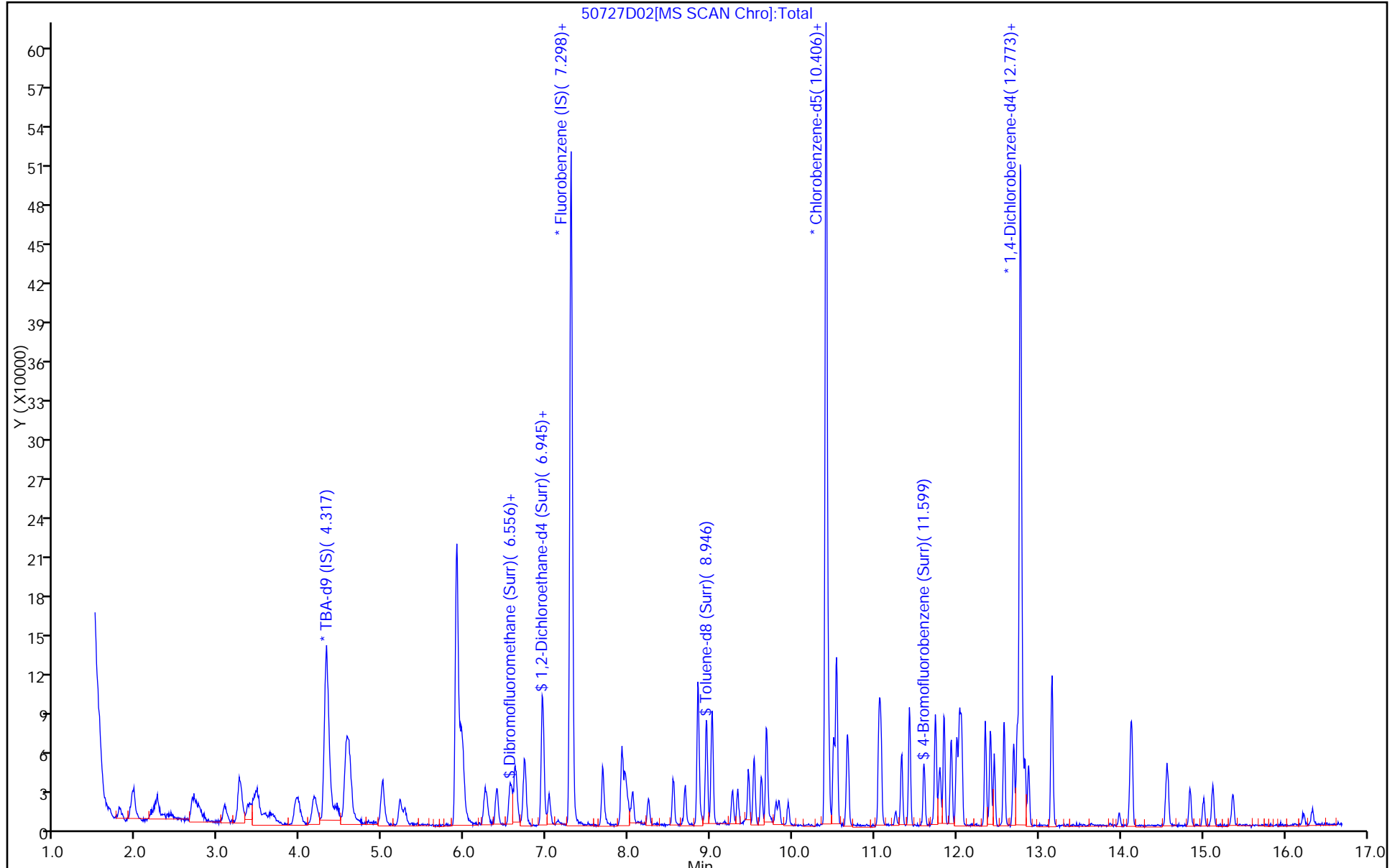
Dil. Factor: 1.0000

ALS Bottle#: 2

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh

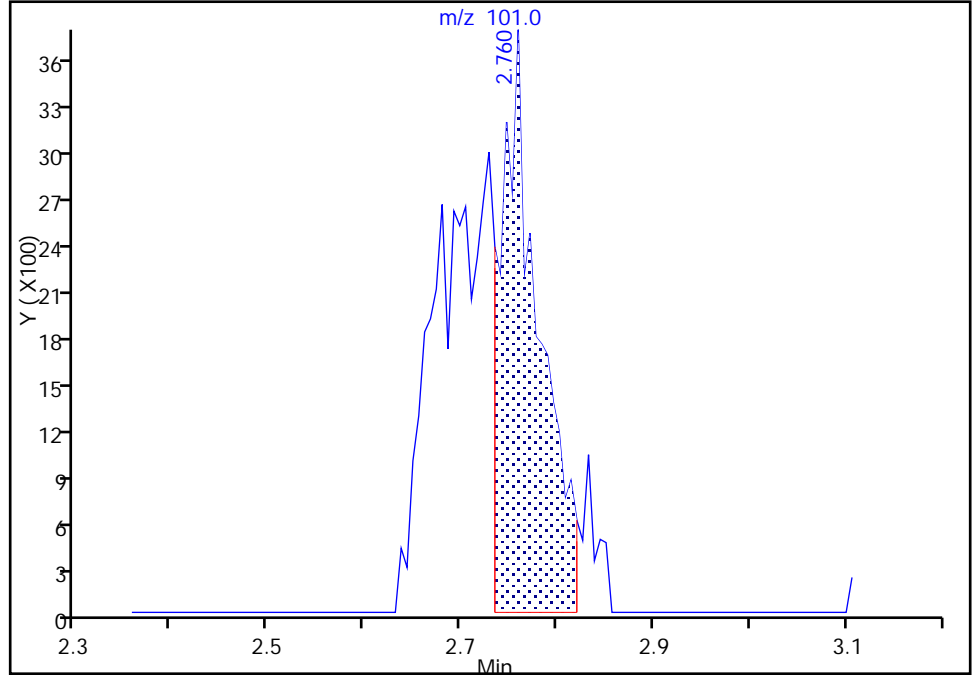
Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D02.D
Injection Date: 27-Jul-2017 00:51:30 Instrument ID: CHHP5
Lims ID: IC VSTD1
Client ID:
Operator ID: 034635 ALS Bottle#: 2 Worklist Smp#: 2
Purge Vol: 5.000 mL Dil. Factor: 1.0000
Method: MSVOA_LL_CHHP5 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm) Detector: MS SCAN

18 Trichlorofluoromethane, CAS: 75-69-4

Signal: 1

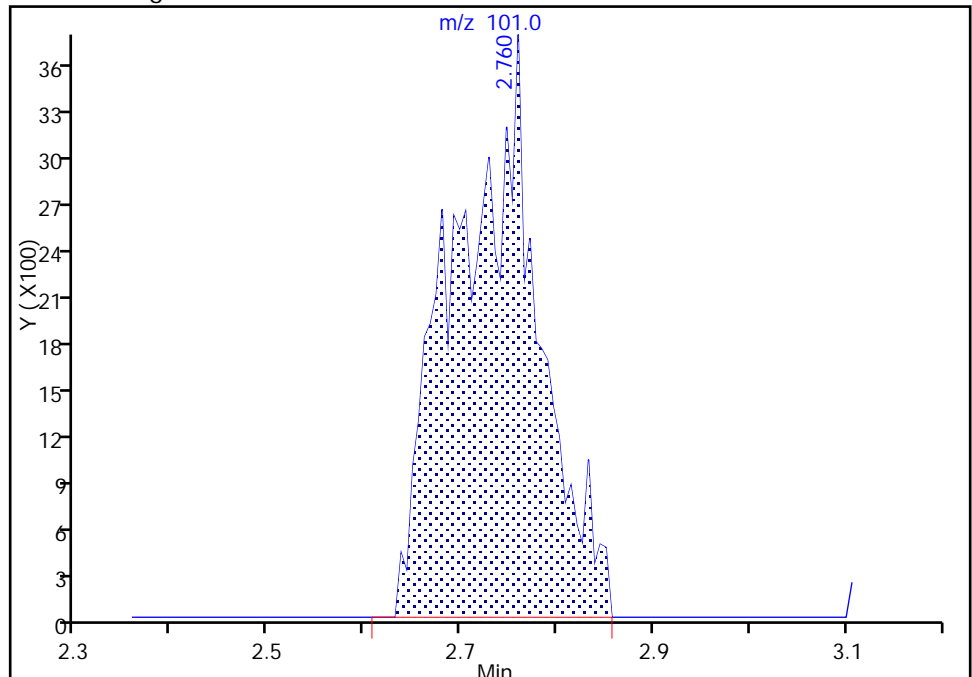
RT: 2.76
Area: 10302
Amount: 3.465076
Amount Units: ng

Processing Integration Results



RT: 2.76
Area: 22371
Amount: 5.667373
Amount Units: ng

Manual Integration Results



Reviewer: bungardf, 27-Jul-2017 03:06:53
Audit Action: Manually Integrated

Audit Reason: Poor chromatography

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D03.D
 Lims ID: IC VSTD5
 Client ID:
 Sample Type: IC Calib Level: 2
 Inject. Date: 27-Jul-2017 01:15:30 ALS Bottle#: 3 Worklist Smp#: 3
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017756-003
 Misc. Info.: IC VSTD5
 Operator ID: 034635 Instrument ID: CHHP5
 Sublist: chrom-MSVOA_LL_CHHP5*sub12
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 28-Jul-2017 01:04:47 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK029

First Level Reviewer: bungardf

Date: 27-Jul-2017 03:14:46

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.319 | 4.323 | -0.004 | 0 | 223811 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.299 | 7.298 | 0.001 | 98 | 538128 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.408 | 10.406 | 0.002 | 85 | 123664 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.775 | 12.773 | 0.002 | 94 | 168910 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.576 | 6.574 | 0.002 | 94 | 65453 | 25.0 | 25.3 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.947 | 6.945 | 0.002 | 0 | 82071 | 25.0 | 26.0 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.948 | 8.946 | 0.002 | 92 | 278432 | 25.0 | 28.3 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.600 | 11.599 | 0.001 | 87 | 94618 | 25.0 | 26.6 | |
| 11 Dichlorodifluoromethane | 85 | 1.648 | 1.646 | 0.002 | 100 | 84559 | 25.0 | 27.0 | |
| 12 Chloromethane | 50 | 1.794 | 1.804 | -0.010 | 99 | 78965 | 25.0 | 25.1 | |
| 13 Vinyl chloride | 62 | 1.946 | 1.944 | 0.002 | 98 | 82670 | 25.0 | 25.9 | |
| 14 Butadiene | 39 | 1.964 | 1.969 | -0.004 | 92 | 74553 | 25.0 | 25.7 | |
| 15 Bromomethane | 94 | 2.262 | 2.254 | 0.008 | 91 | 42224 | 25.0 | 28.0 | |
| 16 Chloroethane | 64 | 2.421 | 2.419 | 0.001 | 98 | 47273 | 25.0 | 26.9 | |
| 17 Dichlorofluoromethane | 67 | 2.700 | 2.699 | 0.001 | 97 | 119855 | 25.0 | 27.0 | |
| 18 Trichlorofluoromethane | 101 | 2.749 | 2.741 | 0.008 | 94 | 104824 | 25.0 | 26.7 | M |
| 20 Ethyl ether | 59 | 3.084 | 3.076 | 0.008 | 87 | 66542 | 25.0 | 26.1 | |
| 21 Acrolein | 56 | 3.266 | 3.252 | 0.014 | 98 | 73476 | 125.0 | 114.3 | |
| 22 1,1-Dichloroethene | 96 | 3.376 | 3.368 | 0.008 | 96 | 67928 | 25.0 | 25.8 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.436 | 3.441 | -0.005 | 93 | 73846 | 25.0 | 25.5 | |
| 24 Acetone | 43 | 3.479 | 3.477 | 0.002 | 96 | 77890 | 50.0 | 55.3 | |
| 25 Iodomethane | 142 | 3.570 | 3.562 | 0.008 | 98 | 103869 | 25.0 | 25.1 | |
| 26 Carbon disulfide | 76 | 3.649 | 3.648 | 0.001 | 99 | 131730 | 25.0 | 22.8 | |
| 28 3-Chloro-1-propene | 76 | 3.954 | 3.946 | 0.008 | 92 | 39946 | 25.0 | 23.5 | |
| 30 Methyl acetate | 43 | 3.978 | 3.976 | 0.002 | 97 | 132543 | 50.0 | 47.6 | |
| 31 Methylene Chloride | 84 | 4.166 | 4.165 | 0.001 | 88 | 84822 | 25.0 | 24.4 | |
| 32 2-Methyl-2-propanol | 59 | 4.446 | 4.451 | -0.005 | 92 | 64738 | 250.0 | 244.6 | |
| 33 Acrylonitrile | 53 | 4.562 | 4.554 | 0.008 | 100 | 336508 | 250.0 | 248.3 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.580 | 4.584 | -0.004 | 98 | 73445 | 25.0 | 24.5 | |
| 35 Methyl tert-butyl ether | 73 | 4.604 | 4.603 | 0.001 | 96 | 196780 | 25.0 | 24.4 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| 36 Hexane | 57 | 5.006 | 4.998 | 0.008 | 92 | 96542 | 25.0 | 25.1 | |
| 37 1,1-Dichloroethane | 63 | 5.219 | 5.217 | 0.002 | 96 | 133976 | 25.0 | 25.7 | |
| 38 Vinyl acetate | 43 | 5.268 | 5.272 | -0.004 | 97 | 115000 | 25.0 | 21.7 | |
| 44 2,2-Dichloropropane | 97 | 5.961 | 5.959 | 0.002 | 57 | 15889 | 25.0 | 23.9 | |
| 45 cis-1,2-Dichloroethene | 96 | 5.961 | 5.965 | -0.004 | 81 | 85931 | 25.0 | 25.0 | |
| 46 2-Butanone (MEK) | 43 | 5.985 | 5.978 | 0.007 | 93 | 105960 | 50.0 | 52.9 | |
| 49 Chlorobromomethane | 128 | 6.253 | 6.245 | 0.008 | 94 | 38047 | 25.0 | 24.9 | |
| 51 Tetrahydrofuran | 42 | 6.271 | 6.263 | 0.008 | 86 | 52866 | 50.0 | 45.3 | |
| 52 Chloroform | 83 | 6.393 | 6.391 | 0.002 | 93 | 134431 | 25.0 | 25.8 | |
| 53 1,1,1-Trichloroethane | 97 | 6.557 | 6.549 | 0.008 | 98 | 98927 | 25.0 | 25.1 | |
| 54 Cyclohexane | 56 | 6.618 | 6.622 | -0.004 | 89 | 124196 | 25.0 | 25.5 | |
| 56 Carbon tetrachloride | 117 | 6.722 | 6.726 | -0.004 | 95 | 80446 | 25.0 | 24.5 | |
| 55 1,1-Dichloropropene | 75 | 6.746 | 6.738 | 0.008 | 98 | 109851 | 25.0 | 25.8 | |
| 57 Isobutyl alcohol | 41 | 6.947 | 6.945 | 0.002 | 82 | 61305 | 625.0 | 572.5 | |
| 58 Benzene | 78 | 6.953 | 6.951 | 0.002 | 97 | 339765 | 25.0 | 26.0 | |
| 59 1,2-Dichloroethane | 62 | 7.032 | 7.030 | 0.002 | 97 | 95627 | 25.0 | 25.1 | |
| 62 n-Heptane | 43 | 7.318 | 7.316 | 0.002 | 90 | 81002 | 25.0 | 26.3 | |
| 64 Trichloroethene | 130 | 7.689 | 7.687 | 0.002 | 98 | 83072 | 25.0 | 25.2 | |
| 66 Methylcyclohexane | 83 | 7.920 | 7.918 | 0.002 | 86 | 125697 | 25.0 | 25.2 | |
| 67 1,2-Dichloropropane | 63 | 7.963 | 7.961 | 0.002 | 94 | 74777 | 25.0 | 24.5 | |
| 68 Dibromomethane | 93 | 8.048 | 8.046 | 0.002 | 95 | 45949 | 25.0 | 25.7 | |
| 70 1,4-Dioxane | 88 | 8.048 | 8.052 | -0.004 | 38 | 15162 | 500.0 | 489.4 | M |
| 71 Dichlorobromomethane | 83 | 8.242 | 8.241 | 0.001 | 98 | 84070 | 25.0 | 24.0 | |
| 73 2-Chloroethyl vinyl ether | 63 | 8.547 | 8.545 | 0.002 | 95 | 103158 | 50.0 | 47.0 | |
| 74 cis-1,3-Dichloropropene | 75 | 8.686 | 8.685 | 0.001 | 96 | 96744 | 25.0 | 22.7 | |
| 75 4-Methyl-2-pentanone (MIBK) | 43 | 8.845 | 8.843 | 0.002 | 95 | 154465 | 50.0 | 48.7 | |
| 76 Toluene | 91 | 9.015 | 9.019 | -0.004 | 98 | 351840 | 25.0 | 28.5 | |
| 77 trans-1,3-Dichloropropene | 75 | 9.270 | 9.269 | 0.001 | 92 | 79122 | 25.0 | 23.6 | |
| 78 Ethyl methacrylate | 69 | 9.325 | 9.330 | -0.005 | 88 | 96602 | 25.0 | 23.9 | |
| 79 1,1,2-Trichloroethane | 97 | 9.465 | 9.457 | 0.008 | 90 | 67966 | 25.0 | 26.5 | |
| 80 Tetrachloroethene | 164 | 9.532 | 9.530 | 0.002 | 97 | 67579 | 25.0 | 28.7 | |
| 81 1,3-Dichloropropane | 76 | 9.617 | 9.615 | 0.002 | 89 | 127957 | 25.0 | 26.9 | |
| 82 2-Hexanone | 43 | 9.678 | 9.682 | -0.004 | 95 | 122936 | 50.0 | 50.5 | |
| 84 Chlorodibromomethane | 129 | 9.836 | 9.834 | 0.002 | 89 | 53302 | 25.0 | 24.5 | |
| 85 Ethylene Dibromide | 107 | 9.946 | 9.944 | 0.002 | 100 | 67745 | 25.0 | 25.7 | |
| 86 3-Chlorobenzotrifluoride | 180 | 10.408 | 10.412 | -0.004 | 95 | 109109 | 25.0 | 25.7 | |
| 87 Chlorobenzene | 112 | 10.432 | 10.437 | -0.005 | 95 | 217561 | 25.0 | 27.1 | |
| 88 4-Chlorobenzotrifluoride | 180 | 10.499 | 10.498 | 0.001 | 95 | 101825 | 25.0 | 26.0 | |
| 89 1,1,1,2-Tetrachloroethane | 131 | 10.530 | 10.528 | 0.002 | 92 | 65901 | 25.0 | 25.8 | |
| 90 Ethylbenzene | 106 | 10.536 | 10.534 | 0.002 | 98 | 120759 | 25.0 | 26.9 | |
| 91 m-Xylene & p-Xylene | 106 | 10.670 | 10.668 | 0.002 | 0 | 151114 | 25.0 | 27.6 | |
| 92 o-Xylene | 106 | 11.053 | 11.051 | 0.002 | 96 | 138375 | 25.0 | 26.5 | |
| 93 Styrene | 104 | 11.071 | 11.069 | 0.002 | 95 | 242031 | 25.0 | 27.4 | |
| 94 Bromoform | 173 | 11.254 | 11.252 | 0.002 | 97 | 30000 | 25.0 | 22.2 | |
| 96 2-Chlorobenzotrifluoride | 180 | 11.327 | 11.325 | 0.002 | 97 | 107103 | 25.0 | 26.3 | |
| 97 Isopropylbenzene | 105 | 11.424 | 11.422 | 0.002 | 96 | 356966 | 25.0 | 28.0 | |
| 100 Bromobenzene | 156 | 11.734 | 11.739 | -0.005 | 95 | 83376 | 25.0 | 25.4 | |
| 99 1,1,2,2-Tetrachloroethane | 83 | 11.740 | 11.745 | -0.005 | 94 | 100341 | 25.0 | 26.4 | |
| 102 trans-1,4-Dichloro-2-buten | 53 | 11.777 | 11.775 | 0.002 | 77 | 23168 | 25.0 | 23.4 | |
| 101 1,2,3-Trichloropropane | 110 | 11.789 | 11.793 | -0.004 | 86 | 32588 | 25.0 | 24.1 | |
| 103 N-Propylbenzene | 120 | 11.838 | 11.842 | -0.004 | 99 | 95261 | 25.0 | 25.4 | |
| 104 2-Chlorotoluene | 126 | 11.929 | 11.927 | 0.002 | 96 | 83234 | 25.0 | 25.7 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 105 3-Chlorotoluene | 126 | 11.996 | 11.994 | 0.002 | 96 | 87067 | 25.0 | 24.7 | |
| 106 1,3,5-Trimethylbenzene | 105 | 12.026 | 12.031 | -0.005 | 95 | 290219 | 25.0 | 27.1 | |
| 107 4-Chlorotoluene | 126 | 12.057 | 12.055 | 0.002 | 96 | 88877 | 25.0 | 25.4 | |
| 108 tert-Butylbenzene | 119 | 12.349 | 12.347 | 0.002 | 93 | 236619 | 25.0 | 26.4 | |
| 110 1,2,4-Trimethylbenzene | 105 | 12.410 | 12.408 | 0.002 | 97 | 288545 | 25.0 | 26.5 | |
| 111 1,2-dichloro-4-(trifluorom | 214 | 12.452 | 12.456 | -0.004 | 96 | 64854 | 25.0 | 23.8 | |
| 112 sec-Butylbenzene | 105 | 12.574 | 12.572 | 0.002 | 94 | 336681 | 25.0 | 26.9 | |
| 113 1,3-Dichlorobenzene | 146 | 12.689 | 12.688 | 0.001 | 97 | 151590 | 25.0 | 25.9 | |
| 114 4-Isopropyltoluene | 119 | 12.732 | 12.730 | 0.002 | 97 | 277710 | 25.0 | 26.7 | |
| 115 1,4-Dichlorobenzene | 146 | 12.799 | 12.797 | 0.002 | 95 | 154714 | 25.0 | 25.7 | |
| 116 2,4-Dichloro-1-(trifluorom | 214 | 12.829 | 12.828 | 0.001 | 96 | 64892 | 25.0 | 25.5 | |
| 118 2,5-Dichlorobenzotrifluori | 214 | 12.872 | 12.870 | 0.002 | 0 | 67486 | 25.0 | 24.6 | |
| 120 n-Butylbenzene | 91 | 13.152 | 13.150 | 0.002 | 98 | 221777 | 25.0 | 26.1 | |
| 121 1,2-Dichlorobenzene | 146 | 13.158 | 13.156 | 0.002 | 98 | 145778 | 25.0 | 26.1 | |
| 122 1,2-Dibromo-3-Chloropropan | 75 | 13.973 | 13.971 | 0.002 | 83 | 14158 | 25.0 | 22.8 | |
| 123 2,4- & 2,5- & 2,6- Dichlor | 125 | 14.119 | 14.117 | 0.002 | 0 | 260387 | 75.0 | 73.5 | |
| 125 2,3- & 3,4- Dichlorotoluen | 125 | 14.557 | 14.555 | 0.002 | 0 | 173187 | 50.0 | 47.3 | |
| 126 1,2,4-Trichlorobenzene | 180 | 14.837 | 14.829 | 0.008 | 94 | 60672 | 25.0 | 23.7 | |
| 127 Hexachlorobutadiene | 225 | 14.995 | 14.993 | 0.002 | 98 | 24054 | 25.0 | 25.7 | |
| 128 Naphthalene | 128 | 15.111 | 15.103 | 0.008 | 97 | 191971 | 25.0 | 22.1 | |
| 129 1,2,3-Trichlorobenzene | 180 | 15.348 | 15.346 | 0.002 | 95 | 57325 | 25.0 | 24.6 | |
| 131 2,4,5-Trichlorotoluene | 159 | 16.200 | 16.198 | 0.002 | 0 | 23799 | 25.0 | 21.5 | |
| 130 2,3,6-Trichlorotoluene | 159 | 16.309 | 16.307 | 0.002 | 95 | 23065 | 25.0 | 22.3 | |
| 149 3,4-Dichlorotoluene | 1 | | 0.000 | | | | ND | ND | |
| S 134 1,2-Dichloroethene, Total | 96 | | | | 0 | | 50.0 | 49.5 | |
| S 133 Xylenes, Total | 106 | | | | 0 | | 50.0 | 54.1 | |
| S 135 1,3-Dichloropropene, Total | 1 | | | | 0 | | 50.0 | 46.3 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

Reagents:

| | | |
|---------------------|--------------------|-----------|
| VOA8260VOAPRI_00263 | Amount Added: 1.00 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 1.00 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 5.00 | Units: uL |
| voaWVA1stRest_00017 | Amount Added: 1.00 | Units: uL |
| voaWEEmix1stR_00009 | Amount Added: 1.00 | Units: uL |
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 1.00 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 1.00 | Units: uL |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D03.D

Injection Date: 27-Jul-2017 01:15:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: IC VSTD5

Worklist Smp#: 3

Client ID:

Purge Vol: 5.000 mL

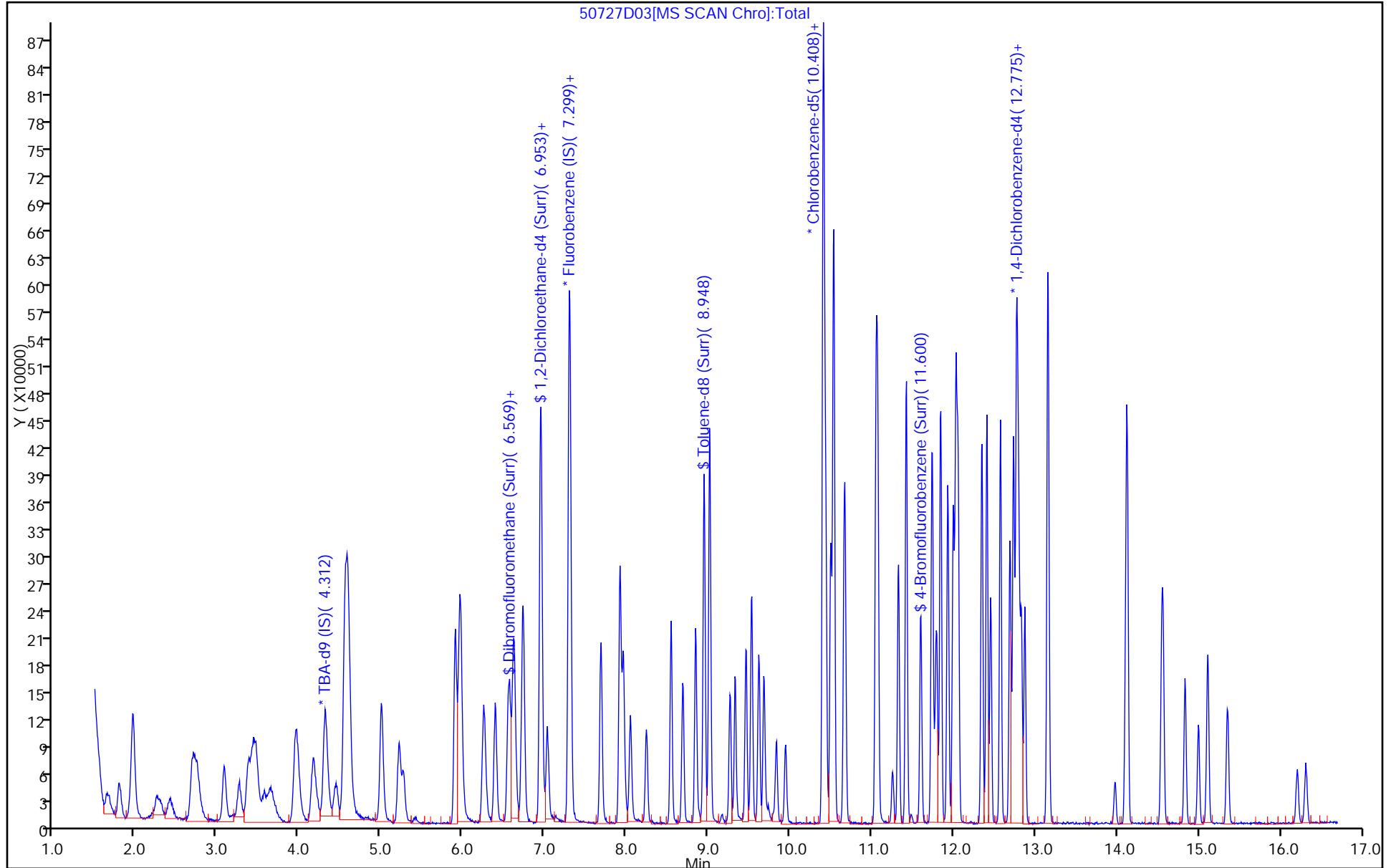
Dil. Factor: 1.0000

ALS Bottle#: 3

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh

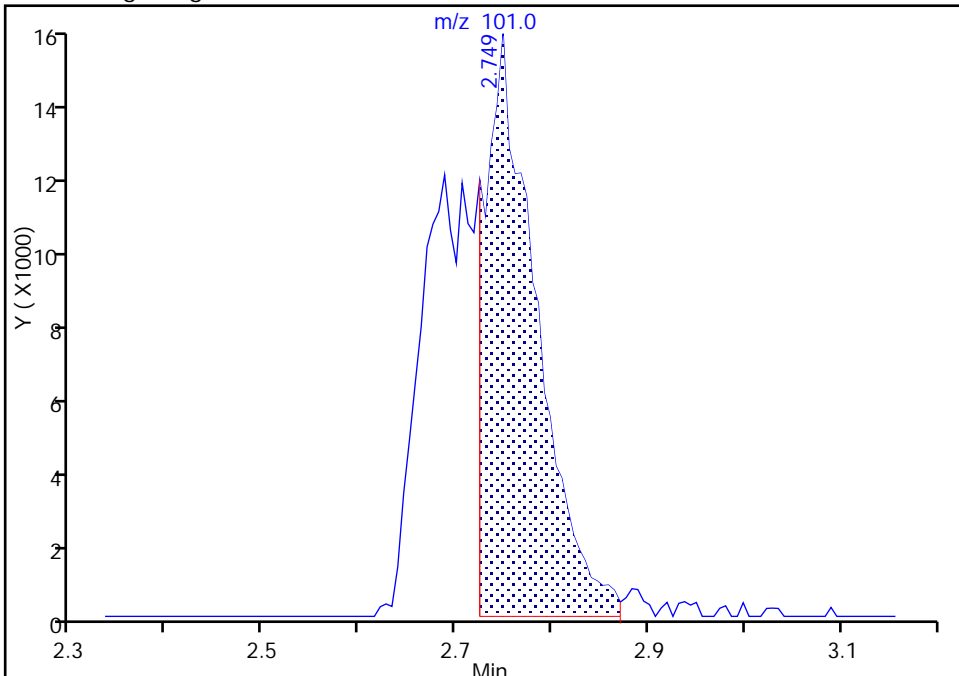
Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D03.D
Injection Date: 27-Jul-2017 01:15:30 Instrument ID: CHHP5
Lims ID: IC VSTD5
Client ID:
Operator ID: 034635 ALS Bottle#: 3 Worklist Smp#: 3
Purge Vol: 5.000 mL Dil. Factor: 1.0000
Method: MSVOA_LL_CHHP5 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm) Detector: MS SCAN

18 Trichlorofluoromethane, CAS: 75-69-4

Signal: 1

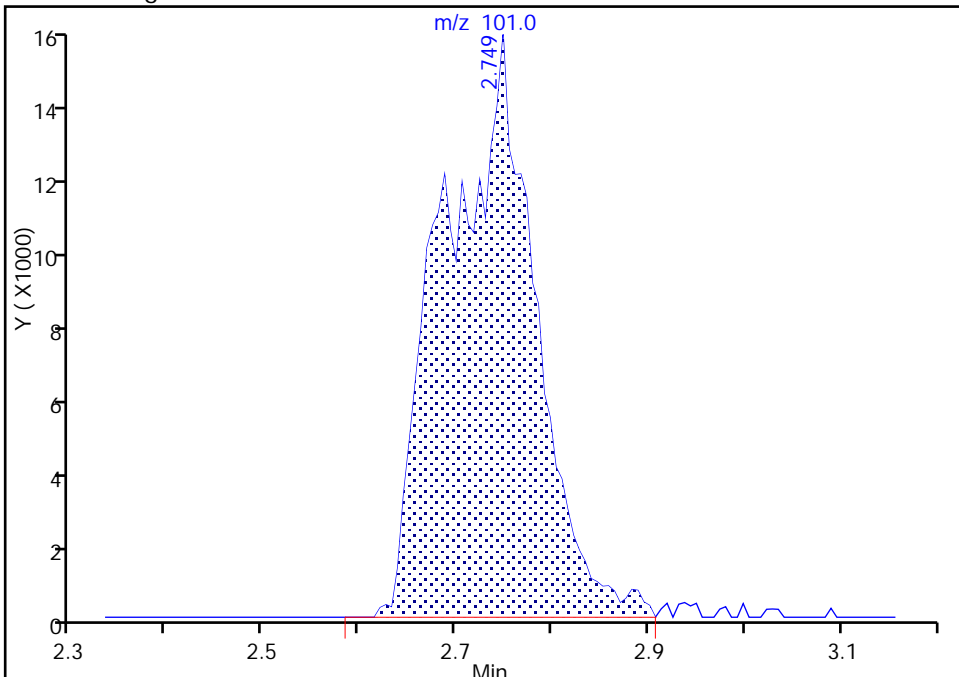
RT: 2.75
Area: 59636
Amount: 17.371088
Amount Units: ng

Processing Integration Results



RT: 2.75
Area: 104824
Amount: 26.731985
Amount Units: ng

Manual Integration Results



Reviewer: bungardf, 27-Jul-2017 03:13:52
Audit Action: Manually Integrated

TestAmerica Pittsburgh

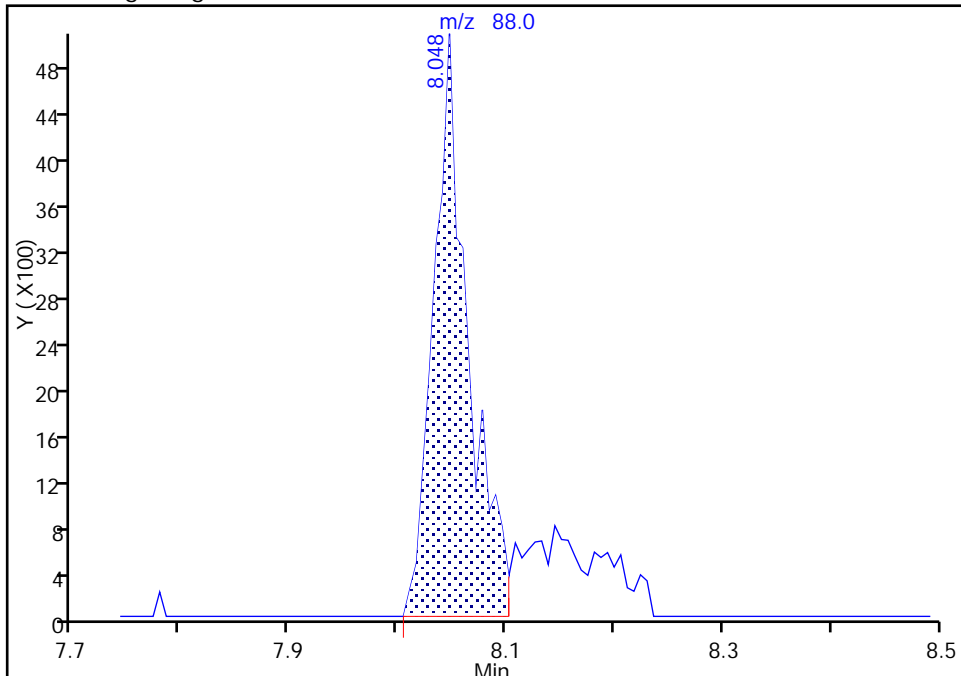
Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D03.D
Injection Date: 27-Jul-2017 01:15:30 Instrument ID: CHHP5
Lims ID: IC VSTD5
Client ID:
Operator ID: 034635 ALS Bottle#: 3 Worklist Smp#: 3
Purge Vol: 5.000 mL Dil. Factor: 1.0000
Method: MSVOA_LL_CHHP5 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm) Detector: MS SCAN

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

Signal: 1

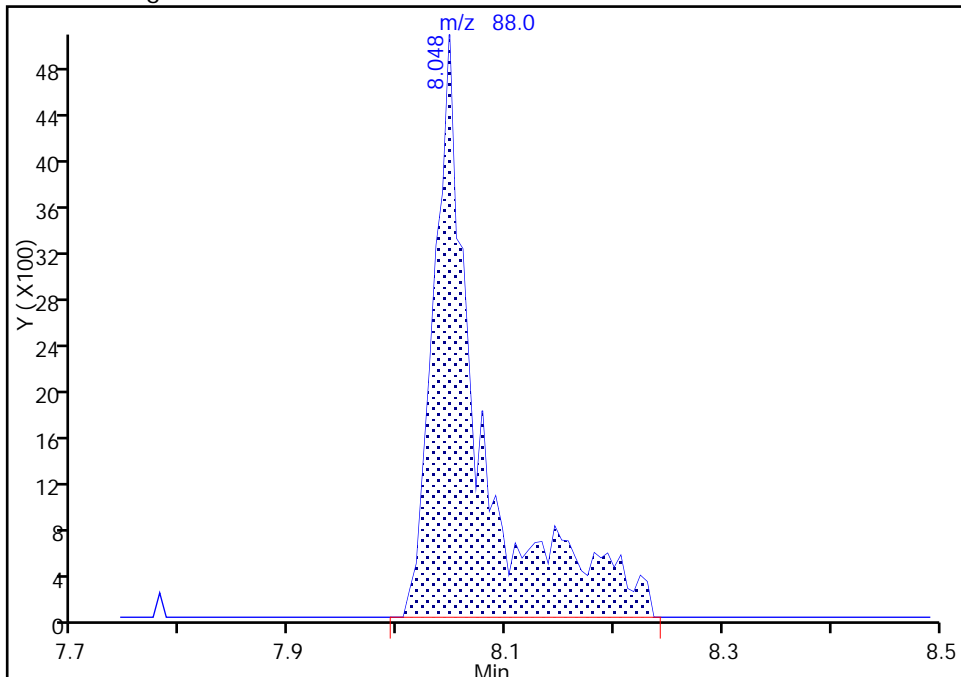
RT: 8.05
Area: 11273
Amount: 403.3803
Amount Units: ng

Processing Integration Results



RT: 8.05
Area: 15162
Amount: 489.3788
Amount Units: ng

Manual Integration Results



Reviewer: bungardf, 27-Jul-2017 03:14:22
Audit Action: Manually Integrated

Audit Reason: Poor chromatography

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D04.D
 Lims ID: ICIS VSTD10
 Client ID:
 Sample Type: ICIS Calib Level: 3
 Inject. Date: 27-Jul-2017 01:39:30 ALS Bottle#: 4 Worklist Smp#: 4
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017756-004
 Misc. Info.: ICIS VSTD10
 Operator ID: 034635 Instrument ID: CHHP5
 Sublist: chrom-MSVOA_LL_CHHP5*sub12
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 28-Jul-2017 01:04:50 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last Ical File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK029

First Level Reviewer: bungardf

Date: 27-Jul-2017 03:16:02

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.328 | 4.328 | 0.000 | 0 | 240414 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.297 | 7.297 | 0.000 | 99 | 539679 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.405 | 10.405 | 0.000 | 86 | 132843 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.772 | 12.772 | 0.000 | 94 | 174621 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.573 | 6.573 | 0.000 | 94 | 127700 | 50.0 | 49.2 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.944 | 6.944 | 0.000 | 0 | 159071 | 50.0 | 50.2 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.951 | 8.951 | 0.000 | 92 | 541748 | 50.0 | 51.2 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.598 | 11.598 | 0.000 | 87 | 191158 | 50.0 | 50.1 | |
| 11 Dichlorodifluoromethane | 85 | 1.663 | 1.663 | 0.000 | 99 | 159957 | 50.0 | 51.0 | |
| 12 Chloromethane | 50 | 1.797 | 1.797 | 0.000 | 99 | 154943 | 50.0 | 49.1 | |
| 13 Vinyl chloride | 62 | 1.955 | 1.955 | 0.000 | 98 | 162634 | 50.0 | 50.8 | |
| 14 Butadiene | 39 | 1.968 | 1.968 | 0.000 | 94 | 143576 | 50.0 | 49.4 | |
| 15 Bromomethane | 94 | 2.272 | 2.272 | 0.000 | 89 | 81346 | 50.0 | 53.8 | |
| 16 Chloroethane | 64 | 2.424 | 2.424 | 0.000 | 98 | 86601 | 50.0 | 49.2 | |
| 17 Dichlorofluoromethane | 67 | 2.710 | 2.710 | 0.000 | 96 | 224450 | 50.0 | 50.4 | |
| 18 Trichlorofluoromethane | 101 | 2.746 | 2.746 | 0.000 | 97 | 205127 | 50.0 | 52.2 | M |
| 20 Ethyl ether | 59 | 3.087 | 3.087 | 0.000 | 89 | 126496 | 50.0 | 49.4 | |
| 21 Acrolein | 56 | 3.269 | 3.269 | 0.000 | 99 | 101829 | 150.0 | 158.0 | |
| 22 1,1-Dichloroethene | 96 | 3.373 | 3.373 | 0.000 | 83 | 131576 | 50.0 | 49.8 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.440 | 3.440 | 0.000 | 91 | 141127 | 50.0 | 48.7 | |
| 24 Acetone | 43 | 3.482 | 3.482 | 0.000 | 100 | 149782 | 100.0 | 106.1 | |
| 25 Iodomethane | 142 | 3.580 | 3.580 | 0.000 | 99 | 200342 | 50.0 | 48.3 | |
| 26 Carbon disulfide | 76 | 3.659 | 3.659 | 0.000 | 98 | 266935 | 50.0 | 46.0 | |
| 28 3-Chloro-1-propene | 76 | 3.951 | 3.951 | 0.000 | 92 | 83167 | 50.0 | 48.7 | |
| 30 Methyl acetate | 43 | 3.975 | 3.975 | 0.000 | 97 | 283974 | 100.0 | 101.6 | |
| 31 Methylene Chloride | 84 | 4.170 | 4.170 | 0.000 | 90 | 164284 | 50.0 | 50.2 | |
| 32 2-Methyl-2-propanol | 59 | 4.450 | 4.450 | 0.000 | 93 | 139891 | 500.0 | 492.0 | |
| 33 Acrylonitrile | 53 | 4.559 | 4.559 | 0.000 | 99 | 708552 | 500.0 | 521.4 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.584 | 4.584 | 0.000 | 97 | 147191 | 50.0 | 48.9 | |
| 35 Methyl tert-butyl ether | 73 | 4.608 | 4.608 | 0.000 | 96 | 390184 | 50.0 | 48.3 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 36 Hexane | 57 | 5.003 | 5.003 | 0.000 | 93 | 186124 | 50.0 | 48.2 | |
| 37 1,1-Dichloroethane | 63 | 5.222 | 5.222 | 0.000 | 96 | 261874 | 50.0 | 50.0 | |
| 38 Vinyl acetate | 43 | 5.271 | 5.271 | 0.000 | 97 | 245879 | 50.0 | 46.2 | |
| 44 2,2-Dichloropropane | 97 | 5.958 | 5.958 | 0.000 | 72 | 31118 | 50.0 | 46.7 | |
| 45 cis-1,2-Dichloroethene | 96 | 5.971 | 5.971 | 0.000 | 79 | 172690 | 50.0 | 50.2 | |
| 46 2-Butanone (MEK) | 43 | 5.977 | 5.977 | 0.000 | 98 | 214731 | 100.0 | 106.9 | |
| 49 Chlorobromomethane | 128 | 6.250 | 6.250 | 0.000 | 95 | 75687 | 50.0 | 49.5 | |
| 51 Tetrahydrofuran | 42 | 6.269 | 6.269 | 0.000 | 89 | 117485 | 100.0 | 100.4 | |
| 52 Chloroform | 83 | 6.396 | 6.396 | 0.000 | 92 | 254354 | 50.0 | 48.7 | |
| 53 1,1,1-Trichloroethane | 97 | 6.555 | 6.555 | 0.000 | 98 | 196286 | 50.0 | 49.6 | |
| 54 Cyclohexane | 56 | 6.621 | 6.621 | 0.000 | 89 | 239333 | 50.0 | 49.0 | |
| 56 Carbon tetrachloride | 117 | 6.719 | 6.719 | 0.000 | 97 | 162849 | 50.0 | 49.5 | |
| 55 1,1-Dichloropropene | 75 | 6.743 | 6.743 | 0.000 | 97 | 215336 | 50.0 | 50.4 | |
| 57 Isobutyl alcohol | 41 | 6.950 | 6.950 | 0.000 | 84 | 136973 | 1250.0 | 1275.5 | |
| 58 Benzene | 78 | 6.950 | 6.950 | 0.000 | 97 | 669098 | 50.0 | 51.0 | |
| 59 1,2-Dichloroethane | 62 | 7.035 | 7.035 | 0.000 | 97 | 190422 | 50.0 | 49.8 | |
| 62 n-Heptane | 43 | 7.315 | 7.315 | 0.000 | 86 | 154370 | 50.0 | 50.0 | |
| 64 Trichloroethene | 130 | 7.692 | 7.692 | 0.000 | 98 | 164695 | 50.0 | 49.9 | |
| 66 Methylcyclohexane | 83 | 7.917 | 7.917 | 0.000 | 86 | 253511 | 50.0 | 50.8 | |
| 67 1,2-Dichloropropane | 63 | 7.960 | 7.960 | 0.000 | 94 | 150135 | 50.0 | 49.1 | |
| 68 Dibromomethane | 93 | 8.045 | 8.045 | 0.000 | 95 | 88395 | 50.0 | 49.4 | |
| 70 1,4-Dioxane | 88 | 8.051 | 8.051 | 0.000 | 40 | 33209 | 1000.0 | 1068.8 | M |
| 71 Dichlorobromomethane | 83 | 8.246 | 8.246 | 0.000 | 99 | 171049 | 50.0 | 48.7 | |
| 73 2-Chloroethyl vinyl ether | 63 | 8.544 | 8.544 | 0.000 | 92 | 219328 | 100.0 | 99.7 | |
| 74 cis-1,3-Dichloropropene | 75 | 8.690 | 8.690 | 0.000 | 95 | 204344 | 50.0 | 47.9 | |
| 75 4-Methyl-2-pentanone (MIBK) | 43 | 8.848 | 8.848 | 0.000 | 96 | 361112 | 100.0 | 106.0 | |
| 76 Toluene | 91 | 9.018 | 9.018 | 0.000 | 99 | 692901 | 50.0 | 52.3 | |
| 77 trans-1,3-Dichloropropene | 75 | 9.268 | 9.268 | 0.000 | 93 | 170710 | 50.0 | 47.4 | |
| 78 Ethyl methacrylate | 69 | 9.329 | 9.329 | 0.000 | 88 | 222171 | 50.0 | 51.1 | |
| 79 1,1,2-Trichloroethane | 97 | 9.456 | 9.456 | 0.000 | 90 | 138196 | 50.0 | 50.1 | |
| 80 Tetrachloroethene | 164 | 9.535 | 9.535 | 0.000 | 97 | 126273 | 50.0 | 50.0 | |
| 81 1,3-Dichloropropane | 76 | 9.621 | 9.621 | 0.000 | 89 | 256477 | 50.0 | 50.3 | |
| 82 2-Hexanone | 43 | 9.681 | 9.681 | 0.000 | 94 | 278579 | 100.0 | 106.6 | |
| 84 Chlorodibromomethane | 129 | 9.834 | 9.834 | 0.000 | 90 | 114911 | 50.0 | 49.3 | |
| 85 Ethylene Dibromide | 107 | 9.943 | 9.943 | 0.000 | 98 | 142489 | 50.0 | 50.3 | |
| 86 3-Chlorobenzotrifluoride | 180 | 10.411 | 10.411 | 0.000 | 93 | 222871 | 50.0 | 48.8 | |
| 87 Chlorobenzene | 112 | 10.436 | 10.436 | 0.000 | 95 | 431311 | 50.0 | 50.0 | |
| 88 4-Chlorobenzotrifluoride | 180 | 10.497 | 10.497 | 0.000 | 96 | 207774 | 50.0 | 49.3 | |
| 89 1,1,1,2-Tetrachloroethane | 131 | 10.533 | 10.533 | 0.000 | 94 | 137710 | 50.0 | 50.2 | |
| 90 Ethylbenzene | 106 | 10.533 | 10.533 | 0.000 | 98 | 249792 | 50.0 | 51.9 | |
| 91 m-Xylene & p-Xylene | 106 | 10.667 | 10.667 | 0.000 | 0 | 306948 | 50.0 | 52.2 | |
| 92 o-Xylene | 106 | 11.050 | 11.050 | 0.000 | 96 | 288885 | 50.0 | 51.5 | |
| 93 Styrene | 104 | 11.068 | 11.068 | 0.000 | 95 | 498873 | 50.0 | 52.6 | |
| 94 Bromoform | 173 | 11.257 | 11.257 | 0.000 | 96 | 67829 | 50.0 | 46.8 | |
| 96 2-Chlorobenzotrifluoride | 180 | 11.324 | 11.324 | 0.000 | 97 | 216286 | 50.0 | 49.5 | |
| 97 Isopropylbenzene | 105 | 11.421 | 11.421 | 0.000 | 95 | 726432 | 50.0 | 53.1 | |
| 100 Bromobenzene | 156 | 11.738 | 11.738 | 0.000 | 94 | 163748 | 50.0 | 48.3 | |
| 99 1,1,2,2-Tetrachloroethane | 83 | 11.738 | 11.738 | 0.000 | 95 | 211912 | 50.0 | 51.9 | |
| 102 trans-1,4-Dichloro-2-buten | 53 | 11.780 | 11.780 | 0.000 | 83 | 49334 | 50.0 | 48.3 | |
| 101 1,2,3-Trichloropropane | 110 | 11.792 | 11.792 | 0.000 | 85 | 72643 | 50.0 | 51.9 | |
| 103 N-Propylbenzene | 120 | 11.841 | 11.841 | 0.000 | 98 | 198029 | 50.0 | 51.1 | |
| 104 2-Chlorotoluene | 126 | 11.926 | 11.926 | 0.000 | 97 | 167713 | 50.0 | 50.1 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 105 3-Chlorotoluene | 126 | 11.999 | 11.999 | 0.000 | 96 | 185343 | 50.0 | 50.9 | |
| 106 1,3,5-Trimethylbenzene | 105 | 12.030 | 12.030 | 0.000 | 94 | 578518 | 50.0 | 52.2 | |
| 107 4-Chlorotoluene | 126 | 12.054 | 12.054 | 0.000 | 96 | 180584 | 50.0 | 50.0 | |
| 108 tert-Butylbenzene | 119 | 12.346 | 12.346 | 0.000 | 93 | 480729 | 50.0 | 51.9 | |
| 110 1,2,4-Trimethylbenzene | 105 | 12.407 | 12.407 | 0.000 | 97 | 588662 | 50.0 | 52.3 | |
| 111 1,2-dichloro-4-(trifluorom | 214 | 12.456 | 12.456 | 0.000 | 97 | 138659 | 50.0 | 49.1 | |
| 112 sec-Butylbenzene | 105 | 12.571 | 12.571 | 0.000 | 94 | 679839 | 50.0 | 52.6 | |
| 113 1,3-Dichlorobenzene | 146 | 12.687 | 12.687 | 0.000 | 97 | 305374 | 50.0 | 50.4 | |
| 114 4-Isopropyltoluene | 119 | 12.735 | 12.735 | 0.000 | 97 | 570403 | 50.0 | 53.0 | |
| 115 1,4-Dichlorobenzene | 146 | 12.796 | 12.796 | 0.000 | 95 | 315614 | 50.0 | 50.8 | |
| 116 2,4-Dichloro-1-(trifluorom | 214 | 12.827 | 12.827 | 0.000 | 95 | 125268 | 50.0 | 47.7 | |
| 118 2,5-Dichlorobenzotrifluori | 214 | 12.875 | 12.875 | 0.000 | 0 | 140272 | 50.0 | 49.4 | |
| 120 n-Butylbenzene | 91 | 13.149 | 13.149 | 0.000 | 98 | 454742 | 50.0 | 51.8 | |
| 121 1,2-Dichlorobenzene | 146 | 13.161 | 13.161 | 0.000 | 98 | 290492 | 50.0 | 50.3 | |
| 122 1,2-Dibromo-3-Chloropropan | 75 | 13.976 | 13.976 | 0.000 | 85 | 30986 | 50.0 | 48.4 | |
| 123 2,4- & 2,5- & 2,6- Dichlor | 125 | 14.122 | 14.122 | 0.000 | 0 | 566788 | 150.0 | 154.8 | |
| 125 2,3- & 3,4- Dichlorotoluen | 125 | 14.554 | 14.554 | 0.000 | 0 | 380181 | 100.0 | 100.4 | |
| 126 1,2,4-Trichlorobenzene | 180 | 14.834 | 14.834 | 0.000 | 93 | 134753 | 50.0 | 51.0 | |
| 127 Hexachlorobutadiene | 225 | 14.992 | 14.992 | 0.000 | 97 | 49048 | 50.0 | 50.8 | |
| 128 Naphthalene | 128 | 15.108 | 15.108 | 0.000 | 97 | 465533 | 50.0 | 51.7 | |
| 129 1,2,3-Trichlorobenzene | 180 | 15.351 | 15.351 | 0.000 | 95 | 117120 | 50.0 | 48.5 | |
| 131 2,4,5-Trichlorotoluene | 159 | 16.203 | 16.203 | 0.000 | 0 | 53498 | 50.0 | 46.6 | |
| 130 2,3,6-Trichlorotoluene | 159 | 16.312 | 16.312 | 0.000 | 97 | 53869 | 50.0 | 50.5 | |
| 149 3,4-Dichlorotoluene | 1 | | 0.000 | | | | ND | ND | |
| S 133 Xylenes, Total | 106 | | | | 0 | | 100.0 | 103.7 | |
| S 134 1,2-Dichloroethene, Total | 96 | | | | 0 | | 100.0 | 99.0 | |
| S 135 1,3-Dichloropropene, Total | 1 | | | | 0 | | 100.0 | 95.2 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

Reagents:

| | | |
|---------------------|--------------------|-----------|
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 2.00 | Units: uL |
| VOA8260VOAPRI_00263 | Amount Added: 2.00 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 2.00 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 6.00 | Units: uL |
| voaWVA1stRest_00017 | Amount Added: 2.00 | Units: uL |
| voaWEEmix1stR_00009 | Amount Added: 2.00 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 2.00 | Units: uL |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D04.D

Injection Date: 27-Jul-2017 01:39:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: ICIS VSTD10

Worklist Smp#: 4

Client ID:

Purge Vol: 5.000 mL

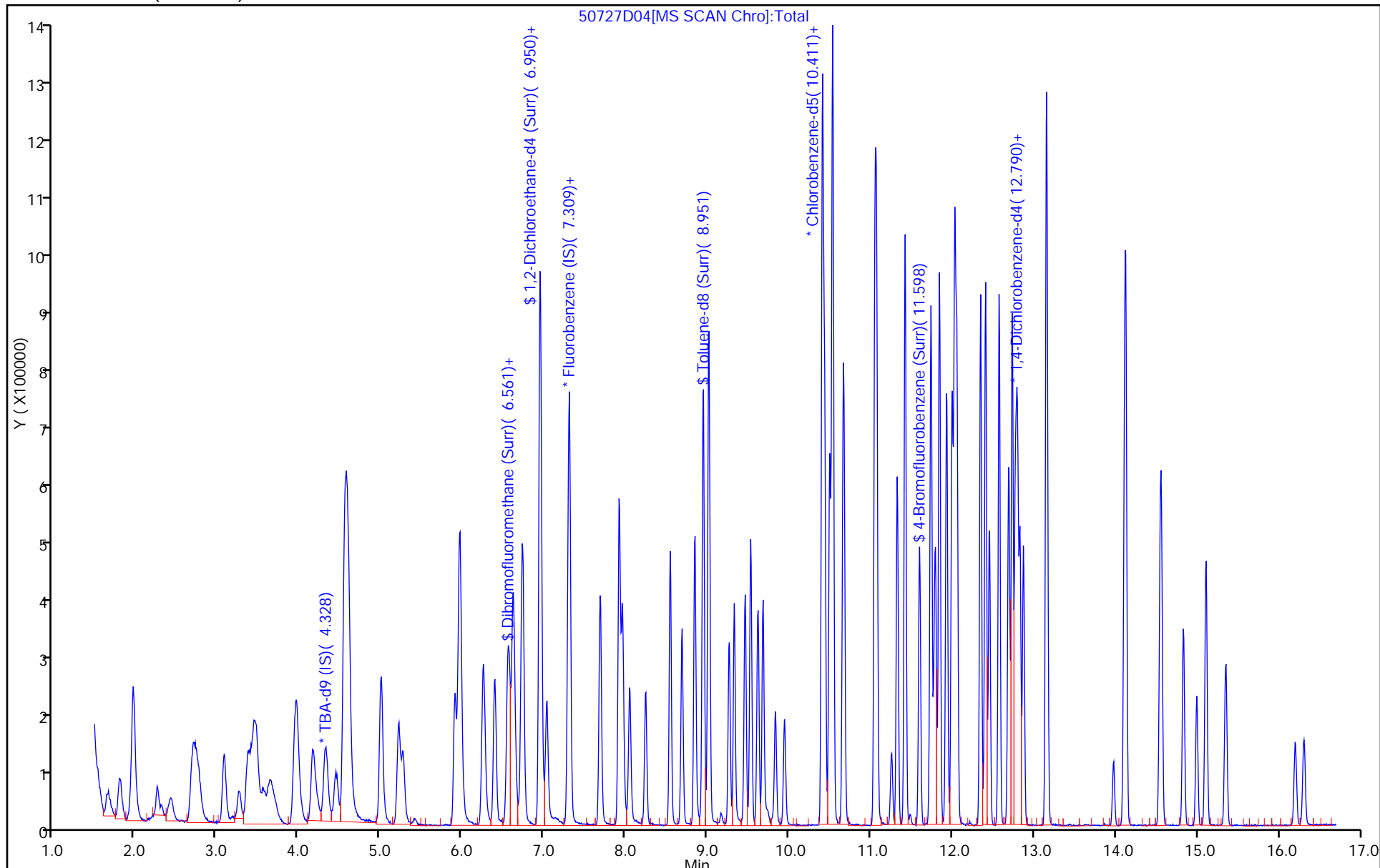
Dil. Factor: 1.0000

ALS Bottle#: 4

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh

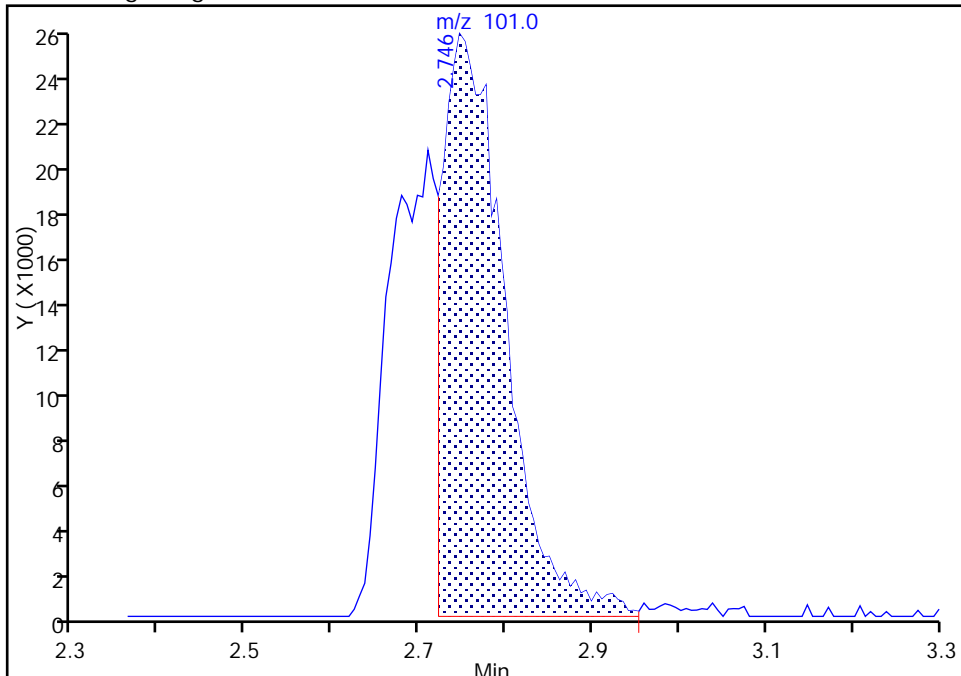
Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D04.D
Injection Date: 27-Jul-2017 01:39:30 Instrument ID: CHHP5
Lims ID: ICIS VSTD10
Client ID:
Operator ID: 034635 ALS Bottle#: 4 Worklist Smp#: 4
Purge Vol: 5.000 mL Dil. Factor: 1.0000
Method: MSVOA_LL_CHHP5 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm) Detector: MS SCAN

18 Trichlorofluoromethane, CAS: 75-69-4

Signal: 1

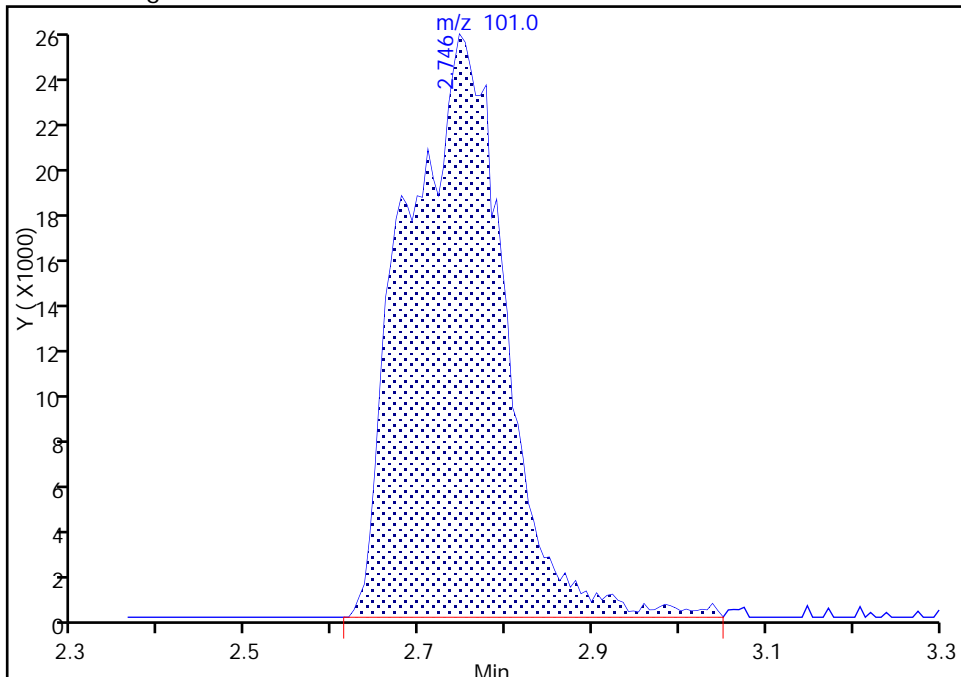
RT: 2.75
Area: 129465
Amount: 34.020484
Amount Units: ng

Processing Integration Results



RT: 2.75
Area: 205127
Amount: 52.160696
Amount Units: ng

Manual Integration Results



Reviewer: bungardf, 27-Jul-2017 03:15:11
Audit Action: Manually Integrated

Audit Reason: Poor chromatography

TestAmerica Pittsburgh

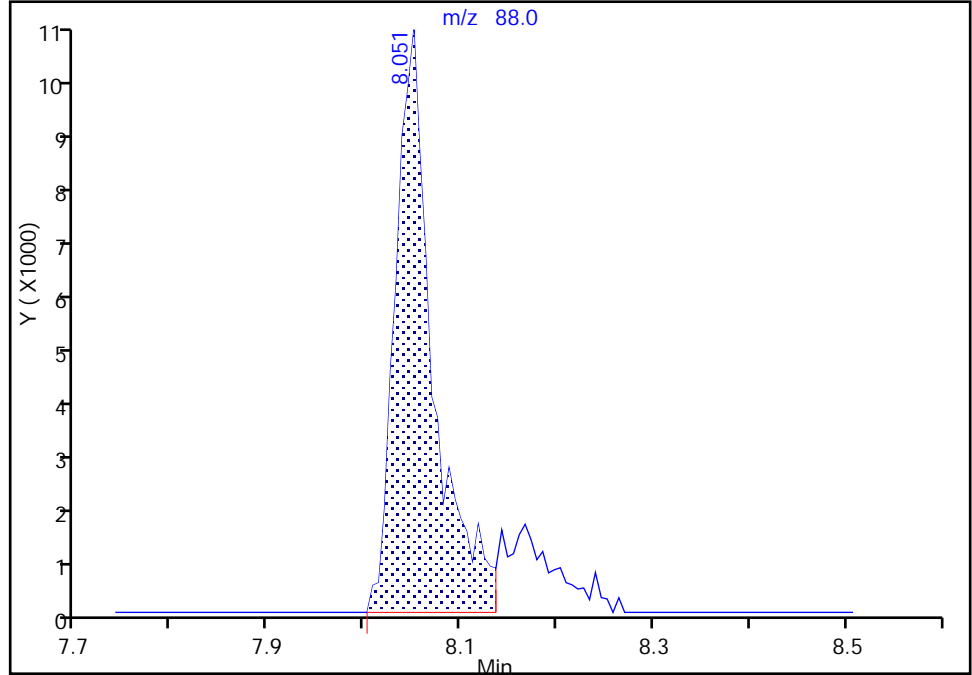
Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D04.D
Injection Date: 27-Jul-2017 01:39:30 Instrument ID: CHHP5
Lims ID: ICIS VSTD10
Client ID:
Operator ID: 034635 ALS Bottle#: 4 Worklist Smp#: 4
Purge Vol: 5.000 mL Dil. Factor: 1.0000
Method: MSVOA_LL_CHHP5 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm) Detector: MS SCAN

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

Signal: 1

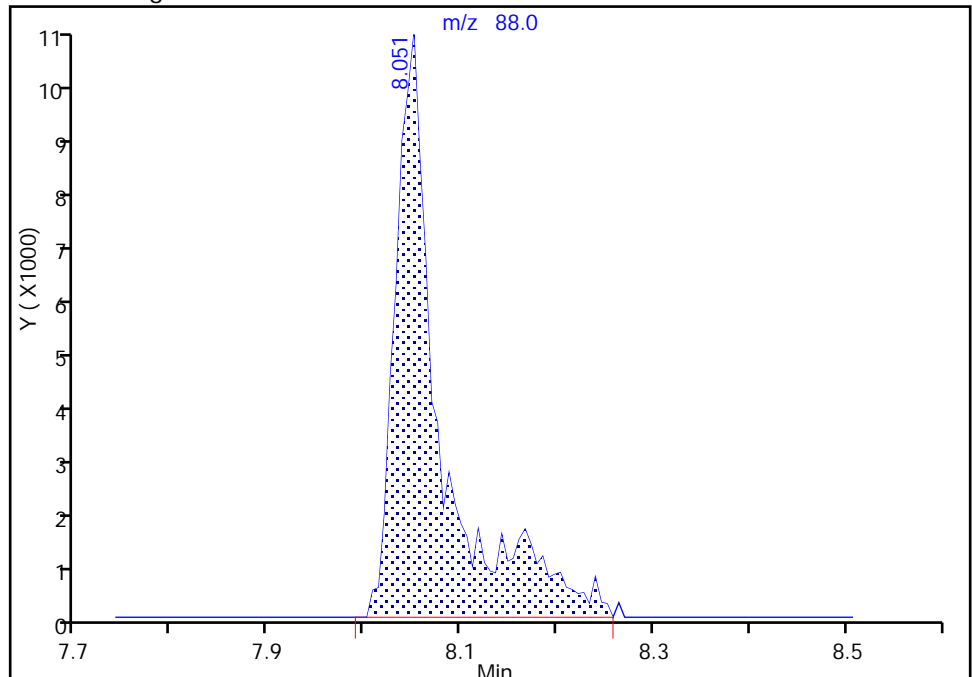
RT: 8.05
Area: 27736
Amount: 937.4398
Amount Units: ng

Processing Integration Results



RT: 8.05
Area: 33209
Amount: 1068.7953
Amount Units: ng

Manual Integration Results



Reviewer: bungardf, 27-Jul-2017 03:15:41
Audit Action: Manually Integrated

Audit Reason: Poor chromatography

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D05.D
 Lims ID: IC VSTD15
 Client ID:
 Sample Type: IC Calib Level: 4
 Inject. Date: 27-Jul-2017 02:02:30 ALS Bottle#: 5 Worklist Smp#: 5
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017756-005
 Misc. Info.: IC VSTD15
 Operator ID: 034635 Instrument ID: CHHP5
 Sublist: chrom-MSVOA_LL_CHHP5*sub12
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 28-Jul-2017 01:04:55 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK029

First Level Reviewer: bungardf

Date: 27-Jul-2017 03:16:54

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.323 | 4.323 | 0.000 | 0 | 240814 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.298 | 7.298 | 0.000 | 98 | 519897 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.406 | 10.406 | 0.000 | 84 | 132905 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.773 | 12.773 | 0.000 | 91 | 174376 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.574 | 6.574 | 0.000 | 93 | 193042 | 75.0 | 77.2 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.945 | 6.945 | 0.000 | 0 | 234269 | 75.0 | 76.8 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.946 | 8.946 | 0.000 | 92 | 780569 | 75.0 | 73.8 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.599 | 11.599 | 0.000 | 88 | 289432 | 75.0 | 75.8 | |
| 11 Dichlorodifluoromethane | 85 | 1.646 | 1.646 | 0.000 | 98 | 226899 | 75.0 | 75.1 | |
| 12 Chloromethane | 50 | 1.804 | 1.804 | 0.000 | 99 | 232300 | 75.0 | 76.5 | |
| 13 Vinyl chloride | 62 | 1.944 | 1.944 | 0.000 | 98 | 221295 | 75.0 | 71.8 | |
| 14 Butadiene | 39 | 1.969 | 1.969 | 0.000 | 96 | 204212 | 75.0 | 72.9 | |
| 15 Bromomethane | 94 | 2.254 | 2.254 | 0.000 | 90 | 112119 | 75.0 | 76.9 | |
| 16 Chloroethane | 64 | 2.419 | 2.419 | 0.000 | 99 | 128899 | 75.0 | 76.1 | |
| 17 Dichlorofluoromethane | 67 | 2.699 | 2.699 | 0.000 | 97 | 327021 | 75.0 | 76.3 | |
| 18 Trichlorofluoromethane | 101 | 2.741 | 2.741 | 0.000 | 94 | 283194 | 75.0 | 74.8 | |
| 20 Ethyl ether | 59 | 3.076 | 3.076 | 0.000 | 87 | 188662 | 75.0 | 76.6 | |
| 21 Acrolein | 56 | 3.252 | 3.252 | 0.000 | 99 | 115103 | 175.0 | 185.4 | |
| 22 1,1-Dichloroethene | 96 | 3.368 | 3.368 | 0.000 | 97 | 190985 | 75.0 | 75.0 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.441 | 3.441 | 0.000 | 92 | 206212 | 75.0 | 73.8 | |
| 24 Acetone | 43 | 3.477 | 3.477 | 0.000 | 100 | 227784 | 150.0 | 167.5 | |
| 25 Iodomethane | 142 | 3.562 | 3.562 | 0.000 | 96 | 304618 | 75.0 | 76.2 | |
| 26 Carbon disulfide | 76 | 3.648 | 3.648 | 0.000 | 98 | 403056 | 75.0 | 72.2 | |
| 28 3-Chloro-1-propene | 76 | 3.946 | 3.946 | 0.000 | 92 | 121734 | 75.0 | 74.0 | |
| 30 Methyl acetate | 43 | 3.976 | 3.976 | 0.000 | 97 | 419273 | 150.0 | 155.7 | |
| 31 Methylene Chloride | 84 | 4.165 | 4.165 | 0.000 | 87 | 242665 | 75.0 | 78.8 | |
| 32 2-Methyl-2-propanol | 59 | 4.451 | 4.451 | 0.000 | 95 | 204334 | 750.0 | 717.5 | |
| 33 Acrylonitrile | 53 | 4.554 | 4.554 | 0.000 | 98 | 1029651 | 750.0 | 786.5 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.584 | 4.584 | 0.000 | 97 | 222245 | 75.0 | 76.6 | |
| 35 Methyl tert-butyl ether | 73 | 4.603 | 4.603 | 0.000 | 95 | 613933 | 75.0 | 78.9 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 36 Hexane | 57 | 4.998 | 4.998 | 0.000 | 93 | 266987 | 75.0 | 71.7 | |
| 37 1,1-Dichloroethane | 63 | 5.217 | 5.217 | 0.000 | 96 | 379320 | 75.0 | 75.2 | |
| 38 Vinyl acetate | 43 | 5.272 | 5.272 | 0.000 | 97 | 400099 | 75.0 | 78.0 | |
| 44 2,2-Dichloropropane | 97 | 5.959 | 5.959 | 0.000 | 93 | 48893 | 75.0 | 76.2 | |
| 45 cis-1,2-Dichloroethene | 96 | 5.965 | 5.965 | 0.000 | 79 | 259385 | 75.0 | 78.2 | |
| 46 2-Butanone (MEK) | 43 | 5.978 | 5.978 | 0.000 | 98 | 321867 | 150.0 | 166.3 | |
| 49 Chlorobromomethane | 128 | 6.245 | 6.245 | 0.000 | 94 | 113290 | 75.0 | 76.8 | |
| 51 Tetrahydrofuran | 42 | 6.263 | 6.263 | 0.000 | 87 | 176266 | 150.0 | 156.4 | |
| 52 Chloroform | 83 | 6.391 | 6.391 | 0.000 | 93 | 389323 | 75.0 | 77.3 | |
| 53 1,1,1-Trichloroethane | 97 | 6.549 | 6.549 | 0.000 | 98 | 285488 | 75.0 | 74.9 | |
| 54 Cyclohexane | 56 | 6.622 | 6.622 | 0.000 | 88 | 345041 | 75.0 | 73.4 | |
| 56 Carbon tetrachloride | 117 | 6.726 | 6.726 | 0.000 | 97 | 238173 | 75.0 | 75.1 | |
| 55 1,1-Dichloropropene | 75 | 6.738 | 6.738 | 0.000 | 98 | 312373 | 75.0 | 75.9 | |
| 57 Isobutyl alcohol | 41 | 6.945 | 6.945 | 0.000 | 61 | 216532 | 1875.0 | 2093.1 | |
| 58 Benzene | 78 | 6.951 | 6.951 | 0.000 | 97 | 981851 | 75.0 | 77.7 | |
| 59 1,2-Dichloroethane | 62 | 7.030 | 7.030 | 0.000 | 98 | 292683 | 75.0 | 79.4 | |
| 62 n-Heptane | 43 | 7.316 | 7.316 | 0.000 | 88 | 214813 | 75.0 | 72.2 | |
| 64 Trichloroethene | 130 | 7.687 | 7.687 | 0.000 | 98 | 241861 | 75.0 | 76.0 | |
| 66 Methylcyclohexane | 83 | 7.918 | 7.918 | 0.000 | 86 | 358781 | 75.0 | 74.6 | |
| 67 1,2-Dichloropropane | 63 | 7.961 | 7.961 | 0.000 | 96 | 227133 | 75.0 | 77.2 | |
| 68 Dibromomethane | 93 | 8.046 | 8.046 | 0.000 | 95 | 135198 | 75.0 | 78.4 | |
| 70 1,4-Dioxane | 88 | 8.052 | 8.052 | 0.000 | 38 | 46920 | 1500.0 | 1567.5 | |
| 71 Dichlorobromomethane | 83 | 8.241 | 8.241 | 0.000 | 99 | 268080 | 75.0 | 79.2 | |
| 73 2-Chloroethyl vinyl ether | 63 | 8.545 | 8.545 | 0.000 | 92 | 343066 | 150.0 | 162.0 | |
| 74 cis-1,3-Dichloropropene | 75 | 8.685 | 8.685 | 0.000 | 96 | 320956 | 75.0 | 78.1 | |
| 75 4-Methyl-2-pentanone (MIBK) | 43 | 8.843 | 8.843 | 0.000 | 95 | 542662 | 150.0 | 159.2 | |
| 76 Toluene | 91 | 9.019 | 9.019 | 0.000 | 99 | 1000479 | 75.0 | 75.5 | |
| 77 trans-1,3-Dichloropropene | 75 | 9.269 | 9.269 | 0.000 | 93 | 278226 | 75.0 | 77.2 | |
| 78 Ethyl methacrylate | 69 | 9.330 | 9.330 | 0.000 | 87 | 352819 | 75.0 | 81.1 | |
| 79 1,1,2-Trichloroethane | 97 | 9.457 | 9.457 | 0.000 | 91 | 209928 | 75.0 | 76.0 | |
| 80 Tetrachloroethene | 164 | 9.530 | 9.530 | 0.000 | 97 | 184171 | 75.0 | 72.9 | |
| 81 1,3-Dichloropropane | 76 | 9.615 | 9.615 | 0.000 | 88 | 397870 | 75.0 | 78.0 | |
| 82 2-Hexanone | 43 | 9.682 | 9.682 | 0.000 | 93 | 419354 | 150.0 | 160.4 | |
| 84 Chlorodibromomethane | 129 | 9.834 | 9.834 | 0.000 | 91 | 181267 | 75.0 | 77.7 | |
| 85 Ethylene Dibromide | 107 | 9.944 | 9.944 | 0.000 | 97 | 223815 | 75.0 | 79.0 | |
| 86 3-Chlorobenzotrifluoride | 180 | 10.412 | 10.412 | 0.000 | 93 | 352260 | 75.0 | 77.1 | |
| 87 Chlorobenzene | 112 | 10.437 | 10.437 | 0.000 | 94 | 660247 | 75.0 | 76.5 | |
| 88 4-Chlorobenzotrifluoride | 180 | 10.498 | 10.498 | 0.000 | 96 | 327327 | 75.0 | 77.7 | |
| 89 1,1,1,2-Tetrachloroethane | 131 | 10.528 | 10.528 | 0.000 | 92 | 212641 | 75.0 | 77.5 | |
| 90 Ethylbenzene | 106 | 10.534 | 10.534 | 0.000 | 98 | 371119 | 75.0 | 77.1 | |
| 91 m-Xylene & p-Xylene | 106 | 10.668 | 10.668 | 0.000 | 0 | 452043 | 75.0 | 76.8 | |
| 92 o-Xylene | 106 | 11.051 | 11.051 | 0.000 | 95 | 440285 | 75.0 | 78.5 | |
| 93 Styrene | 104 | 11.069 | 11.069 | 0.000 | 94 | 745860 | 75.0 | 78.6 | |
| 94 Bromoform | 173 | 11.252 | 11.252 | 0.000 | 96 | 112077 | 75.0 | 77.3 | |
| 96 2-Chlorobenzotrifluoride | 180 | 11.325 | 11.325 | 0.000 | 97 | 348911 | 75.0 | 79.8 | |
| 97 Isopropylbenzene | 105 | 11.422 | 11.422 | 0.000 | 96 | 1080505 | 75.0 | 78.9 | |
| 100 Bromobenzene | 156 | 11.739 | 11.739 | 0.000 | 95 | 261052 | 75.0 | 77.1 | |
| 99 1,1,2,2-Tetrachloroethane | 83 | 11.745 | 11.745 | 0.000 | 95 | 316221 | 75.0 | 77.4 | |
| 102 trans-1,4-Dichloro-2-buten | 53 | 11.775 | 11.775 | 0.000 | 82 | 83561 | 75.0 | 81.9 | |
| 101 1,2,3-Trichloropropane | 110 | 11.793 | 11.793 | 0.000 | 85 | 109372 | 75.0 | 78.3 | |
| 103 N-Propylbenzene | 120 | 11.842 | 11.842 | 0.000 | 98 | 291693 | 75.0 | 75.4 | |
| 104 2-Chlorotoluene | 126 | 11.927 | 11.927 | 0.000 | 97 | 256066 | 75.0 | 76.6 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 105 3-Chlorotoluene | 126 | 11.994 | 11.994 | 0.000 | 97 | 289960 | 75.0 | 79.7 | |
| 106 1,3,5-Trimethylbenzene | 105 | 12.031 | 12.031 | 0.000 | 94 | 866332 | 75.0 | 78.3 | |
| 107 4-Chlorotoluene | 126 | 12.055 | 12.055 | 0.000 | 96 | 269544 | 75.0 | 74.7 | |
| 108 tert-Butylbenzene | 119 | 12.347 | 12.347 | 0.000 | 93 | 721573 | 75.0 | 78.0 | |
| 110 1,2,4-Trimethylbenzene | 105 | 12.408 | 12.408 | 0.000 | 97 | 884487 | 75.0 | 78.6 | |
| 111 1,2-dichloro-4-(trifluorom | 214 | 12.456 | 12.456 | 0.000 | 97 | 219982 | 75.0 | 78.1 | |
| 112 sec-Butylbenzene | 105 | 12.572 | 12.572 | 0.000 | 94 | 993968 | 75.0 | 77.0 | |
| 113 1,3-Dichlorobenzene | 146 | 12.688 | 12.688 | 0.000 | 97 | 462404 | 75.0 | 76.5 | |
| 114 4-Isopropyltoluene | 119 | 12.730 | 12.730 | 0.000 | 96 | 837492 | 75.0 | 77.9 | |
| 115 1,4-Dichlorobenzene | 146 | 12.797 | 12.797 | 0.000 | 96 | 474362 | 75.0 | 76.4 | |
| 116 2,4-Dichloro-1-(trifluorom | 214 | 12.828 | 12.828 | 0.000 | 94 | 206368 | 75.0 | 78.6 | |
| 118 2,5-Dichlorobenzotrifluori | 214 | 12.870 | 12.870 | 0.000 | 0 | 217211 | 75.0 | 76.6 | |
| 120 n-Butylbenzene | 91 | 13.150 | 13.150 | 0.000 | 98 | 671190 | 75.0 | 76.5 | |
| 121 1,2-Dichlorobenzene | 146 | 13.156 | 13.156 | 0.000 | 98 | 437966 | 75.0 | 76.0 | |
| 122 1,2-Dibromo-3-Chloropropan | 75 | 13.971 | 13.971 | 0.000 | 83 | 47827 | 75.0 | 74.7 | |
| 123 2,4- & 2,5- & 2,6- Dichlor | 125 | 14.117 | 14.117 | 0.000 | 0 | 889724 | 225.0 | 243.4 | |
| 125 2,3- & 3,4- Dichlorotoluen | 125 | 14.555 | 14.555 | 0.000 | 0 | 620870 | 150.0 | 164.2 | |
| 126 1,2,4-Trichlorobenzene | 180 | 14.829 | 14.829 | 0.000 | 94 | 200638 | 75.0 | 76.1 | |
| 127 Hexachlorobutadiene | 225 | 14.993 | 14.993 | 0.000 | 98 | 73984 | 75.0 | 76.7 | |
| 128 Naphthalene | 128 | 15.103 | 15.103 | 0.000 | 97 | 733996 | 75.0 | 81.7 | |
| 129 1,2,3-Trichlorobenzene | 180 | 15.346 | 15.346 | 0.000 | 96 | 184932 | 75.0 | 76.8 | |
| 131 2,4,5-Trichlorotoluene | 159 | 16.198 | 16.198 | 0.000 | 0 | 91488 | 75.0 | 79.9 | |
| 130 2,3,6-Trichlorotoluene | 159 | 16.307 | 16.307 | 0.000 | 98 | 89402 | 75.0 | 83.9 | |
| 149 3,4-Dichlorotoluene | 1 | | 0.000 | | | | ND | ND | |
| S 134 1,2-Dichloroethene, Total | 96 | | | | 0 | | 150.0 | 154.8 | |
| S 133 Xylenes, Total | 106 | | | | 0 | | 150.0 | 155.3 | |
| S 135 1,3-Dichloropropene, Total | 1 | | | | 0 | | 150.0 | 155.2 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Reagents:

| | | |
|---------------------|--------------------|-----------|
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 3.00 | Units: uL |
| VOA8260VOAPRI_00263 | Amount Added: 3.00 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 3.00 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 7.00 | Units: uL |
| voaWVA1stRest_00017 | Amount Added: 3.00 | Units: uL |
| voaWEEmix1stR_00009 | Amount Added: 3.00 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 3.00 | Units: uL |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D05.D

Injection Date: 27-Jul-2017 02:02:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: IC VSTD15

Worklist Smp#: 5

Client ID:

Purge Vol: 5.000 mL

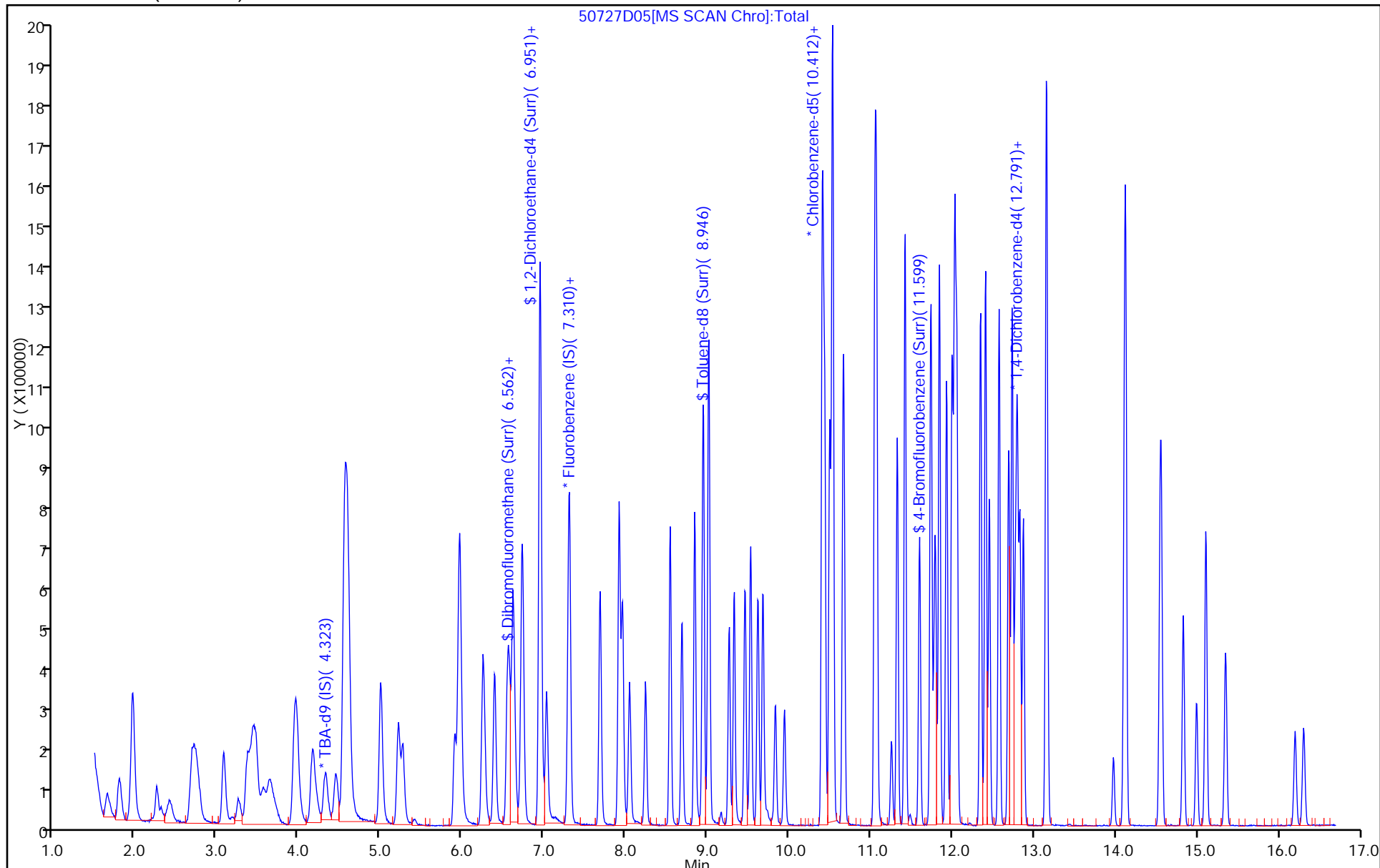
Dil. Factor: 1.0000

ALS Bottle#: 5

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D06.D
 Lims ID: IC VSTD20
 Client ID:
 Sample Type: IC Calib Level: 5
 Inject. Date: 27-Jul-2017 02:26:30 ALS Bottle#: 6 Worklist Smp#: 6
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017756-006
 Misc. Info.: IC VSTD20
 Operator ID: 034635 Instrument ID: CHHP5
 Sublist: chrom-MSVOA_LL_CHHP5*sub12
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 28-Jul-2017 01:04:58 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK029

First Level Reviewer: bungardf

Date: 27-Jul-2017 03:06:29

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.316 | 4.323 | -0.007 | 0 | 252187 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.297 | 7.298 | -0.001 | 98 | 520193 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.406 | 10.406 | 0.000 | 85 | 132635 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.772 | 12.773 | -0.001 | 95 | 171832 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.573 | 6.574 | -0.001 | 93 | 257355 | 100.0 | 102.8 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.944 | 6.945 | -0.001 | 0 | 307676 | 100.0 | 100.8 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.945 | 8.946 | -0.001 | 92 | 1040595 | 100.0 | 98.6 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.598 | 11.599 | -0.001 | 87 | 390879 | 100.0 | 102.5 | |
| 11 Dichlorodifluoromethane | 85 | 1.651 | 1.646 | 0.005 | 99 | 286388 | 100.0 | 94.7 | |
| 12 Chloromethane | 50 | 1.797 | 1.804 | -0.007 | 99 | 302276 | 100.0 | 99.4 | |
| 13 Vinyl chloride | 62 | 1.949 | 1.944 | 0.005 | 98 | 291558 | 100.0 | 94.5 | |
| 14 Butadiene | 39 | 1.962 | 1.969 | -0.006 | 92 | 260580 | 100.0 | 93.0 | |
| 15 Bromomethane | 94 | 2.260 | 2.254 | 0.006 | 90 | 161865 | 100.0 | 111.0 | |
| 16 Chloroethane | 64 | 2.412 | 2.419 | -0.007 | 99 | 172552 | 100.0 | 101.8 | |
| 17 Dichlorofluoromethane | 67 | 2.710 | 2.699 | 0.011 | 97 | 436022 | 100.0 | 101.7 | |
| 18 Trichlorofluoromethane | 101 | 2.734 | 2.741 | -0.007 | 96 | 371684 | 100.0 | 98.1 | |
| 20 Ethyl ether | 59 | 3.081 | 3.076 | 0.005 | 89 | 262150 | 100.0 | 106.3 | |
| 21 Acrolein | 56 | 3.264 | 3.252 | 0.012 | 99 | 130923 | 200.0 | 210.7 | |
| 22 1,1-Dichloroethene | 96 | 3.373 | 3.368 | 0.005 | 98 | 247279 | 100.0 | 97.1 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.446 | 3.441 | 0.005 | 93 | 263603 | 100.0 | 94.3 | |
| 24 Acetone | 43 | 3.476 | 3.477 | -0.001 | 100 | 316026 | 200.0 | 232.3 | |
| 25 Iodomethane | 142 | 3.562 | 3.562 | 0.000 | 98 | 408622 | 100.0 | 102.2 | |
| 26 Carbon disulfide | 76 | 3.647 | 3.648 | -0.001 | 99 | 561008 | 100.0 | 100.4 | |
| 28 3-Chloro-1-propene | 76 | 3.951 | 3.946 | 0.005 | 92 | 164305 | 100.0 | 99.8 | |
| 30 Methyl acetate | 43 | 3.969 | 3.976 | -0.007 | 97 | 558912 | 200.0 | 207.5 | |
| 31 Methylene Chloride | 84 | 4.164 | 4.165 | -0.001 | 93 | 323324 | 100.0 | 106.0 | |
| 32 2-Methyl-2-propanol | 59 | 4.444 | 4.451 | -0.007 | 94 | 283777 | 1000.0 | 951.5 | |
| 33 Acrylonitrile | 53 | 4.553 | 4.554 | -0.001 | 99 | 1387354 | 1000.0 | 1059.2 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.584 | 4.584 | 0.000 | 98 | 296608 | 100.0 | 102.2 | |
| 35 Methyl tert-butyl ether | 73 | 4.602 | 4.603 | -0.001 | 95 | 822838 | 100.0 | 105.8 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 36 Hexane | 57 | 5.003 | 4.998 | 0.005 | 92 | 337300 | 100.0 | 90.6 | |
| 37 1,1-Dichloroethane | 63 | 5.210 | 5.217 | -0.007 | 96 | 510811 | 100.0 | 101.2 | |
| 38 Vinyl acetate | 43 | 5.265 | 5.272 | -0.007 | 97 | 532250 | 100.0 | 103.7 | |
| 44 2,2-Dichloropropane | 97 | 5.959 | 5.959 | -0.001 | 57 | 65750 | 100.0 | 102.4 | |
| 45 cis-1,2-Dichloroethene | 96 | 5.965 | 5.965 | 0.000 | 79 | 347303 | 100.0 | 104.6 | |
| 46 2-Butanone (MEK) | 43 | 5.983 | 5.978 | 0.005 | 98 | 426755 | 200.0 | 220.4 | |
| 49 Chlorobromomethane | 128 | 6.251 | 6.245 | 0.005 | 94 | 155416 | 100.0 | 105.4 | |
| 51 Tetrahydrofuran | 42 | 6.263 | 6.263 | 0.000 | 86 | 224432 | 200.0 | 199.0 | |
| 52 Chloroform | 83 | 6.390 | 6.391 | -0.001 | 92 | 517765 | 100.0 | 102.8 | |
| 53 1,1,1-Trichloroethane | 97 | 6.555 | 6.549 | 0.006 | 98 | 383868 | 100.0 | 100.7 | |
| 54 Cyclohexane | 56 | 6.622 | 6.622 | 0.000 | 89 | 446560 | 100.0 | 94.9 | |
| 56 Carbon tetrachloride | 117 | 6.725 | 6.726 | -0.001 | 96 | 317033 | 100.0 | 99.9 | |
| 55 1,1-Dichloropropene | 75 | 6.737 | 6.738 | -0.001 | 98 | 408627 | 100.0 | 99.2 | |
| 58 Benzene | 78 | 6.956 | 6.951 | 0.005 | 97 | 1307056 | 100.0 | 103.3 | |
| 57 Isobutyl alcohol | 41 | 6.944 | 6.945 | -0.001 | 91 | 290317 | 2500.0 | 2804.8 | |
| 59 1,2-Dichloroethane | 62 | 7.029 | 7.030 | -0.001 | 97 | 385206 | 100.0 | 104.5 | |
| 62 n-Heptane | 43 | 7.315 | 7.316 | -0.001 | 89 | 279216 | 100.0 | 93.8 | |
| 64 Trichloroethene | 130 | 7.686 | 7.687 | -0.001 | 98 | 329499 | 100.0 | 103.5 | |
| 66 Methylcyclohexane | 83 | 7.917 | 7.918 | -0.001 | 87 | 467268 | 100.0 | 97.1 | |
| 67 1,2-Dichloropropane | 63 | 7.960 | 7.961 | -0.001 | 96 | 309491 | 100.0 | 105.1 | |
| 68 Dibromomethane | 93 | 8.051 | 8.046 | 0.005 | 96 | 184529 | 100.0 | 106.9 | |
| 70 1,4-Dioxane | 88 | 8.045 | 8.052 | -0.007 | 39 | 65688 | 2000.0 | 2193.3 | |
| 71 Dichlorobromomethane | 83 | 8.240 | 8.241 | -0.001 | 99 | 366097 | 100.0 | 108.1 | |
| 73 2-Chloroethyl vinyl ether | 63 | 8.544 | 8.545 | -0.001 | 92 | 467677 | 200.0 | 220.7 | |
| 74 cis-1,3-Dichloropropene | 75 | 8.684 | 8.685 | -0.001 | 96 | 447138 | 100.0 | 108.7 | |
| 75 4-Methyl-2-pentanone (MIBK) | 43 | 8.842 | 8.843 | -0.001 | 95 | 738839 | 200.0 | 217.2 | |
| 76 Toluene | 91 | 9.018 | 9.019 | -0.001 | 99 | 1332783 | 100.0 | 100.8 | |
| 77 trans-1,3-Dichloropropene | 75 | 9.268 | 9.269 | -0.001 | 92 | 396221 | 100.0 | 110.1 | |
| 78 Ethyl methacrylate | 69 | 9.329 | 9.330 | -0.001 | 87 | 483364 | 100.0 | 111.4 | |
| 79 1,1,2-Trichloroethane | 97 | 9.456 | 9.457 | -0.001 | 90 | 283688 | 100.0 | 103.0 | |
| 80 Tetrachloroethene | 164 | 9.529 | 9.530 | -0.001 | 97 | 244346 | 100.0 | 96.9 | |
| 81 1,3-Dichloropropane | 76 | 9.615 | 9.615 | 0.000 | 89 | 518120 | 100.0 | 101.7 | |
| 82 2-Hexanone | 43 | 9.676 | 9.682 | -0.006 | 94 | 581383 | 200.0 | 222.8 | |
| 84 Chlorodibromomethane | 129 | 9.834 | 9.834 | 0.000 | 90 | 254603 | 100.0 | 109.3 | |
| 85 Ethylene Dibromide | 107 | 9.943 | 9.944 | -0.001 | 99 | 294438 | 100.0 | 104.2 | |
| 86 3-Chlorobenzotrifluoride | 180 | 10.412 | 10.412 | 0.000 | 94 | 461082 | 100.0 | 101.2 | |
| 87 Chlorobenzene | 112 | 10.436 | 10.437 | -0.001 | 95 | 877804 | 100.0 | 102.0 | |
| 88 4-Chlorobenzotrifluoride | 180 | 10.497 | 10.498 | -0.001 | 96 | 420704 | 100.0 | 100.0 | |
| 90 Ethylbenzene | 106 | 10.533 | 10.534 | -0.001 | 98 | 499116 | 100.0 | 103.8 | |
| 89 1,1,1,2-Tetrachloroethane | 131 | 10.527 | 10.528 | -0.001 | 92 | 289044 | 100.0 | 105.6 | |
| 91 m-Xylene & p-Xylene | 106 | 10.667 | 10.668 | -0.001 | 0 | 610286 | 100.0 | 103.9 | |
| 92 o-Xylene | 106 | 11.050 | 11.051 | -0.001 | 95 | 592117 | 100.0 | 105.8 | |
| 93 Styrene | 104 | 11.075 | 11.069 | 0.006 | 94 | 1002147 | 100.0 | 105.8 | |
| 94 Bromoform | 173 | 11.251 | 11.252 | -0.001 | 97 | 157509 | 100.0 | 108.8 | |
| 96 2-Chlorobenzotrifluoride | 180 | 11.324 | 11.325 | -0.001 | 97 | 454842 | 100.0 | 104.3 | |
| 97 Isopropylbenzene | 105 | 11.421 | 11.422 | -0.001 | 96 | 1415676 | 100.0 | 103.6 | |
| 99 1,1,2,2-Tetrachloroethane | 83 | 11.738 | 11.745 | -0.007 | 95 | 412534 | 100.0 | 101.1 | |
| 100 Bromobenzene | 156 | 11.738 | 11.739 | -0.001 | 95 | 348475 | 100.0 | 104.5 | |
| 102 trans-1,4-Dichloro-2-buten | 53 | 11.774 | 11.775 | -0.001 | 82 | 104361 | 100.0 | 103.8 | |
| 101 1,2,3-Trichloropropane | 110 | 11.793 | 11.793 | 0.000 | 85 | 144469 | 100.0 | 105.0 | |
| 103 N-Propylbenzene | 120 | 11.841 | 11.842 | -0.001 | 98 | 387234 | 100.0 | 101.6 | |
| 104 2-Chlorotoluene | 126 | 11.926 | 11.927 | -0.001 | 97 | 344800 | 100.0 | 104.7 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 105 3-Chlorotoluene | 126 | 11.999 | 11.994 | 0.005 | 96 | 381649 | 100.0 | 106.5 | |
| 106 1,3,5-Trimethylbenzene | 105 | 12.030 | 12.031 | -0.001 | 94 | 1140888 | 100.0 | 104.6 | |
| 107 4-Chlorotoluene | 126 | 12.054 | 12.055 | -0.001 | 96 | 369832 | 100.0 | 104.0 | |
| 108 tert-Butylbenzene | 119 | 12.346 | 12.347 | -0.001 | 93 | 931884 | 100.0 | 102.2 | |
| 110 1,2,4-Trimethylbenzene | 105 | 12.407 | 12.408 | -0.001 | 97 | 1156912 | 100.0 | 104.4 | |
| 111 1,2-dichloro-4-(trifluorom | 214 | 12.456 | 12.456 | 0.000 | 97 | 277157 | 100.0 | 99.8 | |
| 112 sec-Butylbenzene | 105 | 12.571 | 12.572 | -0.001 | 94 | 1298722 | 100.0 | 102.1 | |
| 113 1,3-Dichlorobenzene | 146 | 12.687 | 12.688 | -0.001 | 97 | 613101 | 100.0 | 102.9 | |
| 114 4-Isopropyltoluene | 119 | 12.729 | 12.730 | -0.001 | 96 | 1086140 | 100.0 | 102.5 | |
| 115 1,4-Dichlorobenzene | 146 | 12.796 | 12.797 | -0.001 | 94 | 622850 | 100.0 | 101.8 | |
| 116 2,4-Dichloro-1-(trifluorom | 214 | 12.827 | 12.828 | -0.001 | 96 | 267418 | 100.0 | 103.4 | |
| 118 2,5-Dichlorobenzotrifluori | 214 | 12.869 | 12.870 | -0.001 | 0 | 279514 | 100.0 | 100.1 | |
| 120 n-Butylbenzene | 91 | 13.149 | 13.150 | -0.001 | 97 | 885288 | 100.0 | 102.4 | |
| 121 1,2-Dichlorobenzene | 146 | 13.155 | 13.156 | -0.001 | 97 | 577962 | 100.0 | 101.8 | |
| 122 1,2-Dibromo-3-Chloropropan | 75 | 13.970 | 13.971 | -0.001 | 85 | 68470 | 100.0 | 108.6 | |
| 123 2,4- & 2,5- & 2,6- Dichlor | 125 | 14.116 | 14.117 | -0.001 | 0 | 1151252 | 300.0 | 319.5 | |
| 125 2,3- & 3,4- Dichlorotoluen | 125 | 14.548 | 14.555 | -0.007 | 0 | 814032 | 200.0 | 218.5 | |
| 126 1,2,4-Trichlorobenzene | 180 | 14.828 | 14.829 | -0.001 | 95 | 266863 | 100.0 | 102.7 | |
| 127 Hexachlorobutadiene | 225 | 14.992 | 14.993 | -0.001 | 97 | 94134 | 100.0 | 99.0 | |
| 128 Naphthalene | 128 | 15.102 | 15.103 | -0.001 | 97 | 990398 | 100.0 | 111.9 | |
| 129 1,2,3-Trichlorobenzene | 180 | 15.345 | 15.346 | -0.001 | 97 | 247660 | 100.0 | 104.3 | |
| 131 2,4,5-Trichlorotoluene | 159 | 16.197 | 16.198 | -0.001 | 0 | 122498 | 100.0 | 108.5 | |
| 130 2,3,6-Trichlorotoluene | 159 | 16.306 | 16.307 | -0.001 | 96 | 115009 | 100.0 | 109.5 | |
| 149 3,4-Dichlorotoluene | 1 | | 0.000 | | | | ND | ND | |
| S 133 Xylenes, Total | 106 | | | | 0 | | 200.0 | 209.7 | |
| S 134 1,2-Dichloroethene, Total | 96 | | | | 0 | | 200.0 | 206.9 | |
| S 135 1,3-Dichloropropene, Total | 1 | | | | 0 | | 200.0 | 218.8 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Reagents:

| | | |
|---------------------|--------------------|-----------|
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 4.00 | Units: uL |
| VOA8260VOAPRI_00263 | Amount Added: 4.00 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 4.00 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 8.00 | Units: uL |
| voaWVA1stRest_00017 | Amount Added: 4.00 | Units: uL |
| voaWEEmix1stR_00009 | Amount Added: 4.00 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 4.00 | Units: uL |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D06.D

Injection Date: 27-Jul-2017 02:26:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: IC VSTD20

Worklist Smp#: 6

Client ID:

Purge Vol: 5.000 mL

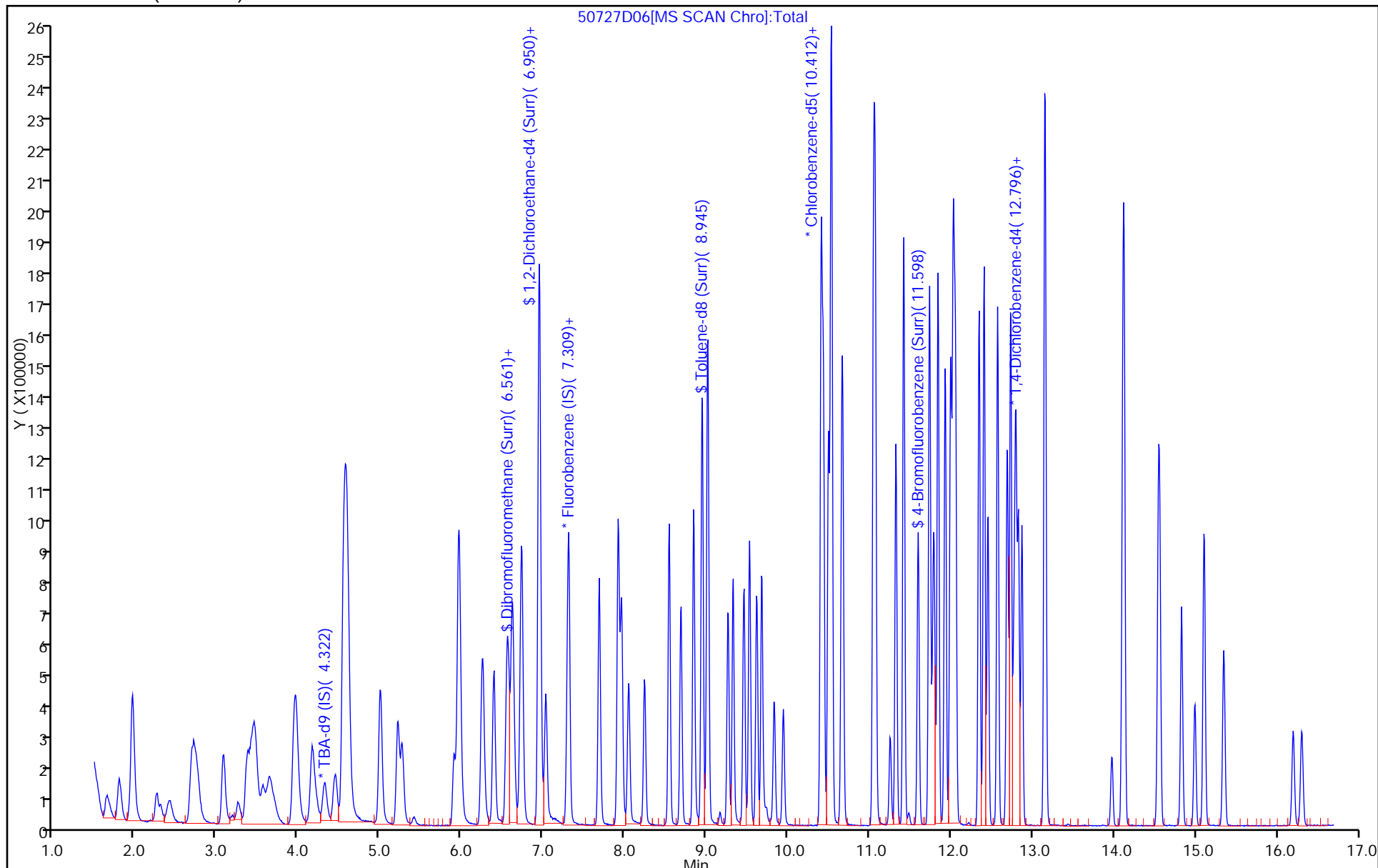
Dil. Factor: 1.0000

ALS Bottle#: 6

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D08.D
 Lims ID: IC VSTD40
 Client ID:
 Sample Type: IC Calib Level: 7
 Inject. Date: 27-Jul-2017 03:13:30 ALS Bottle#: 8 Worklist Smp#: 8
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017756-008
 Misc. Info.: IC VSTD40
 Operator ID: 034635 Instrument ID: CHHP5
 Sublist: chrom-MSVOA_LL_CHHP5*sub12
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 28-Jul-2017 01:05:02 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK029

First Level Reviewer: bungardf

Date: 27-Jul-2017 03:34:06

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.337 | 4.323 | 0.013 | 0 | 252542 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.299 | 7.298 | 0.001 | 99 | 561296 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.408 | 10.406 | 0.002 | 56 | 150914 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.768 | 12.773 | -0.005 | 90 | 189484 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.575 | 6.574 | 0.001 | 94 | 522323 | 200.0 | 193.4 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.946 | 6.945 | 0.001 | 0 | 628942 | 200.0 | 190.9 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.948 | 8.946 | 0.002 | 92 | 2000995 | 200.0 | 166.6 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.594 | 11.599 | -0.005 | 92 | 793129 | 200.0 | 182.8 | |
| 11 Dichlorodifluoromethane | 85 | 1.654 | 1.646 | 0.008 | 99 | 569791 | 200.0 | 174.6 | |
| 12 Chloromethane | 50 | 1.812 | 1.804 | 0.008 | 99 | 580608 | 200.0 | 177.0 | |
| 13 Vinyl chloride | 62 | 1.958 | 1.944 | 0.014 | 97 | 577090 | 200.0 | 173.4 | |
| 14 Butadiene | 39 | 1.970 | 1.969 | 0.002 | 94 | 512032 | 200.0 | 169.3 | |
| 15 Bromomethane | 94 | 2.268 | 2.254 | 0.014 | 91 | 289712 | 200.0 | 184.1 | |
| 16 Chloroethane | 64 | 2.426 | 2.419 | 0.007 | 99 | 322589 | 200.0 | 176.3 | |
| 17 Dichlorofluoromethane | 67 | 2.706 | 2.699 | 0.007 | 97 | 819020 | 200.0 | 177.0 | |
| 18 Trichlorofluoromethane | 101 | 2.761 | 2.741 | 0.020 | 97 | 710415 | 200.0 | 173.7 | |
| 20 Ethyl ether | 59 | 3.077 | 3.076 | 0.001 | 88 | 510033 | 200.0 | 191.7 | |
| 21 Acrolein | 56 | 3.260 | 3.252 | 0.008 | 100 | 179414 | 250.0 | 267.6 | |
| 22 1,1-Dichloroethene | 96 | 3.369 | 3.368 | 0.001 | 96 | 489503 | 200.0 | 178.1 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.442 | 3.441 | 0.001 | 93 | 534815 | 200.0 | 177.3 | |
| 24 Acetone | 43 | 3.485 | 3.477 | 0.008 | 100 | 522287 | 400.0 | 355.8 | |
| 25 Iodomethane | 142 | 3.576 | 3.562 | 0.014 | 98 | 834240 | 200.0 | 193.3 | |
| 26 Carbon disulfide | 76 | 3.649 | 3.648 | 0.001 | 99 | 1211678 | 200.0 | 200.9 | |
| 28 3-Chloro-1-propene | 76 | 3.947 | 3.946 | 0.001 | 92 | 366340 | 200.0 | 206.3 | |
| 30 Methyl acetate | 43 | 3.978 | 3.976 | 0.002 | 97 | 1173609 | 400.0 | 403.7 | |
| 31 Methylene Chloride | 84 | 4.166 | 4.165 | 0.001 | 88 | 653341 | 200.0 | 201.5 | |
| 32 2-Methyl-2-propanol | 59 | 4.464 | 4.451 | 0.013 | 93 | 519054 | 2000.0 | 1737.9 | |
| 33 Acrylonitrile | 53 | 4.562 | 4.554 | 0.008 | 99 | 2794353 | 2000.0 | 1977.2 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.580 | 4.584 | -0.004 | 97 | 571864 | 200.0 | 182.6 | |
| 35 Methyl tert-butyl ether | 73 | 4.604 | 4.603 | 0.001 | 95 | 1751345 | 200.0 | 208.6 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 36 Hexane | 57 | 5.000 | 4.998 | 0.002 | 92 | 708650 | 200.0 | 176.3 | |
| 37 1,1-Dichloroethane | 63 | 5.219 | 5.217 | 0.002 | 96 | 1041269 | 200.0 | 191.3 | |
| 38 Vinyl acetate | 43 | 5.273 | 5.272 | 0.001 | 97 | 1200052 | 200.0 | 216.8 | |
| 44 2,2-Dichloropropane | 97 | 5.961 | 5.959 | 0.002 | 88 | 125406 | 200.0 | 180.9 | |
| 45 cis-1,2-Dichloroethene | 96 | 5.967 | 5.965 | 0.002 | 80 | 687049 | 200.0 | 191.8 | |
| 46 2-Butanone (MEK) | 43 | 5.979 | 5.978 | 0.001 | 98 | 795793 | 400.0 | 380.9 | |
| 49 Chlorobromomethane | 128 | 6.247 | 6.245 | 0.002 | 94 | 313977 | 200.0 | 197.3 | |
| 51 Tetrahydrofuran | 42 | 6.265 | 6.263 | 0.002 | 86 | 488432 | 400.0 | 401.4 | |
| 52 Chloroform | 83 | 6.393 | 6.391 | 0.002 | 93 | 1037446 | 200.0 | 190.8 | |
| 53 1,1,1-Trichloroethane | 97 | 6.551 | 6.549 | 0.002 | 98 | 777880 | 200.0 | 189.0 | |
| 54 Cyclohexane | 56 | 6.618 | 6.622 | -0.004 | 90 | 922281 | 200.0 | 181.6 | |
| 56 Carbon tetrachloride | 117 | 6.721 | 6.726 | -0.005 | 97 | 646700 | 200.0 | 188.8 | |
| 55 1,1-Dichloropropene | 75 | 6.739 | 6.738 | 0.001 | 97 | 825970 | 200.0 | 185.8 | |
| 57 Isobutyl alcohol | 41 | 6.946 | 6.945 | 0.001 | 51 | 587752 | 5000.0 | 5262.5 | |
| 58 Benzene | 78 | 6.952 | 6.951 | 0.001 | 97 | 2487856 | 200.0 | 182.3 | |
| 59 1,2-Dichloroethane | 62 | 7.031 | 7.030 | 0.001 | 97 | 767974 | 200.0 | 193.0 | |
| 62 n-Heptane | 43 | 7.311 | 7.316 | -0.005 | 87 | 573064 | 200.0 | 178.3 | |
| 64 Trichloroethene | 130 | 7.682 | 7.687 | -0.005 | 98 | 647404 | 200.0 | 188.5 | |
| 66 Methylcyclohexane | 83 | 7.920 | 7.918 | 0.002 | 87 | 950167 | 200.0 | 183.0 | |
| 67 1,2-Dichloropropane | 63 | 7.962 | 7.961 | 0.001 | 96 | 624637 | 200.0 | 196.5 | |
| 68 Dibromomethane | 93 | 8.047 | 8.046 | 0.001 | 95 | 374289 | 200.0 | 201.0 | |
| 70 1,4-Dioxane | 88 | 8.041 | 8.052 | -0.011 | 39 | 135844 | 4000.0 | 4203.6 | |
| 71 Dichlorobromomethane | 83 | 8.242 | 8.241 | 0.001 | 99 | 752352 | 200.0 | 205.8 | |
| 73 2-Chloroethyl vinyl ether | 63 | 8.546 | 8.545 | 0.001 | 93 | 977190 | 400.0 | 427.3 | |
| 74 cis-1,3-Dichloropropene | 75 | 8.686 | 8.685 | 0.001 | 96 | 933591 | 200.0 | 210.3 | |
| 75 4-Methyl-2-pentanone (MIBK) | 43 | 8.844 | 8.843 | 0.001 | 95 | 1476808 | 400.0 | 381.5 | |
| 76 Toluene | 91 | 9.015 | 9.019 | -0.004 | 98 | 2540251 | 200.0 | 168.8 | |
| 77 trans-1,3-Dichloropropene | 75 | 9.264 | 9.269 | -0.005 | 92 | 850338 | 200.0 | 207.7 | |
| 78 Ethyl methacrylate | 69 | 9.325 | 9.330 | -0.005 | 88 | 1001550 | 200.0 | 202.8 | |
| 79 1,1,2-Trichloroethane | 97 | 9.459 | 9.457 | 0.002 | 91 | 569083 | 200.0 | 181.5 | |
| 80 Tetrachloroethene | 164 | 9.532 | 9.530 | 0.002 | 97 | 486427 | 200.0 | 169.5 | |
| 81 1,3-Dichloropropane | 76 | 9.617 | 9.615 | 0.002 | 89 | 1058308 | 200.0 | 182.6 | |
| 82 2-Hexanone | 43 | 9.678 | 9.682 | -0.004 | 93 | 1109580 | 400.0 | 373.7 | |
| 84 Chlorodibromomethane | 129 | 9.830 | 9.834 | -0.004 | 89 | 540065 | 200.0 | 203.8 | |
| 85 Ethylene Dibromide | 107 | 9.945 | 9.944 | 0.001 | 98 | 607203 | 200.0 | 188.9 | |
| 86 3-Chlorobenzotrifluoride | 180 | 10.408 | 10.412 | -0.004 | 93 | 869071 | 200.0 | 167.6 | |
| 87 Chlorobenzene | 112 | 10.432 | 10.437 | -0.005 | 93 | 1704167 | 200.0 | 174.0 | |
| 88 4-Chlorobenzotrifluoride | 180 | 10.499 | 10.498 | 0.001 | 96 | 810848 | 200.0 | 169.4 | |
| 89 1,1,1,2-Tetrachloroethane | 131 | 10.529 | 10.528 | 0.001 | 94 | 590452 | 200.0 | 189.5 | |
| 90 Ethylbenzene | 106 | 10.536 | 10.534 | 0.002 | 98 | 972676 | 200.0 | 177.9 | |
| 91 m-Xylene & p-Xylene | 106 | 10.669 | 10.668 | 0.001 | 0 | 1217768 | 200.0 | 182.2 | |
| 92 o-Xylene | 106 | 11.053 | 11.051 | 0.002 | 95 | 1159372 | 200.0 | 182.1 | |
| 93 Styrene | 104 | 11.071 | 11.069 | 0.002 | 94 | 1967591 | 200.0 | 182.6 | |
| 94 Bromoform | 173 | 11.253 | 11.252 | 0.001 | 96 | 350923 | 200.0 | 213.1 | |
| 96 2-Chlorobenzotrifluoride | 180 | 11.326 | 11.325 | 0.001 | 96 | 875687 | 200.0 | 176.5 | |
| 97 Isopropylbenzene | 105 | 11.418 | 11.422 | -0.004 | 96 | 2665903 | 200.0 | 171.5 | |
| 100 Bromobenzene | 156 | 11.734 | 11.739 | -0.005 | 95 | 711710 | 200.0 | 193.5 | |
| 99 1,1,2,2-Tetrachloroethane | 83 | 11.740 | 11.745 | -0.005 | 93 | 870164 | 200.0 | 187.5 | |
| 102 trans-1,4-Dichloro-2-buten | 53 | 11.777 | 11.775 | 0.002 | 85 | 225821 | 200.0 | 203.6 | |
| 101 1,2,3-Trichloropropane | 110 | 11.795 | 11.793 | 0.002 | 85 | 299299 | 200.0 | 197.2 | |
| 103 N-Propylbenzene | 120 | 11.844 | 11.842 | 0.002 | 97 | 774184 | 200.0 | 184.2 | |
| 104 2-Chlorotoluene | 126 | 11.929 | 11.927 | 0.002 | 97 | 700158 | 200.0 | 192.7 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 105 3-Chlorotoluene | 126 | 11.996 | 11.994 | 0.002 | 96 | 742625 | 200.0 | 187.9 | |
| 106 1,3,5-Trimethylbenzene | 105 | 12.026 | 12.031 | -0.005 | 94 | 2188229 | 200.0 | 182.0 | |
| 107 4-Chlorotoluene | 126 | 12.056 | 12.055 | 0.001 | 95 | 738280 | 200.0 | 188.2 | |
| 108 tert-Butylbenzene | 119 | 12.342 | 12.347 | -0.005 | 93 | 1809964 | 200.0 | 180.0 | |
| 110 1,2,4-Trimethylbenzene | 105 | 12.403 | 12.408 | -0.005 | 97 | 2260604 | 200.0 | 184.9 | |
| 111 1,2-dichloro-4-(trifluorom | 214 | 12.452 | 12.456 | -0.004 | 97 | 542681 | 200.0 | 177.2 | |
| 112 sec-Butylbenzene | 105 | 12.574 | 12.572 | 0.002 | 95 | 2474312 | 200.0 | 176.4 | |
| 113 1,3-Dichlorobenzene | 146 | 12.689 | 12.688 | 0.001 | 97 | 1215884 | 200.0 | 185.0 | |
| 114 4-Isopropyltoluene | 119 | 12.732 | 12.730 | 0.002 | 96 | 2107989 | 200.0 | 180.4 | |
| 115 1,4-Dichlorobenzene | 146 | 12.799 | 12.797 | 0.002 | 95 | 1249173 | 200.0 | 185.1 | |
| 116 2,4-Dichloro-1-(trifluorom | 214 | 12.829 | 12.828 | 0.001 | 95 | 497225 | 200.0 | 174.4 | |
| 118 2,5-Dichlorobenzotrifluori | 214 | 12.872 | 12.870 | 0.002 | 0 | 580659 | 200.0 | 188.5 | |
| 120 n-Butylbenzene | 91 | 13.151 | 13.150 | 0.001 | 96 | 1729209 | 200.0 | 181.5 | |
| 121 1,2-Dichlorobenzene | 146 | 13.158 | 13.156 | 0.002 | 97 | 1161072 | 200.0 | 185.4 | |
| 122 1,2-Dibromo-3-Chloropropan | 75 | 13.973 | 13.971 | 0.002 | 85 | 151695 | 200.0 | 218.1 | |
| 123 2,4- & 2,5- & 2,6- Dichlor | 125 | 14.119 | 14.117 | 0.002 | 0 | 2228710 | 600.0 | 561.0 | |
| 125 2,3- & 3,4- Dichlorotoluen | 125 | 14.551 | 14.555 | -0.004 | 0 | 1589536 | 400.0 | 386.9 | |
| 126 1,2,4-Trichlorobenzene | 180 | 14.830 | 14.829 | 0.001 | 94 | 552245 | 200.0 | 192.7 | |
| 127 Hexachlorobutadiene | 225 | 14.995 | 14.993 | 0.002 | 98 | 180140 | 200.0 | 171.8 | |
| 128 Naphthalene | 128 | 15.104 | 15.103 | 0.001 | 97 | 2008065 | 200.0 | 205.7 | |
| 129 1,2,3-Trichlorobenzene | 180 | 15.348 | 15.346 | 0.002 | 96 | 497473 | 200.0 | 190.0 | |
| 131 2,4,5-Trichlorotoluene | 159 | 16.199 | 16.198 | 0.001 | 0 | 253594 | 200.0 | 203.8 | |
| 130 2,3,6-Trichlorotoluene | 159 | 16.303 | 16.307 | -0.004 | 97 | 237299 | 200.0 | 205.0 | |
| 149 3,4-Dichlorotoluene | 1 | | 0.000 | | | | ND | ND | |
| S 134 1,2-Dichloroethene, Total | 96 | | | | 0 | | 400.0 | 374.5 | |
| S 133 Xylenes, Total | 106 | | | | 0 | | 400.0 | 364.3 | |
| S 135 1,3-Dichloropropene, Total | 1 | | | | 0 | | 400.0 | 418.0 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Reagents:

| | | |
|---------------------|---------------------|-----------|
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 8.00 | Units: uL |
| VOA8260VOAPRI_00263 | Amount Added: 8.00 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 8.00 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 10.00 | Units: uL |
| voaWVA1stRest_00017 | Amount Added: 8.00 | Units: uL |
| voaWEEmix1stR_00009 | Amount Added: 8.00 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 8.00 | Units: uL |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D08.D

Injection Date: 27-Jul-2017 03:13:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: IC VSTD40

Worklist Smp#: 8

Client ID:

Purge Vol: 5.000 mL

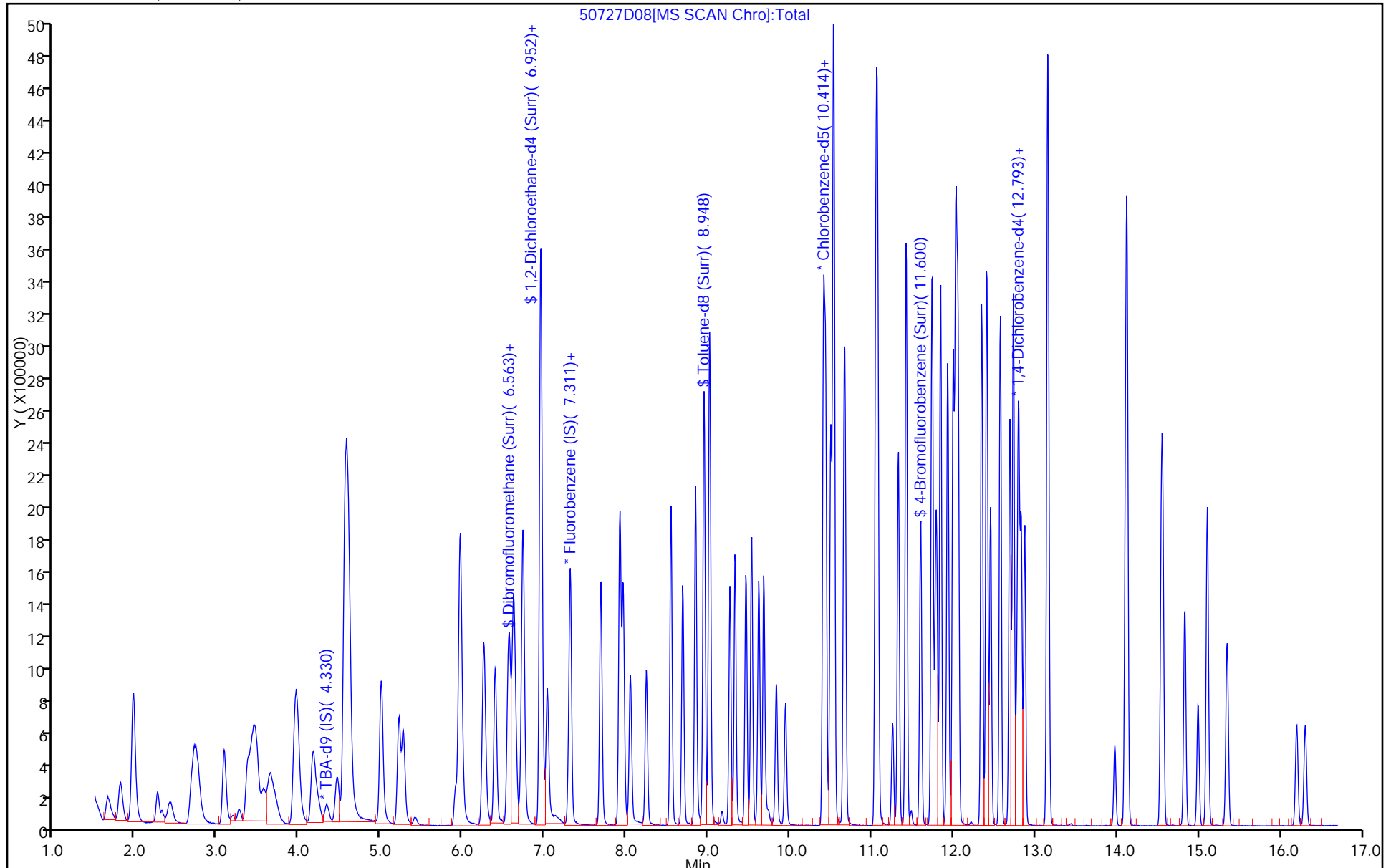
Dil. Factor: 1.0000

ALS Bottle#: 8

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D10.D
 Lims ID: IC VSTD35
 Client ID:
 Sample Type: IC Calib Level: 6
 Inject. Date: 27-Jul-2017 04:00:30 ALS Bottle#: 10 Worklist Smp#: 10
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017756-010
 Misc. Info.: IC VSTD35
 Operator ID: 034635 Instrument ID: CHHP5
 Sublist: chrom-MSVOA_LL_CHHP5*sub12
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 28-Jul-2017 01:05:06 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK029

First Level Reviewer: bungardf

Date: 27-Jul-2017 04:42:28

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.328 | 4.323 | 0.005 | 0 | 232894 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.296 | 7.298 | -0.002 | 94 | 610088 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.405 | 10.406 | -0.001 | 86 | 155120 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.771 | 12.773 | -0.002 | 90 | 193547 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.572 | 6.574 | -0.002 | 94 | 505019 | 175.0 | 172.0 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.943 | 6.945 | -0.002 | 0 | 575099 | 175.0 | 160.6 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.951 | 8.946 | 0.005 | 92 | 1992609 | 175.0 | 161.4 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.597 | 11.599 | -0.002 | 87 | 748217 | 175.0 | 167.8 | |
| 11 Dichlorodifluoromethane | 85 | 1.651 | 1.646 | 0.005 | 99 | 647803 | 175.0 | 182.6 | |
| 12 Chloromethane | 50 | 1.809 | 1.804 | 0.005 | 99 | 595751 | 175.0 | 167.1 | |
| 13 Vinyl chloride | 62 | 1.961 | 1.944 | 0.017 | 98 | 632153 | 175.0 | 174.7 | |
| 14 Butadiene | 39 | 1.967 | 1.969 | -0.001 | 93 | 579584 | 175.0 | 176.3 | |
| 15 Bromomethane | 94 | 2.265 | 2.254 | 0.011 | 91 | 285707 | 175.0 | 167.0 | |
| 16 Chloroethane | 64 | 2.417 | 2.419 | -0.002 | 99 | 340168 | 175.0 | 171.1 | |
| 17 Dichlorofluoromethane | 67 | 2.703 | 2.699 | 0.004 | 97 | 845136 | 175.0 | 168.0 | |
| 18 Trichlorofluoromethane | 101 | 2.746 | 2.741 | 0.005 | 96 | 769762 | 175.0 | 173.1 | |
| 20 Ethyl ether | 59 | 3.074 | 3.076 | -0.002 | 88 | 475422 | 175.0 | 164.4 | |
| 21 Acrolein | 56 | 3.269 | 3.252 | 0.017 | 99 | 154738 | 225.0 | 212.3 | |
| 22 1,1-Dichloroethene | 96 | 3.372 | 3.368 | 0.004 | 96 | 540044 | 175.0 | 180.8 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.452 | 3.441 | 0.011 | 92 | 571742 | 175.0 | 174.4 | |
| 24 Acetone | 43 | 3.482 | 3.477 | 0.005 | 99 | 447756 | 350.0 | 280.6 | |
| 25 Iodomethane | 142 | 3.561 | 3.562 | -0.001 | 96 | 811997 | 175.0 | 173.1 | |
| 26 Carbon disulfide | 76 | 3.646 | 3.648 | -0.002 | 99 | 1310811 | 175.0 | 200.0 | |
| 28 3-Chloro-1-propene | 76 | 3.944 | 3.946 | -0.002 | 93 | 365237 | 175.0 | 189.2 | |
| 30 Methyl acetate | 43 | 3.975 | 3.976 | -0.001 | 97 | 1009713 | 350.0 | 319.6 | |
| 31 Methylene Chloride | 84 | 4.163 | 4.165 | -0.002 | 89 | 602402 | 175.0 | 170.4 | |
| 32 2-Methyl-2-propanol | 59 | 4.455 | 4.451 | 0.004 | 93 | 524619 | 1750.0 | 1904.7 | |
| 33 Acrylonitrile | 53 | 4.553 | 4.554 | -0.001 | 99 | 2362587 | 1750.0 | 1538.0 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.577 | 4.584 | -0.007 | 98 | 595572 | 175.0 | 175.0 | |
| 35 Methyl tert-butyl ether | 73 | 4.601 | 4.603 | -0.002 | 96 | 1597553 | 175.0 | 175.1 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| 36 Hexane | 57 | 4.997 | 4.998 | -0.001 | 91 | 760411 | 175.0 | 174.1 | |
| 37 1,1-Dichloroethane | 63 | 5.216 | 5.217 | -0.001 | 96 | 1024340 | 175.0 | 173.1 | |
| 38 Vinyl acetate | 43 | 5.270 | 5.272 | -0.002 | 97 | 1068205 | 175.0 | 177.5 | |
| 44 2,2-Dichloropropane | 97 | 5.958 | 5.959 | -0.001 | 91 | 136605 | 175.0 | 181.3 | |
| 45 cis-1,2-Dichloroethene | 96 | 5.964 | 5.965 | -0.001 | 79 | 671208 | 175.0 | 172.4 | |
| 46 2-Butanone (MEK) | 43 | 5.982 | 5.978 | 0.004 | 100 | 686266 | 350.0 | 302.2 | |
| 49 Chlorobromomethane | 128 | 6.250 | 6.245 | 0.005 | 95 | 291754 | 175.0 | 168.6 | |
| 51 Tetrahydrofuran | 42 | 6.262 | 6.263 | -0.001 | 87 | 396477 | 350.0 | 299.8 | |
| 52 Chloroform | 83 | 6.396 | 6.391 | 0.005 | 92 | 989929 | 175.0 | 167.5 | |
| 53 1,1,1-Trichloroethane | 97 | 6.554 | 6.549 | 0.005 | 98 | 811476 | 175.0 | 181.4 | |
| 54 Cyclohexane | 56 | 6.621 | 6.622 | -0.001 | 90 | 1012965 | 175.0 | 183.5 | |
| 56 Carbon tetrachloride | 117 | 6.718 | 6.726 | -0.008 | 97 | 682784 | 175.0 | 183.4 | |
| 55 1,1-Dichloropropene | 75 | 6.737 | 6.738 | -0.001 | 97 | 866715 | 175.0 | 179.4 | |
| 57 Isobutyl alcohol | 41 | 6.950 | 6.945 | 0.005 | 91 | 452876 | 4375.0 | 3730.6 | |
| 58 Benzene | 78 | 6.956 | 6.951 | 0.005 | 97 | 2459963 | 175.0 | 165.8 | |
| 59 1,2-Dichloroethane | 62 | 7.029 | 7.030 | -0.001 | 97 | 708898 | 175.0 | 163.9 | |
| 62 n-Heptane | 43 | 7.315 | 7.316 | -0.001 | 88 | 633483 | 175.0 | 181.4 | |
| 64 Trichloroethene | 130 | 7.686 | 7.687 | -0.001 | 98 | 648262 | 175.0 | 173.7 | |
| 66 Methylcyclohexane | 83 | 7.917 | 7.918 | -0.001 | 87 | 1041060 | 175.0 | 184.4 | |
| 67 1,2-Dichloropropane | 63 | 7.959 | 7.961 | -0.002 | 95 | 596512 | 175.0 | 172.7 | |
| 68 Dibromomethane | 93 | 8.045 | 8.046 | -0.001 | 96 | 342853 | 175.0 | 169.4 | |
| 70 1,4-Dioxane | 88 | 8.045 | 8.052 | -0.007 | 39 | 115916 | 3500.0 | 3300.1 | |
| 71 Dichlorobromomethane | 83 | 8.239 | 8.241 | -0.002 | 100 | 712434 | 175.0 | 179.3 | |
| 73 2-Chloroethyl vinyl ether | 63 | 8.543 | 8.545 | -0.002 | 92 | 864836 | 350.0 | 347.9 | |
| 74 cis-1,3-Dichloropropene | 75 | 8.689 | 8.685 | 0.004 | 96 | 881560 | 175.0 | 182.7 | |
| 75 4-Methyl-2-pentanone (MIBK) | 43 | 8.841 | 8.843 | -0.002 | 95 | 1265241 | 350.0 | 318.0 | |
| 76 Toluene | 91 | 9.018 | 9.019 | -0.001 | 98 | 2496911 | 175.0 | 161.4 | |
| 77 trans-1,3-Dichloropropene | 75 | 9.267 | 9.269 | -0.002 | 93 | 781619 | 175.0 | 185.7 | |
| 78 Ethyl methacrylate | 69 | 9.328 | 9.330 | -0.002 | 88 | 905216 | 175.0 | 178.4 | |
| 79 1,1,2-Trichloroethane | 97 | 9.462 | 9.457 | 0.005 | 90 | 523017 | 175.0 | 162.3 | |
| 80 Tetrachloroethene | 164 | 9.529 | 9.530 | -0.001 | 97 | 498519 | 175.0 | 169.0 | |
| 81 1,3-Dichloropropane | 76 | 9.620 | 9.615 | 0.005 | 89 | 969241 | 175.0 | 162.7 | |
| 82 2-Hexanone | 43 | 9.681 | 9.682 | -0.001 | 94 | 977068 | 350.0 | 320.2 | |
| 84 Chlorodibromomethane | 129 | 9.833 | 9.834 | -0.001 | 90 | 489506 | 175.0 | 179.7 | |
| 85 Ethylene Dibromide | 107 | 9.943 | 9.944 | -0.001 | 99 | 550826 | 175.0 | 166.7 | |
| 86 3-Chlorobenzotrifluoride | 180 | 10.411 | 10.412 | -0.001 | 93 | 874266 | 175.0 | 164.0 | |
| 87 Chlorobenzene | 112 | 10.435 | 10.437 | -0.002 | 94 | 1645967 | 175.0 | 163.5 | |
| 88 4-Chlorobenzotrifluoride | 180 | 10.496 | 10.498 | -0.002 | 95 | 826850 | 175.0 | 168.1 | |
| 89 1,1,1,2-Tetrachloroethane | 131 | 10.527 | 10.528 | -0.001 | 93 | 554351 | 175.0 | 173.1 | |
| 90 Ethylbenzene | 106 | 10.533 | 10.534 | -0.001 | 97 | 962208 | 175.0 | 171.2 | |
| 91 m-Xylene & p-Xylene | 106 | 10.667 | 10.668 | -0.001 | 0 | 1197380 | 175.0 | 174.3 | |
| 92 o-Xylene | 106 | 11.050 | 11.051 | -0.001 | 95 | 1130677 | 175.0 | 172.8 | |
| 93 Styrene | 104 | 11.068 | 11.069 | -0.001 | 94 | 1866053 | 175.0 | 168.4 | |
| 94 Bromoform | 173 | 11.257 | 11.252 | 0.005 | 97 | 310948 | 175.0 | 183.7 | |
| 96 2-Chlorobenzotrifluoride | 180 | 11.324 | 11.325 | -0.001 | 96 | 840920 | 175.0 | 164.9 | |
| 97 Isopropylbenzene | 105 | 11.421 | 11.422 | -0.001 | 96 | 2681266 | 175.0 | 167.8 | |
| 100 Bromobenzene | 156 | 11.737 | 11.739 | -0.002 | 95 | 659984 | 175.0 | 175.7 | |
| 99 1,1,2,2-Tetrachloroethane | 83 | 11.737 | 11.745 | -0.008 | 94 | 762601 | 175.0 | 159.9 | |
| 102 trans-1,4-Dichloro-2-buten | 53 | 11.774 | 11.775 | -0.001 | 86 | 199800 | 175.0 | 176.4 | |
| 101 1,2,3-Trichloropropane | 110 | 11.792 | 11.793 | -0.001 | 85 | 255265 | 175.0 | 164.7 | |
| 103 N-Propylbenzene | 120 | 11.841 | 11.842 | -0.001 | 97 | 786064 | 175.0 | 183.1 | |
| 104 2-Chlorotoluene | 126 | 11.926 | 11.927 | -0.001 | 97 | 666236 | 175.0 | 179.6 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 105 3-Chlorotoluene | 126 | 11.993 | 11.994 | -0.001 | 96 | 680717 | 175.0 | 168.7 | |
| 106 1,3,5-Trimethylbenzene | 105 | 12.029 | 12.031 | -0.002 | 94 | 2153457 | 175.0 | 175.3 | |
| 107 4-Chlorotoluene | 126 | 12.054 | 12.055 | -0.001 | 95 | 719035 | 175.0 | 179.5 | |
| 108 tert-Butylbenzene | 119 | 12.346 | 12.347 | -0.001 | 93 | 1844417 | 175.0 | 179.6 | |
| 110 1,2,4-Trimethylbenzene | 105 | 12.406 | 12.408 | -0.002 | 97 | 2182090 | 175.0 | 174.8 | |
| 111 1,2-dichloro-4-(trifluorom | 214 | 12.455 | 12.456 | -0.001 | 97 | 525922 | 175.0 | 168.1 | |
| 112 sec-Butylbenzene | 105 | 12.571 | 12.572 | -0.001 | 94 | 2514051 | 175.0 | 175.5 | |
| 113 1,3-Dichlorobenzene | 146 | 12.692 | 12.688 | 0.004 | 96 | 1146674 | 175.0 | 170.8 | |
| 114 4-Isopropyltoluene | 119 | 12.729 | 12.730 | -0.001 | 96 | 2114911 | 175.0 | 177.2 | |
| 115 1,4-Dichlorobenzene | 146 | 12.796 | 12.797 | -0.001 | 95 | 1174377 | 175.0 | 170.4 | |
| 116 2,4-Dichloro-1-(trifluorom | 214 | 12.826 | 12.828 | -0.002 | 96 | 501975 | 175.0 | 172.4 | |
| 118 2,5-Dichlorobenzotrifluori | 214 | 12.875 | 12.870 | 0.005 | 0 | 541324 | 175.0 | 172.1 | |
| 120 n-Butylbenzene | 91 | 13.149 | 13.150 | -0.001 | 96 | 1748217 | 175.0 | 179.6 | |
| 121 1,2-Dichlorobenzene | 146 | 13.161 | 13.156 | 0.005 | 97 | 1081541 | 175.0 | 169.1 | |
| 122 1,2-Dibromo-3-Chloropropan | 75 | 13.970 | 13.971 | -0.001 | 86 | 125814 | 175.0 | 177.1 | |
| 123 2,4- & 2,5- & 2,6- Dichlor | 125 | 14.116 | 14.117 | -0.001 | 0 | 2069215 | 525.0 | 509.9 | |
| 125 2,3- & 3,4- Dichlorotoluen | 125 | 14.548 | 14.555 | -0.007 | 0 | 1443949 | 350.0 | 344.1 | |
| 126 1,2,4-Trichlorobenzene | 180 | 14.828 | 14.829 | -0.001 | 95 | 511830 | 175.0 | 174.8 | |
| 127 Hexachlorobutadiene | 225 | 14.992 | 14.993 | -0.001 | 98 | 182711 | 175.0 | 170.6 | |
| 128 Naphthalene | 128 | 15.101 | 15.103 | -0.002 | 97 | 1761559 | 175.0 | 176.7 | |
| 129 1,2,3-Trichlorobenzene | 180 | 15.345 | 15.346 | -0.001 | 96 | 453926 | 175.0 | 169.7 | |
| 131 2,4,5-Trichlorotoluene | 159 | 16.196 | 16.198 | -0.002 | 0 | 235417 | 175.0 | 185.2 | |
| 130 2,3,6-Trichlorotoluene | 159 | 16.306 | 16.307 | -0.001 | 97 | 211883 | 175.0 | 179.2 | |
| 149 3,4-Dichlorotoluene | 1 | | 0.000 | | | | ND | ND | |
| S 133 Xylenes, Total | 106 | | | | 0 | | 350.0 | 347.1 | |
| S 134 1,2-Dichloroethene, Total | 96 | | | | 0 | | 350.0 | 347.4 | |
| S 135 1,3-Dichloropropene, Total | 1 | | | | 0 | | 350.0 | 368.4 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Reagents:

| | | |
|---------------------|--------------------|-----------|
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260VOAPRI_00263 | Amount Added: 7.00 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 9.00 | Units: uL |
| voaWVA1stRest_00017 | Amount Added: 7.00 | Units: uL |
| voaWEEmix1stR_00009 | Amount Added: 7.00 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 7.00 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 7.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 7.00 | Units: uL |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D10.D

Injection Date: 27-Jul-2017 04:00:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: IC VSTD35

Worklist Smp#: 10

Client ID:

Purge Vol: 5.000 mL

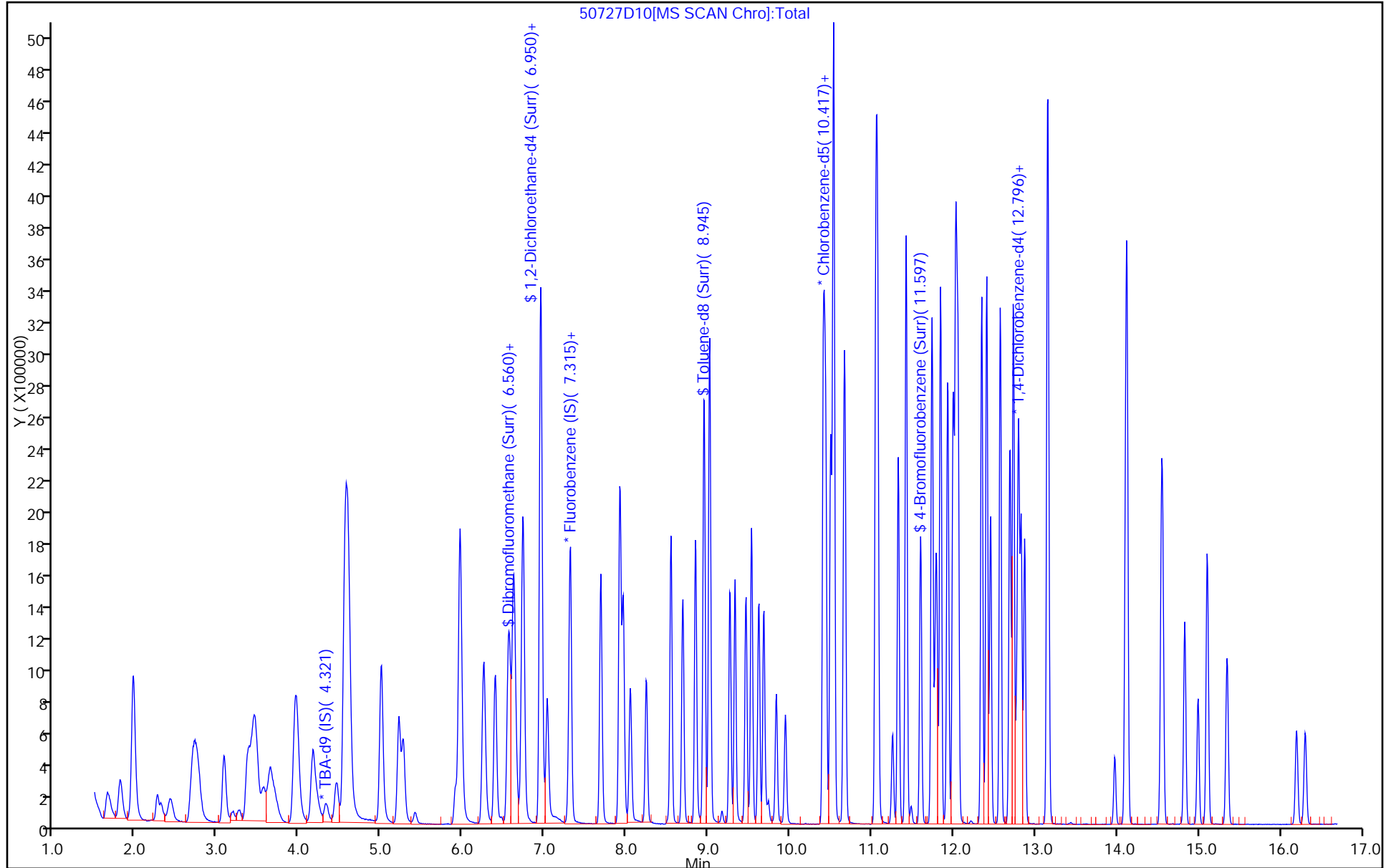
Dil. Factor: 1.0000

ALS Bottle#: 10

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Lims ID: IC VSTD50
 Client ID:
 Sample Type: IC Calib Level: 8
 Inject. Date: 27-Jul-2017 04:24:30 ALS Bottle#: 11 Worklist Smp#: 11
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0017756-011
 Misc. Info.: IC VSTD50
 Operator ID: 034635 Instrument ID: CHHP5
 Sublist: chrom-MSVOA_LL_CHHP5*sub12
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 28-Jul-2017 01:05:08 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK029

First Level Reviewer: bungardf

Date: 27-Jul-2017 05:09:00

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.341 | 4.323 | 0.018 | 0 | 184114 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.297 | 7.298 | -0.001 | 99 | 607808 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.406 | 10.406 | 0.000 | 85 | 161595 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.772 | 12.773 | -0.001 | 89 | 194624 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.573 | 6.574 | -0.001 | 94 | 681339 | 250.0 | 233.0 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.944 | 6.945 | -0.001 | 0 | 795993 | 250.0 | 223.2 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.946 | 8.946 | 0.000 | 92 | 2678162 | 250.0 | 208.2 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.598 | 11.599 | -0.001 | 87 | 1033645 | 250.0 | 222.5 | |
| 11 Dichlorodifluoromethane | 85 | 1.652 | 1.646 | 0.006 | 99 | 857078 | 250.0 | 242.5 | |
| 12 Chloromethane | 50 | 1.804 | 1.804 | 0.000 | 99 | 811941 | 250.0 | 228.6 | |
| 13 Vinyl chloride | 62 | 1.956 | 1.944 | 0.012 | 98 | 867536 | 250.0 | 240.7 | |
| 14 Butadiene | 39 | 1.968 | 1.969 | 0.000 | 94 | 815610 | 250.0 | 249.1 | |
| 15 Bromomethane | 94 | 2.266 | 2.254 | 0.012 | 90 | 377950 | 250.0 | 221.8 | |
| 16 Chloroethane | 64 | 2.406 | 2.419 | -0.013 | 99 | 414342 | 250.0 | 209.1 | |
| 17 Dichlorofluoromethane | 67 | 2.698 | 2.699 | -0.001 | 97 | 1057272 | 250.0 | 211.0 | |
| 18 Trichlorofluoromethane | 101 | 2.728 | 2.741 | -0.013 | 97 | 1017488 | 250.0 | 229.7 | |
| 20 Ethyl ether | 59 | 3.069 | 3.076 | -0.007 | 88 | 612640 | 250.0 | 212.6 | |
| 21 Acrolein | 56 | 3.264 | 3.252 | 0.012 | 98 | 183852 | 275.0 | 253.2 | |
| 22 1,1-Dichloroethene | 96 | 3.367 | 3.368 | -0.001 | 97 | 745282 | 250.0 | 250.5 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.428 | 3.441 | -0.013 | 92 | 774058 | 250.0 | 237.0 | |
| 24 Acetone | 43 | 3.483 | 3.477 | 0.006 | 100 | 630881 | 500.0 | 396.9 | |
| 25 Iodomethane | 142 | 3.580 | 3.562 | 0.018 | 97 | 1099819 | 250.0 | 235.3 | |
| 26 Carbon disulfide | 76 | 3.647 | 3.648 | -0.001 | 99 | 1856339 | 250.0 | 284.2 | |
| 28 3-Chloro-1-propene | 76 | 3.939 | 3.946 | -0.007 | 93 | 500032 | 250.0 | 260.0 | |
| 30 Methyl acetate | 43 | 3.976 | 3.976 | 0.000 | 97 | 1447736 | 500.0 | 459.9 | |
| 31 Methylene Chloride | 84 | 4.164 | 4.165 | -0.001 | 88 | 813282 | 250.0 | 232.1 | |
| 32 2-Methyl-2-propanol | 59 | 4.468 | 4.451 | 0.017 | 94 | 568135 | 2500.0 | 2609.2 | |
| 33 Acrylonitrile | 53 | 4.553 | 4.554 | -0.001 | 98 | 3495451 | 2500.0 | 2284.0 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.578 | 4.584 | -0.006 | 98 | 806194 | 250.0 | 237.8 | |
| 35 Methyl tert-butyl ether | 73 | 4.602 | 4.603 | -0.001 | 96 | 2170401 | 250.0 | 238.7 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| 36 Hexane | 57 | 4.991 | 4.998 | -0.007 | 92 | 1101558 | 250.0 | 253.1 | |
| 37 1,1-Dichloroethane | 63 | 5.210 | 5.217 | -0.007 | 96 | 1376176 | 250.0 | 233.4 | |
| 38 Vinyl acetate | 43 | 5.271 | 5.272 | -0.001 | 97 | 1523056 | 250.0 | 254.0 | |
| 44 2,2-Dichloropropane | 97 | 5.959 | 5.959 | 0.000 | 91 | 188250 | 250.0 | 250.8 | |
| 45 cis-1,2-Dichloroethene | 96 | 5.959 | 5.965 | -0.006 | 79 | 900432 | 250.0 | 232.2 | |
| 46 2-Butanone (MEK) | 43 | 5.983 | 5.978 | 0.005 | 98 | 962704 | 500.0 | 425.5 | |
| 49 Chlorobromomethane | 128 | 6.245 | 6.245 | 0.000 | 94 | 394763 | 250.0 | 229.0 | |
| 51 Tetrahydrofuran | 42 | 6.263 | 6.263 | 0.000 | 87 | 609910 | 500.0 | 462.9 | |
| 52 Chloroform | 83 | 6.391 | 6.391 | 0.000 | 92 | 1319564 | 250.0 | 224.1 | |
| 53 1,1,1-Trichloroethane | 97 | 6.549 | 6.549 | 0.000 | 98 | 1097196 | 250.0 | 246.2 | |
| 54 Cyclohexane | 56 | 6.616 | 6.622 | -0.006 | 90 | 1394833 | 250.0 | 253.7 | |
| 56 Carbon tetrachloride | 117 | 6.719 | 6.726 | -0.007 | 97 | 923177 | 250.0 | 248.9 | |
| 55 1,1-Dichloropropene | 75 | 6.737 | 6.738 | -0.001 | 96 | 1178056 | 250.0 | 244.7 | |
| 57 Isobutyl alcohol | 41 | 6.950 | 6.945 | 0.005 | 68 | 715201 | 6250.0 | 5913.6 | |
| 58 Benzene | 78 | 6.950 | 6.951 | -0.001 | 97 | 3249284 | 250.0 | 219.9 | |
| 59 1,2-Dichloroethane | 62 | 7.029 | 7.030 | -0.001 | 97 | 969148 | 250.0 | 225.0 | |
| 62 n-Heptane | 43 | 7.309 | 7.316 | -0.007 | 89 | 922592 | 250.0 | 265.1 | |
| 64 Trichloroethene | 130 | 7.686 | 7.687 | -0.001 | 98 | 887332 | 250.0 | 238.6 | |
| 66 Methylcyclohexane | 83 | 7.918 | 7.918 | 0.000 | 87 | 1432791 | 250.0 | 254.8 | |
| 67 1,2-Dichloropropane | 63 | 7.960 | 7.961 | -0.001 | 95 | 793667 | 250.0 | 230.6 | |
| 68 Dibromomethane | 93 | 8.045 | 8.046 | -0.001 | 97 | 470836 | 250.0 | 233.5 | |
| 70 1,4-Dioxane | 88 | 8.039 | 8.052 | -0.013 | 38 | 187034 | 5000.0 | 5344.8 | |
| 71 Dichlorobromomethane | 83 | 8.240 | 8.241 | -0.001 | 100 | 945026 | 250.0 | 238.8 | |
| 73 2-Chloroethyl vinyl ether | 63 | 8.544 | 8.545 | -0.001 | 92 | 1234429 | 500.0 | 498.5 | |
| 74 cis-1,3-Dichloropropene | 75 | 8.684 | 8.685 | -0.001 | 96 | 1203144 | 250.0 | 250.3 | |
| 75 4-Methyl-2-pentanone (MIBK) | 43 | 8.842 | 8.843 | -0.001 | 94 | 1863520 | 500.0 | 449.6 | |
| 76 Toluene | 91 | 9.019 | 9.019 | 0.000 | 97 | 3254284 | 250.0 | 202.0 | |
| 77 trans-1,3-Dichloropropene | 75 | 9.268 | 9.269 | -0.001 | 93 | 1070347 | 250.0 | 244.1 | |
| 78 Ethyl methacrylate | 69 | 9.329 | 9.330 | -0.001 | 88 | 1271580 | 250.0 | 240.5 | |
| 79 1,1,2-Trichloroethane | 97 | 9.457 | 9.457 | 0.000 | 91 | 718069 | 250.0 | 213.9 | |
| 80 Tetrachloroethene | 164 | 9.530 | 9.530 | 0.000 | 97 | 683462 | 250.0 | 222.4 | |
| 81 1,3-Dichloropropane | 76 | 9.621 | 9.615 | 0.006 | 89 | 1320887 | 250.0 | 212.9 | |
| 82 2-Hexanone | 43 | 9.676 | 9.682 | -0.006 | 93 | 1418811 | 500.0 | 446.3 | |
| 84 Chlorodibromomethane | 129 | 9.834 | 9.834 | 0.000 | 90 | 672369 | 250.0 | 237.0 | |
| 85 Ethylene Dibromide | 107 | 9.943 | 9.944 | -0.001 | 99 | 773664 | 250.0 | 224.7 | |
| 86 3-Chlorobenzotrifluoride | 180 | 10.412 | 10.412 | 0.000 | 93 | 1290067 | 250.0 | 232.3 | |
| 87 Chlorobenzene | 112 | 10.436 | 10.437 | -0.001 | 95 | 2170926 | 250.0 | 207.0 | |
| 88 4-Chlorobenzotrifluoride | 180 | 10.497 | 10.498 | -0.001 | 96 | 1226371 | 250.0 | 239.3 | |
| 89 1,1,1,2-Tetrachloroethane | 131 | 10.527 | 10.528 | -0.001 | 94 | 751692 | 250.0 | 225.4 | |
| 90 Ethylbenzene | 106 | 10.533 | 10.534 | -0.001 | 97 | 1304914 | 250.0 | 222.8 | |
| 91 m-Xylene & p-Xylene | 106 | 10.667 | 10.668 | -0.001 | 0 | 1614353 | 250.0 | 225.6 | |
| 92 o-Xylene | 106 | 11.051 | 11.051 | 0.000 | 95 | 1518391 | 250.0 | 222.7 | |
| 93 Styrene | 104 | 11.069 | 11.069 | 0.000 | 94 | 2462559 | 250.0 | 213.4 | |
| 94 Bromoform | 173 | 11.257 | 11.252 | 0.005 | 98 | 443094 | 250.0 | 251.3 | |
| 96 2-Chlorobenzotrifluoride | 180 | 11.324 | 11.325 | -0.001 | 95 | 1244752 | 250.0 | 234.2 | |
| 97 Isopropylbenzene | 105 | 11.422 | 11.422 | 0.000 | 96 | 3502176 | 250.0 | 210.4 | |
| 100 Bromobenzene | 156 | 11.738 | 11.739 | -0.001 | 95 | 889999 | 250.0 | 235.6 | |
| 99 1,1,2,2-Tetrachloroethane | 83 | 11.738 | 11.745 | -0.007 | 95 | 1078742 | 250.0 | 217.1 | |
| 102 trans-1,4-Dichloro-2-buten | 53 | 11.781 | 11.775 | 0.006 | 84 | 299994 | 250.0 | 263.4 | |
| 101 1,2,3-Trichloropropane | 110 | 11.793 | 11.793 | 0.000 | 84 | 371250 | 250.0 | 238.1 | |
| 103 N-Propylbenzene | 120 | 11.841 | 11.842 | -0.001 | 96 | 1069171 | 250.0 | 247.7 | |
| 104 2-Chlorotoluene | 126 | 11.927 | 11.927 | 0.000 | 97 | 907016 | 250.0 | 243.1 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 105 3-Chlorotoluene | 126 | 11.994 | 11.994 | 0.000 | 96 | 1010916 | 250.0 | 249.1 | |
| 106 1,3,5-Trimethylbenzene | 105 | 12.030 | 12.031 | -0.001 | 95 | 2828999 | 250.0 | 229.0 | |
| 107 4-Chlorotoluene | 126 | 12.054 | 12.055 | -0.001 | 96 | 970169 | 250.0 | 240.8 | |
| 108 tert-Butylbenzene | 119 | 12.346 | 12.347 | -0.001 | 92 | 2446270 | 250.0 | 236.9 | |
| 110 1,2,4-Trimethylbenzene | 105 | 12.407 | 12.408 | -0.001 | 97 | 2860516 | 250.0 | 227.8 | |
| 111 1,2-dichloro-4-(trifluorom | 214 | 12.456 | 12.456 | 0.000 | 96 | 801099 | 250.0 | 254.7 | |
| 112 sec-Butylbenzene | 105 | 12.571 | 12.572 | -0.001 | 95 | 3330508 | 250.0 | 231.2 | |
| 113 1,3-Dichlorobenzene | 146 | 12.687 | 12.688 | -0.001 | 96 | 1545747 | 250.0 | 229.0 | |
| 114 4-Isopropyltoluene | 119 | 12.730 | 12.730 | 0.000 | 95 | 2809716 | 250.0 | 234.1 | |
| 115 1,4-Dichlorobenzene | 146 | 12.797 | 12.797 | 0.000 | 95 | 1574222 | 250.0 | 227.2 | |
| 116 2,4-Dichloro-1-(trifluorom | 214 | 12.827 | 12.828 | -0.001 | 94 | 771761 | 250.0 | 263.5 | |
| 118 2,5-Dichlorobenzotrifluori | 214 | 12.870 | 12.870 | 0.000 | 0 | 797256 | 250.0 | 252.0 | |
| 120 n-Butylbenzene | 91 | 13.149 | 13.150 | -0.001 | 95 | 2372703 | 250.0 | 242.4 | |
| 121 1,2-Dichlorobenzene | 146 | 13.155 | 13.156 | -0.001 | 96 | 1435184 | 250.0 | 223.1 | |
| 122 1,2-Dibromo-3-Chloropropan | 75 | 13.971 | 13.971 | 0.000 | 86 | 182290 | 250.0 | 255.2 | |
| 123 2,4- & 2,5- & 2,6- Dichlor | 125 | 14.117 | 14.117 | 0.000 | 0 | 3049908 | 750.0 | 747.4 | |
| 125 2,3- & 3,4- Dichlorotoluen | 125 | 14.555 | 14.555 | 0.000 | 0 | 2191624 | 500.0 | 519.4 | |
| 126 1,2,4-Trichlorobenzene | 180 | 14.828 | 14.829 | -0.001 | 95 | 755690 | 250.0 | 256.7 | |
| 127 Hexachlorobutadiene | 225 | 14.993 | 14.993 | 0.000 | 98 | 282046 | 250.0 | 261.8 | |
| 128 Naphthalene | 128 | 15.102 | 15.103 | -0.001 | 98 | 2561966 | 250.0 | 255.5 | |
| 129 1,2,3-Trichlorobenzene | 180 | 15.346 | 15.346 | 0.000 | 96 | 693791 | 250.0 | 258.0 | |
| 131 2,4,5-Trichlorotoluene | 159 | 16.197 | 16.198 | -0.001 | 0 | 452516 | 250.0 | 354.0 | |
| 130 2,3,6-Trichlorotoluene | 159 | 16.301 | 16.307 | -0.006 | 98 | 417201 | 250.0 | 350.8 | |
| 149 3,4-Dichlorotoluene | 1 | | 0.000 | | | | ND | ND | |
| S 134 1,2-Dichloroethene, Total | 96 | | | | 0 | | 500.0 | 470.0 | |
| S 133 Xylenes, Total | 106 | | | | 0 | | 500.0 | 448.3 | |
| S 135 1,3-Dichloropropene, Total | 1 | | | | 0 | | 500.0 | 494.4 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Reagents:

| | | |
|---------------------|---------------------|-----------|
| VOA8260VOAPRI_00263 | Amount Added: 10.00 | Units: uL |
| voaWAcro1stRe_00016 | Amount Added: 11.00 | Units: uL |
| voaWVA1stRest_00017 | Amount Added: 10.00 | Units: uL |
| voaWEEmix1stR_00009 | Amount Added: 10.00 | Units: uL |
| voaW2clev1stR_00013 | Amount Added: 10.00 | Units: uL |
| voaWKetmix1st_00004 | Amount Added: 10.00 | Units: uL |
| VOA8260INT_00072 | Amount Added: 2.00 | Units: uL |
| VOA8260SURR_00071 | Amount Added: 10.00 | Units: uL |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D

Injection Date: 27-Jul-2017 04:24:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: IC VSTD50

Worklist Smp#: 11

Client ID:

Purge Vol: 5.000 mL

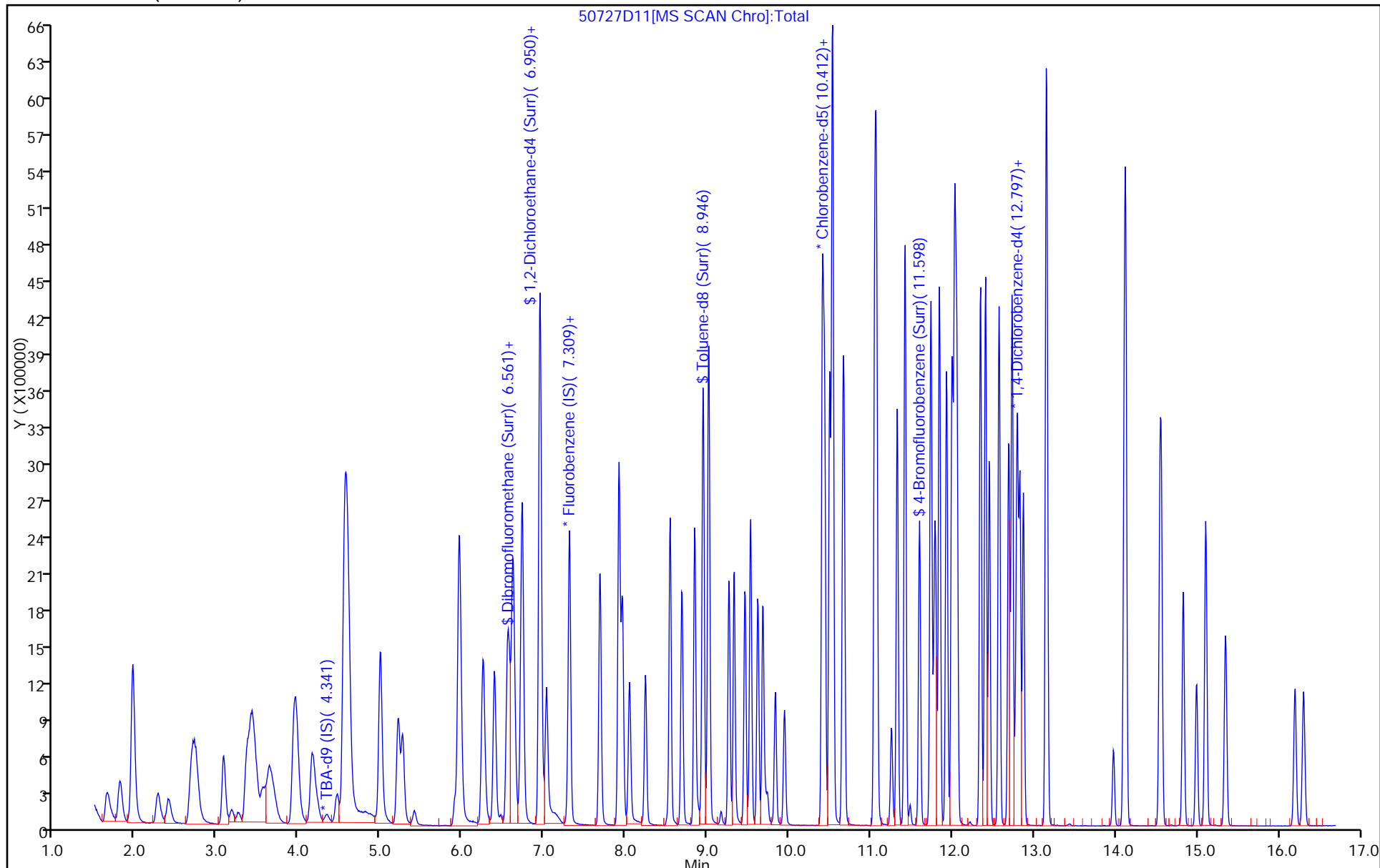
Dil. Factor: 1.0000

ALS Bottle#: 11

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Lab Sample ID: CCVIS 180-224792/2 Calibration Date: 10/04/2017 00:22
 Instrument ID: CHHP5 Calib Start Date: 07/27/2017 00:51
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 07/27/2017 04:24
 Lab File ID: 51003D02.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 | 0.2907 | 0.3094 | 0.1000 | 10.6 | 10.0 | 6.4 | 20.0 |
| Chloromethane | Ave | 0.2922 | 0.3610 | 0.1000 | 12.4 | 10.0 | 23.5* | 20.0 |
| Vinyl chloride | Ave | 0.2965 | 0.3404 | 0.1000 | 11.5 | 10.0 | 14.8 | 20.0 |
| 1,3-Butadiene | Ave | 0.2694 | 0.3290 | 0.0100 | 12.2 | 10.0 | 22.1* | 20.0 |
| Bromomethane | Ave | 0.1402 | 0.1396 | 0.0500 | 9.96 | 10.0 | -0.4 | 20.0 |
| Chloroethane | Ave | 0.1630 | 0.1689 | 0.0500 | 10.4 | 10.0 | 3.6 | 20.0 |
| Trichlorofluoromethane | Ave | 0.3643 | 0.3595 | 0.1000 | 9.87 | 10.0 | -1.3 | 20.0 |
| Ethyl ether | Ave | 0.2370 | 0.2377 | 0.0100 | 10.0 | 10.0 | 0.3 | 20.0 |
| Acrolein | Ave | 0.0597 | 0.0604 | 0.0100 | 30.3 | 30.0 | 1.1 | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2448 | 0.2670 | 0.1000 | 10.9 | 10.0 | 9.1 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.2686 | 0.2909 | 0.1000 | 10.8 | 10.0 | 8.3 | 20.0 |
| Acetone | Ave | 0.1308 | 0.1351 | 0.0500 | 20.7 | 20.0 | 3.3 | 20.0 |
| Iodomethane | Ave | 0.3845 | 0.3797 | 0.0100 | 9.88 | 10.0 | -1.2 | 20.0 |
| Carbon disulfide | Ave | 0.5372 | 0.4405 | 0.1000 | 8.20 | 10.0 | -18.0 | 20.0 |
| Allyl chloride | Ave | 0.1582 | 0.1436 | 0.0100 | 9.08 | 10.0 | -9.2 | 20.0 |
| Methyl acetate | Ave | 0.2589 | 0.2478 | 0.1000 | 19.1 | 20.0 | -4.3 | 20.0 |
| Methylene Chloride | Lin2 | | 0.2845 | 0.1000 | 9.34 | 10.0 | -6.6 | 20.0 |
| tert-Butyl alcohol | Ave | 1.183 | 1.134 | 0.0100 | 95.9 | 100 | -4.1 | 20.0 |
| Acrylonitrile | Ave | 0.1259 | 0.1224 | 0.0100 | 97.2 | 100 | -2.8 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2789 | 0.2795 | 0.1000 | 10.0 | 10.0 | 0.2 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.7479 | 0.6362 | 0.1000 | 8.51 | 10.0 | -14.9 | 20.0 |
| Hexane | Ave | 0.3580 | 0.3661 | 0.0100 | 10.2 | 10.0 | 2.3 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.4850 | 0.4550 | 0.2000 | 9.38 | 10.0 | -6.2 | 20.0 |
| Vinyl acetate | Ave | 0.4932 | 0.4318 | 0.0100 | 8.76 | 10.0 | -12.4 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.0617 | 0.0570 | 0.0100 | 9.23 | 10.0 | -7.7 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.3190 | 0.2941 | 0.1000 | 9.22 | 10.0 | -7.8 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.1861 | 0.1739 | 0.0500 | 18.7 | 20.0 | -6.6 | 20.0 |
| Bromochloromethane | Ave | 0.1418 | 0.1253 | 0.0100 | 8.83 | 10.0 | -11.7 | 20.0 |
| Tetrahydrofuran | Ave | 0.1084 | 0.0881 | 0.0100 | 16.3 | 20.0 | -18.7 | 20.0 |
| Chloroform | Ave | 0.4843 | 0.4505 | 0.2000 | 9.30 | 10.0 | -7.0 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.3666 | 0.3561 | 0.1000 | 9.72 | 10.0 | -2.8 | 20.0 |
| Cyclohexane | Ave | 0.4524 | 0.4360 | 0.1000 | 9.64 | 10.0 | -3.6 | 20.0 |
| Carbon tetrachloride | Ave | 0.3051 | 0.2986 | 0.1000 | 9.79 | 10.0 | -2.1 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3961 | 0.3696 | 0.0100 | 9.33 | 10.0 | -6.7 | 20.0 |
| Isobutyl alcohol | Ave | 0.0099 | 0.0092* | 0.0100 | 231 | 250 | -7.5 | 20.0 |
| Benzene | Ave | 1.216 | 1.124 | 0.5000 | 9.24 | 10.0 | -7.6 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.3544 | 0.3311 | 0.1000 | 9.34 | 10.0 | -6.6 | 20.0 |
| n-Heptane | Ave | 0.2863 | 0.2918 | 0.0100 | 10.2 | 10.0 | 1.9 | 20.0 |
| Trichloroethene | Ave | 0.3059 | 0.2755 | 0.2000 | 9.00 | 10.0 | -10.0 | 20.0 |
| Methylcyclohexane | Ave | 0.4626 | 0.4029 | 0.1000 | 8.71 | 10.0 | -12.9 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Lab Sample ID: CCVIS 180-224792/2 Calibration Date: 10/04/2017 00:22
 Instrument ID: CHHP5 Calib Start Date: 07/27/2017 00:51
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 07/27/2017 04:24
 Lab File ID: 51003D02.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 | 0.2831 | 0.2370 | 0.1000 | 8.37 | 10.0 | -16.3 | 20.0 |
| 1,4-Dioxane | Ave | 0.0029 | 0.0027* | 0.0100 | 184 | 200 | -8.1 | 20.0 |
| Dibromomethane | Ave | 0.1659 | 0.1353 | 0.0100 | 8.15 | 10.0 | -18.5 | 20.0 |
| Bromodichloromethane | Ave | 0.3256 | 0.2505 | 0.2000 | 7.69 | 10.0 | -23.1* | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.2037 | 0.1249 | 0.0100 | 12.3 | 20.0 | -38.7* | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3955 | 0.3042 | 0.2000 | 7.69 | 10.0 | -23.1* | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 1.282 | 1.219 | 0.1000 | 19.0 | 20.0 | -5.0 | 20.0 |
| Toluene | Ave | 4.986 | 5.324 | 0.4000 | 10.7 | 10.0 | 6.8 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 1.357 | 1.226 | 0.1000 | 9.04 | 10.0 | -9.6 | 20.0 |
| Ethyl methacrylate | Ave | 1.636 | 1.280 | 0.0100 | 7.82 | 10.0 | -21.8* | 20.0 |
| 1,1,2-Trichloroethane | Ave | 1.039 | 1.051 | 0.1000 | 10.1 | 10.0 | 1.2 | 20.0 |
| Tetrachloroethene | Ave | 0.9508 | 1.036 | 0.2000 | 10.9 | 10.0 | 8.9 | 20.0 |
| 1,3-Dichloropropane | Ave | 1.920 | 1.737 | 0.0100 | 9.05 | 10.0 | -9.5 | 20.0 |
| 2-Hexanone | Ave | 0.9836 | 0.8982 | 0.1000 | 18.3 | 20.0 | -8.7 | 20.0 |
| Dibromochloromethane | Ave | 0.8779 | 0.7440 | 0.1000 | 8.47 | 10.0 | -15.3 | 20.0 |
| 1,2-Dibromoethane (EDB) | Ave | 1.065 | 0.9904 | 0.1000 | 9.30 | 10.0 | -7.0 | 20.0 |
| 3-Chlorobenzotrifluoride | Ave | 1.718 | 1.982 | 0.0100 | 11.5 | 10.0 | 15.4 | 20.0 |
| Chlorobenzene | Ave | 3.246 | 3.120 | 0.5000 | 9.61 | 10.0 | -3.9 | 20.0 |
| 4-Chlorobenzotrifluoride | Ave | 1.586 | 1.861 | 0.0100 | 11.7 | 10.0 | 17.4 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 1.032 | 0.997 | 0.0100 | 9.66 | 10.0 | -3.4 | 20.0 |
| Ethylbenzene | Ave | 1.812 | 1.771 | 0.1000 | 9.77 | 10.0 | -2.3 | 20.0 |
| m-Xylene & p-Xylene | Ave | 2.214 | 2.198 | 0.1000 | 9.93 | 10.0 | -0.7 | 20.0 |
| o-Xylene | Ave | 2.110 | 2.025 | 0.3000 | 9.60 | 10.0 | -4.0 | 20.0 |
| Styrene | Ave | 3.571 | 3.364 | 0.3000 | 9.42 | 10.0 | -5.8 | 20.0 |
| Bromoform | Ave | 0.5456 | 0.4263 | 0.1000 | 7.81 | 10.0 | -21.9* | 20.0 |
| 2-Chlorobenzotrifluoride | Ave | 1.644 | 1.910 | 0.0100 | 11.6 | 10.0 | 16.2 | 20.0 |
| Isopropylbenzene | Ave | 5.150 | 5.302 | 0.1000 | 10.3 | 10.0 | 3.0 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 1.538 | 1.423 | 0.3000 | 9.26 | 10.0 | -7.4 | 20.0 |
| Bromobenzene | Ave | 0.9704 | 0.8525 | 0.0100 | 8.78 | 10.0 | -12.2 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2926 | 0.2888 | 0.0100 | 9.87 | 10.0 | -1.3 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.4005 | 0.3616 | 0.0100 | 9.03 | 10.0 | -9.7 | 20.0 |
| N-Propylbenzene | Ave | 1.109 | 1.053 | 0.0100 | 9.49 | 10.0 | -5.1 | 20.0 |
| 2-Chlorotoluene | Ave | 0.9585 | 0.9022 | 0.0100 | 9.41 | 10.0 | -5.9 | 20.0 |
| 3-Chlorotoluene | Ave | 1.043 | 1.087 | 0.0100 | 10.4 | 10.0 | 4.3 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.173 | 3.086 | 0.0100 | 9.73 | 10.0 | -2.7 | 20.0 |
| 4-Chlorotoluene | Ave | 1.035 | 0.9831 | 0.0100 | 9.50 | 10.0 | -5.0 | 20.0 |
| tert-Butylbenzene | Ave | 2.653 | 2.393 | 0.0100 | 9.02 | 10.0 | -9.8 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.226 | 2.957 | 0.0100 | 9.17 | 10.0 | -8.3 | 20.0 |
| 3,4-Dichlorobenzotrifluoride | Ave | 0.8081 | 0.8405 | 0.0100 | 10.4 | 10.0 | 4.0 | 20.0 |
| sec-Butylbenzene | Ave | 3.701 | 3.501 | 0.0100 | 9.46 | 10.0 | -5.4 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.734 | 1.586 | 0.6000 | 9.14 | 10.0 | -8.6 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Lab Sample ID: CCVIS 180-224792/2 Calibration Date: 10/04/2017 00:22
 Instrument ID: CHHP5 Calib Start Date: 07/27/2017 00:51
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 07/27/2017 04:24
 Lab File ID: 51003D02.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 | 3.083 | 2.898 | 0.0100 | 9.40 | 10.0 | -6.0 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.780 | 1.602 | 0.5000 | 9.00 | 10.0 | -10.0 | 20.0 |
| 2,4-Dichlorobenzotrifluoride | Ave | 0.7524 | 0.7666 | 0.0100 | 10.2 | 10.0 | 1.9 | 20.0 |
| 2,5-Dichlorobenzotrifluoride | Ave | 0.8127 | 0.8669 | 0.0100 | 10.7 | 10.0 | 6.7 | 20.0 |
| n-Butylbenzene | Ave | 2.514 | 2.277 | 0.0100 | 9.06 | 10.0 | -9.4 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.653 | 1.446 | 0.4000 | 8.75 | 10.0 | -12.5 | 20.0 |
| 1,2-Dibromo-3-Chloropropane | Ave | 0.1835 | 0.1233 | 0.0500 | 6.72 | 10.0 | -32.8* | 20.0 |
| 2,4- & 2,5- & 2,6-Dichlorotoluene | Ave | 1.048 | 1.008 | 0.0100 | 28.8 | 30.0 | -3.8 | 20.0 |
| 2,3- & 3,4- Dichlorotoluene | Ave | 1.084 | 0.9482 | 0.0100 | 17.5 | 20.0 | -12.5 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.7563 | 0.5369 | 0.2000 | 7.10 | 10.0 | -29.0* | 20.0 |
| Hexachlorobutadiene | Ave | 0.2767 | 0.2375 | 0.0100 | 8.58 | 10.0 | -14.2 | 20.0 |
| Naphthalene | Ave | 2.576 | 1.618 | 0.0100 | 6.28 | 10.0 | -37.2* | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 0.6909 | 0.4791 | 0.0100 | 6.94 | 10.0 | -30.6* | 20.0 |
| 2,4,5-Trichlorotoluene | Ave | 0.3284 | 0.2032 | 0.0100 | 6.19 | 10.0 | -38.1* | 20.0 |
| 2,3,6-Trichlorotoluene | Ave | 0.3055 | 0.1917 | 0.0100 | 6.28 | 10.0 | -37.2* | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2406 | 0.2144 | | 8.91 | 10.0 | -10.9 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.2934 | 0.2766 | | 9.43 | 10.0 | -5.7 | 20.0 |
| Toluene-d8 (Surr) | Ave | 3.979 | 4.338 | | 10.9 | 10.0 | 9.0 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 1.437 | 1.433 | | 9.97 | 10.0 | -0.3 | 20.0 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D02.D
 Lims ID: CCVIS
 Client ID:
 Sample Type: CCVIS
 Inject. Date: 04-Oct-2017 00:22:30 ALS Bottle#: 2 Worklist Smp#: 2
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018710-002
 Misc. Info.: CCVIS
 Operator ID: 034635 Instrument ID: CHHP5
 Sublist: chrom-MSVOA_LL_CHHP5*sub29
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 04-Oct-2017 21:10:17 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK005

First Level Reviewer: bungardf

Date: 04-Oct-2017 00:55:51

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.378 | 4.378 | 0.000 | 0 | 166571 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.334 | 7.334 | 0.000 | 97 | 398262 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.431 | 10.431 | 0.000 | 85 | 81838 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.767 | 12.767 | 0.000 | 95 | 115198 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.616 | 6.616 | 0.000 | 93 | 85382 | 50.0 | 44.6 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.981 | 6.981 | 0.000 | 0 | 110172 | 50.0 | 47.1 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.977 | 8.977 | 0.000 | 93 | 355030 | 50.0 | 54.5 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.611 | 11.611 | 0.000 | 87 | 117308 | 50.0 | 49.9 | |
| 11 Dichlorodifluoromethane | 85 | 1.689 | 1.689 | 0.000 | 99 | 123214 | 50.0 | 53.2 | |
| 12 Chloromethane | 50 | 1.823 | 1.823 | 0.000 | 99 | 143761 | 50.0 | 61.8 | |
| 13 Vinyl chloride | 62 | 1.963 | 1.963 | 0.000 | 98 | 135565 | 50.0 | 57.4 | |
| 14 Butadiene | 39 | 1.993 | 1.993 | 0.000 | 92 | 131019 | 50.0 | 61.1 | |
| 15 Bromomethane | 94 | 2.291 | 2.291 | 0.000 | 92 | 55590 | 50.0 | 49.8 | |
| 16 Chloroethane | 64 | 2.461 | 2.461 | 0.000 | 100 | 67252 | 50.0 | 51.8 | |
| 17 Dichlorofluoromethane | 67 | 2.735 | 2.735 | 0.000 | 97 | 171703 | 50.0 | 52.3 | |
| 18 Trichlorofluoromethane | 101 | 2.790 | 2.790 | 0.000 | 95 | 143154 | 50.0 | 49.3 | |
| 20 Ethyl ether | 59 | 3.124 | 3.124 | 0.000 | 93 | 94654 | 50.0 | 50.1 | |
| 21 Acrolein | 56 | 3.301 | 3.301 | 0.000 | 98 | 72146 | 150.0 | 151.7 | |
| 22 1,1-Dichloroethene | 96 | 3.423 | 3.423 | 0.000 | 97 | 106325 | 50.0 | 54.5 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.508 | 3.508 | 0.000 | 91 | 115846 | 50.0 | 54.1 | |
| 24 Acetone | 43 | 3.526 | 3.526 | 0.000 | 98 | 107626 | 100.0 | 103.3 | |
| 25 Iodomethane | 142 | 3.617 | 3.617 | 0.000 | 94 | 151214 | 50.0 | 49.4 | |
| 26 Carbon disulfide | 76 | 3.708 | 3.708 | 0.000 | 99 | 175428 | 50.0 | 41.0 | |
| 28 3-Chloro-1-propene | 76 | 3.994 | 3.994 | 0.000 | 92 | 57196 | 50.0 | 45.4 | |
| 30 Methyl acetate | 43 | 4.031 | 4.031 | 0.000 | 98 | 197400 | 100.0 | 95.7 | |
| 31 Methylene Chloride | 84 | 4.232 | 4.232 | 0.000 | 93 | 113303 | 50.0 | 46.7 | |
| 32 2-Methyl-2-propanol | 59 | 4.493 | 4.493 | 0.000 | 91 | 94463 | 500.0 | 479.5 | |
| 33 Acrylonitrile | 53 | 4.609 | 4.609 | 0.000 | 100 | 487528 | 500.0 | 486.2 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.633 | 4.633 | 0.000 | 98 | 111324 | 50.0 | 50.1 | |
| 35 Methyl tert-butyl ether | 73 | 4.657 | 4.657 | 0.000 | 97 | 253392 | 50.0 | 42.5 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 36 Hexane | 57 | 5.053 | 5.053 | 0.000 | 95 | 145820 | 50.0 | 51.1 | |
| 37 1,1-Dichloroethane | 63 | 5.272 | 5.272 | 0.000 | 96 | 181217 | 50.0 | 46.9 | |
| 38 Vinyl acetate | 43 | 5.315 | 5.315 | 0.000 | 98 | 171984 | 50.0 | 43.8 | |
| 45 cis-1,2-Dichloroethene | 96 | 6.008 | 6.008 | 0.000 | 81 | 117125 | 50.0 | 46.1 | |
| 44 2,2-Dichloropropane | 97 | 6.008 | 6.008 | 0.000 | 65 | 22698 | 50.0 | 46.2 | M |
| 46 2-Butanone (MEK) | 43 | 6.026 | 6.026 | 0.000 | 92 | 138518 | 100.0 | 93.4 | |
| 49 Chlorobromomethane | 128 | 6.294 | 6.294 | 0.000 | 96 | 49889 | 50.0 | 44.2 | |
| 51 Tetrahydrofuran | 42 | 6.306 | 6.306 | 0.000 | 88 | 70164 | 100.0 | 81.3 | |
| 52 Chloroform | 83 | 6.440 | 6.440 | 0.000 | 93 | 179415 | 50.0 | 46.5 | |
| 53 1,1,1-Trichloroethane | 97 | 6.592 | 6.592 | 0.000 | 98 | 141832 | 50.0 | 48.6 | |
| 54 Cyclohexane | 56 | 6.659 | 6.659 | 0.000 | 89 | 173654 | 50.0 | 48.2 | |
| 56 Carbon tetrachloride | 117 | 6.762 | 6.762 | 0.000 | 96 | 118933 | 50.0 | 48.9 | |
| 55 1,1-Dichloropropene | 75 | 6.781 | 6.781 | 0.000 | 95 | 147215 | 50.0 | 46.7 | |
| 57 Isobutyl alcohol | 41 | 6.981 | 6.981 | 0.000 | 90 | 91604 | 1250.0 | 1155.9 | |
| 58 Benzene | 78 | 6.994 | 6.994 | 0.000 | 97 | 447540 | 50.0 | 46.2 | |
| 59 1,2-Dichloroethane | 62 | 7.067 | 7.067 | 0.000 | 97 | 131845 | 50.0 | 46.7 | |
| 62 n-Heptane | 43 | 7.352 | 7.352 | 0.000 | 89 | 116211 | 50.0 | 51.0 | |
| 64 Trichloroethene | 130 | 7.724 | 7.724 | 0.000 | 98 | 109711 | 50.0 | 45.0 | |
| 66 Methylcyclohexane | 83 | 7.961 | 7.961 | 0.000 | 88 | 160467 | 50.0 | 43.5 | |
| 67 1,2-Dichloropropane | 63 | 7.997 | 7.997 | 0.000 | 93 | 94395 | 50.0 | 41.9 | |
| 68 Dibromomethane | 93 | 8.082 | 8.082 | 0.000 | 96 | 53869 | 50.0 | 40.8 | |
| 70 1,4-Dioxane | 88 | 8.082 | 8.082 | 0.000 | 44 | 21077 | 1000.0 | 919.2 | |
| 71 Dichlorobromomethane | 83 | 8.277 | 8.277 | 0.000 | 99 | 99744 | 50.0 | 38.5 | |
| 73 2-Chloroethyl vinyl ether | 63 | 8.575 | 8.575 | 0.000 | 93 | 99449 | 100.0 | 61.3 | |
| 74 cis-1,3-Dichloropropene | 75 | 8.721 | 8.721 | 0.000 | 95 | 121133 | 50.0 | 38.5 | |
| 75 4-Methyl-2-pentanone (MIBK) | 43 | 8.873 | 8.873 | 0.000 | 97 | 199462 | 100.0 | 95.0 | |
| 76 Toluene | 91 | 9.044 | 9.044 | 0.000 | 99 | 435712 | 50.0 | 53.4 | |
| 77 trans-1,3-Dichloropropene | 75 | 9.293 | 9.293 | 0.000 | 93 | 100373 | 50.0 | 45.2 | |
| 78 Ethyl methacrylate | 69 | 9.354 | 9.354 | 0.000 | 90 | 104724 | 50.0 | 39.1 | |
| 79 1,1,2-Trichloroethane | 97 | 9.488 | 9.488 | 0.000 | 91 | 86022 | 50.0 | 50.6 | |
| 80 Tetrachloroethene | 164 | 9.561 | 9.561 | 0.000 | 97 | 84768 | 50.0 | 54.5 | |
| 81 1,3-Dichloropropane | 76 | 9.646 | 9.646 | 0.000 | 92 | 142133 | 50.0 | 45.2 | |
| 82 2-Hexanone | 43 | 9.707 | 9.707 | 0.000 | 96 | 147017 | 100.0 | 91.3 | |
| 84 Chlorodibromomethane | 129 | 9.853 | 9.853 | 0.000 | 92 | 60886 | 50.0 | 42.4 | |
| 85 Ethylene Dibromide | 107 | 9.974 | 9.974 | 0.000 | 97 | 81053 | 50.0 | 46.5 | |
| 86 3-Chlorobenzotrifluoride | 180 | 10.431 | 10.431 | 0.000 | 89 | 162192 | 50.0 | 57.7 | |
| 87 Chlorobenzene | 112 | 10.455 | 10.455 | 0.000 | 94 | 255369 | 50.0 | 48.1 | |
| 88 4-Chlorobenzotrifluoride | 180 | 10.522 | 10.522 | 0.000 | 96 | 152320 | 50.0 | 58.7 | |
| 89 1,1,1,2-Tetrachloroethane | 131 | 10.552 | 10.552 | 0.000 | 92 | 81615 | 50.0 | 48.3 | |
| 90 Ethylbenzene | 106 | 10.558 | 10.558 | 0.000 | 98 | 144895 | 50.0 | 48.9 | |
| 91 m-Xylene & p-Xylene | 106 | 10.692 | 10.692 | 0.000 | 0 | 179890 | 50.0 | 49.6 | |
| 92 o-Xylene | 106 | 11.069 | 11.069 | 0.000 | 96 | 165716 | 50.0 | 48.0 | |
| 93 Styrene | 104 | 11.094 | 11.094 | 0.000 | 95 | 275305 | 50.0 | 47.1 | |
| 94 Bromoform | 173 | 11.270 | 11.270 | 0.000 | 95 | 34884 | 50.0 | 39.1 | |
| 96 2-Chlorobenzotrifluoride | 180 | 11.343 | 11.343 | 0.000 | 97 | 156336 | 50.0 | 58.1 | |
| 97 Isopropylbenzene | 105 | 11.441 | 11.441 | 0.000 | 96 | 433891 | 50.0 | 51.5 | |
| 100 Bromobenzene | 156 | 11.751 | 11.751 | 0.000 | 96 | 98207 | 50.0 | 43.9 | |
| 99 1,1,2,2-Tetrachloroethane | 83 | 11.751 | 11.751 | 0.000 | 81 | 116469 | 50.0 | 46.3 | |
| 102 trans-1,4-Dichloro-2-buten | 53 | 11.787 | 11.787 | 0.000 | 81 | 33264 | 50.0 | 49.3 | |
| 101 1,2,3-Trichloropropane | 110 | 11.799 | 11.799 | 0.000 | 85 | 41652 | 50.0 | 45.1 | |
| 103 N-Propylbenzene | 120 | 11.854 | 11.854 | 0.000 | 98 | 121289 | 50.0 | 47.5 | |
| 104 2-Chlorotoluene | 126 | 11.939 | 11.939 | 0.000 | 97 | 103932 | 50.0 | 47.1 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 105 3-Chlorotoluene | 126 | 12.006 | 12.006 | 0.000 | 97 | 125273 | 50.0 | 52.1 | |
| 106 1,3,5-Trimethylbenzene | 105 | 12.037 | 12.037 | 0.000 | 94 | 355533 | 50.0 | 48.6 | |
| 107 4-Chlorotoluene | 126 | 12.061 | 12.061 | 0.000 | 96 | 113255 | 50.0 | 47.5 | |
| 108 tert-Butylbenzene | 119 | 12.353 | 12.353 | 0.000 | 94 | 275632 | 50.0 | 45.1 | |
| 110 1,2,4-Trimethylbenzene | 105 | 12.408 | 12.408 | 0.000 | 97 | 340653 | 50.0 | 45.8 | |
| 111 1,2-dichloro-4-(trifluorom | 214 | 12.450 | 12.450 | 0.000 | 97 | 96828 | 50.0 | 52.0 | |
| 112 sec-Butylbenzene | 105 | 12.572 | 12.572 | 0.000 | 94 | 403274 | 50.0 | 47.3 | |
| 113 1,3-Dichlorobenzene | 146 | 12.694 | 12.694 | 0.000 | 98 | 182667 | 50.0 | 45.7 | |
| 114 4-Isopropyltoluene | 119 | 12.730 | 12.730 | 0.000 | 97 | 333864 | 50.0 | 47.0 | |
| 115 1,4-Dichlorobenzene | 146 | 12.797 | 12.797 | 0.000 | 96 | 184494 | 50.0 | 45.0 | |
| 116 2,4-Dichloro-1-(trifluorom | 214 | 12.822 | 12.822 | 0.000 | 95 | 88307 | 50.0 | 50.9 | |
| 118 2,5-Dichlorobenzotrifluori | 214 | 12.864 | 12.864 | 0.000 | 0 | 99861 | 50.0 | 53.3 | |
| 120 n-Butylbenzene | 91 | 13.138 | 13.138 | 0.000 | 98 | 262344 | 50.0 | 45.3 | |
| 121 1,2-Dichlorobenzene | 146 | 13.150 | 13.150 | 0.000 | 97 | 166519 | 50.0 | 43.7 | |
| 122 1,2-Dibromo-3-Chloropropan | 75 | 13.947 | 13.947 | 0.000 | 77 | 14209 | 50.0 | 33.6 | |
| 123 2,4- & 2,5- & 2,6- Dichlor | 125 | 14.081 | 14.081 | 0.000 | 0 | 348400 | 150.0 | 144.2 | |
| 125 2,3- & 3,4- Dichlorotoluen | 125 | 14.501 | 14.501 | 0.000 | 0 | 218459 | 100.0 | 87.5 | |
| 126 1,2,4-Trichlorobenzene | 180 | 14.768 | 14.768 | 0.000 | 94 | 61849 | 50.0 | 35.5 | |
| 127 Hexachlorobutadiene | 225 | 14.908 | 14.908 | 0.000 | 96 | 27358 | 50.0 | 42.9 | |
| 128 Naphthalene | 128 | 15.030 | 15.030 | 0.000 | 97 | 186435 | 50.0 | 31.4 | |
| 129 1,2,3-Trichlorobenzene | 180 | 15.255 | 15.255 | 0.000 | 96 | 55196 | 50.0 | 34.7 | |
| 131 2,4,5-Trichlorotoluene | 159 | 16.027 | 16.027 | 0.000 | 0 | 23409 | 50.0 | 30.9 | |
| 130 2,3,6-Trichlorotoluene | 159 | 16.125 | 16.125 | 0.000 | 94 | 22084 | 50.0 | 31.4 | |
| 149 3,4-Dichlorotoluene | 1 | | 0.000 | | | | ND | ND | |
| S 133 Xylenes, Total | 106 | | | | 0 | | 100.0 | 97.6 | |
| S 134 1,2-Dichloroethene, Total | 96 | | | | 0 | | 100.0 | 96.2 | |
| S 135 1,3-Dichloropropene, Total | 1 | | | | 0 | | 100.0 | 83.7 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

Reagents:

| | | | |
|---------------------|--------------------|-----------|-------------|
| voaWEEmix1stR_00011 | Amount Added: 2.00 | Units: uL | |
| voaWKetmix1st_00006 | Amount Added: 2.00 | Units: uL | |
| voaWVA1stRest_00021 | Amount Added: 2.00 | Units: uL | |
| VOA8260VOA2ND_00267 | Amount Added: 2.00 | Units: uL | |
| VOA2CEVE2ND_00008 | Amount Added: 2.00 | Units: uL | |
| voaWAcro1stRe_00021 | Amount Added: 6.00 | Units: uL | |
| VOA8260INT_00074 | Amount Added: 2.00 | Units: uL | Run Reagent |
| VOA8260SURR_00073 | Amount Added: 2.00 | Units: uL | Run Reagent |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D02.D

Injection Date: 04-Oct-2017 00:22:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: CCVIS

Worklist Smp#: 2

Client ID:

Purge Vol: 5.000 mL

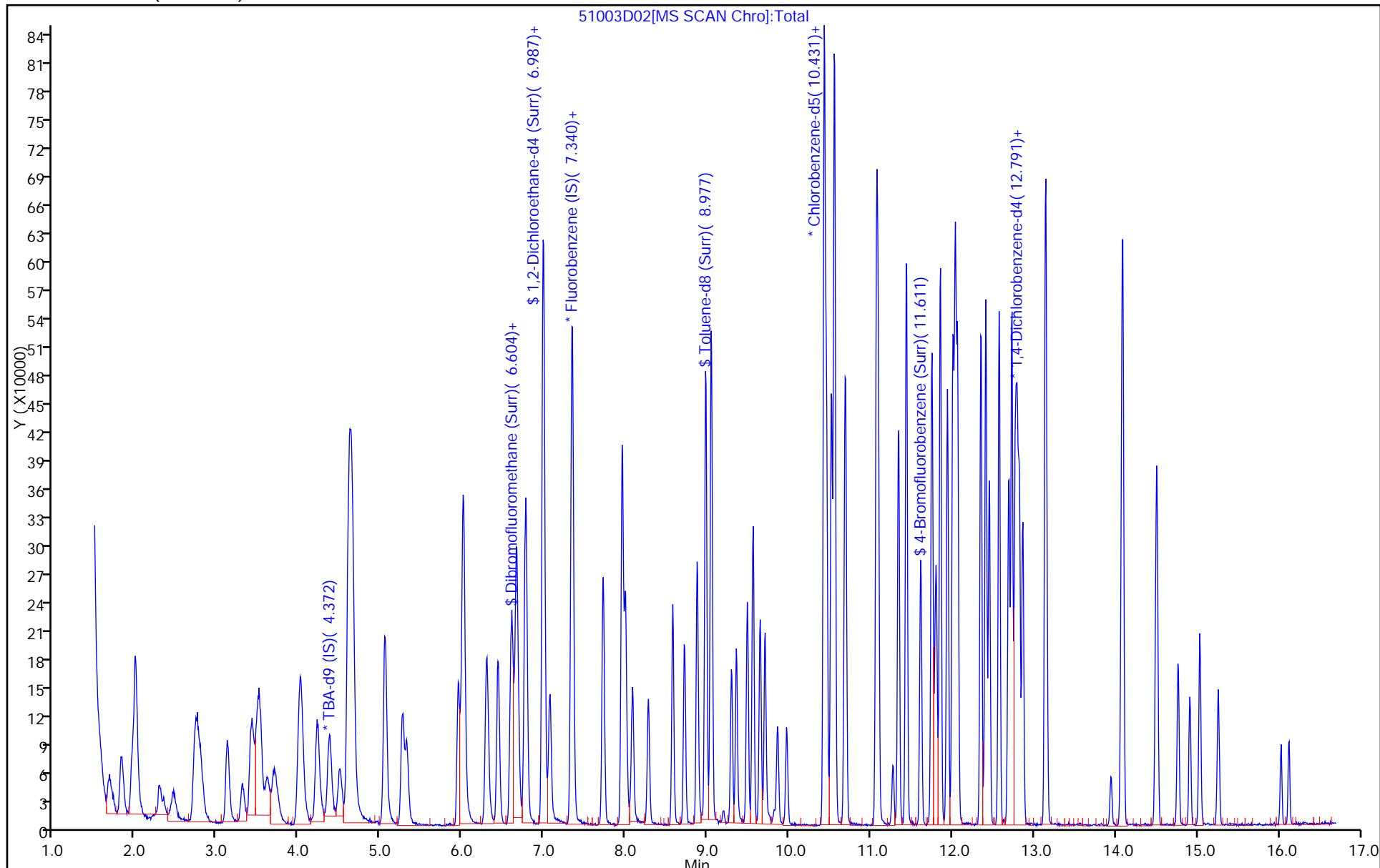
Dil. Factor: 1.0000

ALS Bottle#: 2

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh

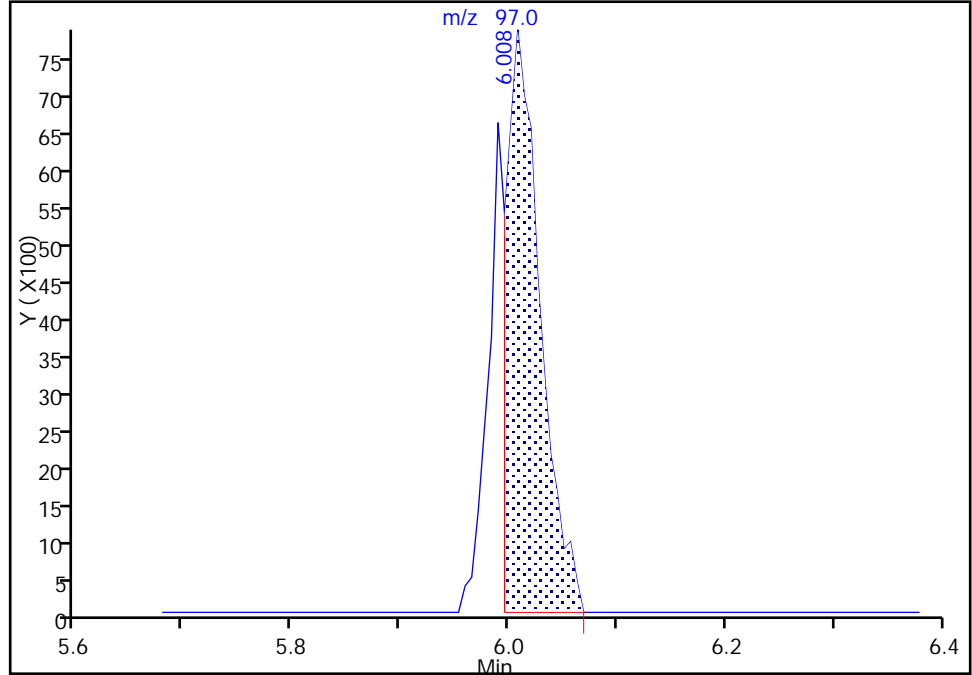
Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D02.D
Injection Date: 04-Oct-2017 00:22:30 Instrument ID: CHHP5
Lims ID: CCVIS
Client ID:
Operator ID: 034635 ALS Bottle#: 2 Worklist Smp#: 2
Purge Vol: 5.000 mL Dil. Factor: 1.0000
Method: MSVOA_LL_CHHP5 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm) Detector: MS SCAN

44 2,2-Dichloropropane, CAS: 594-20-7

Signal: 1

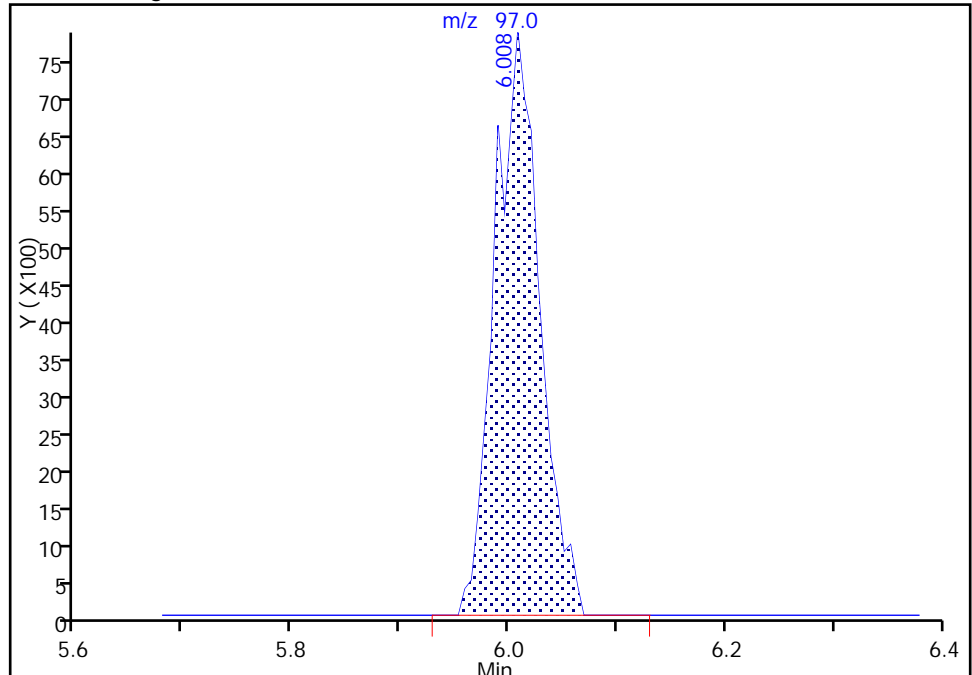
RT: 6.01
Area: 17197
Amount: 34.966668
Amount Units: ng

Processing Integration Results



RT: 6.01
Area: 22698
Amount: 46.151853
Amount Units: ng

Manual Integration Results



Reviewer: bungardf, 04-Oct-2017 00:54:24
Audit Action: Manually Integrated

Audit Reason: Poor chromatography

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Lab Sample ID: CCVIS 180-224919/2 Calibration Date: 10/04/2017 23:29
 Instrument ID: CHHP5 Calib Start Date: 07/27/2017 00:51
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 07/27/2017 04:24
 Lab File ID: 51004D02.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 | 0.2907 | 0.3067 | 0.1000 | 10.5 | 10.0 | 5.5 | 20.0 |
| Chloromethane | Ave | 0.2922 | 0.4296 | 0.1000 | 14.7 | 10.0 | 47.0* | 20.0 |
| Vinyl chloride | Ave | 0.2965 | 0.4118 | 0.1000 | 13.9 | 10.0 | 38.9* | 20.0 |
| 1,3-Butadiene | Ave | 0.2694 | 0.4008 | 0.0100 | 14.9 | 10.0 | 48.8* | 20.0 |
| Bromomethane | Ave | 0.1402 | 0.1596 | 0.0500 | 11.4 | 10.0 | 13.8 | 20.0 |
| Chloroethane | Ave | 0.1630 | 0.2001 | 0.0500 | 12.3 | 10.0 | 22.8* | 20.0 |
| Trichlorofluoromethane | Ave | 0.3643 | 0.4135 | 0.1000 | 11.3 | 10.0 | 13.5 | 20.0 |
| Ethyl ether | Ave | 0.2370 | 0.2491 | 0.0100 | 10.5 | 10.0 | 5.1 | 20.0 |
| Acrolein | Ave | 0.0597 | 0.0650 | 0.0100 | 32.6 | 30.0 | 8.8 | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2448 | 0.2691 | 0.1000 | 11.0 | 10.0 | 9.9 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.2686 | 0.2942 | 0.1000 | 11.0 | 10.0 | 9.5 | 20.0 |
| Acetone | Ave | 0.1308 | 0.1330 | 0.0500 | 20.3 | 20.0 | 1.7 | 20.0 |
| Iodomethane | Ave | 0.3845 | 0.3881 | 0.0100 | 10.1 | 10.0 | 0.9 | 20.0 |
| Carbon disulfide | Ave | 0.5372 | 0.4843 | 0.1000 | 9.01 | 10.0 | -9.9 | 20.0 |
| Allyl chloride | Ave | 0.1582 | 0.1517 | 0.0100 | 9.59 | 10.0 | -4.1 | 20.0 |
| Methyl acetate | Ave | 0.2589 | 0.2776 | 0.1000 | 21.4 | 20.0 | 7.2 | 20.0 |
| Methylene Chloride | Lin2 | | 0.3195 | 0.1000 | 10.6 | 10.0 | 5.8 | 20.0 |
| tert-Butyl alcohol | Ave | 1.183 | 1.183 | 0.0100 | 100 | 100 | 0.0 | 20.0 |
| Acrylonitrile | Ave | 0.1259 | 0.1319 | 0.0100 | 105 | 100 | 4.7 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2789 | 0.2920 | 0.1000 | 10.5 | 10.0 | 4.7 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.7479 | 0.7085 | 0.1000 | 9.47 | 10.0 | -5.3 | 20.0 |
| Hexane | Ave | 0.3580 | 0.3784 | 0.0100 | 10.6 | 10.0 | 5.7 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.4850 | 0.4837 | 0.2000 | 9.97 | 10.0 | -0.3 | 20.0 |
| Vinyl acetate | Ave | 0.4932 | 0.4764 | 0.0100 | 9.66 | 10.0 | -3.4 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.0617 | 0.0627 | 0.0100 | 10.2 | 10.0 | 1.6 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.3190 | 0.3099 | 0.1000 | 9.71 | 10.0 | -2.9 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.1861 | 0.1711 | 0.0500 | 18.4 | 20.0 | -8.1 | 20.0 |
| Bromochloromethane | Ave | 0.1418 | 0.1366 | 0.0100 | 9.64 | 10.0 | -3.6 | 20.0 |
| Tetrahydrofuran | Ave | 0.1084 | 0.0983 | 0.0100 | 18.1 | 20.0 | -9.3 | 20.0 |
| Chloroform | Ave | 0.4843 | 0.4886 | 0.2000 | 10.1 | 10.0 | 0.9 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.3666 | 0.3863 | 0.1000 | 10.5 | 10.0 | 5.4 | 20.0 |
| Cyclohexane | Ave | 0.4524 | 0.4651 | 0.1000 | 10.3 | 10.0 | 2.8 | 20.0 |
| Carbon tetrachloride | Ave | 0.3051 | 0.3037 | 0.1000 | 9.96 | 10.0 | -0.4 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3961 | 0.3779 | 0.0100 | 9.54 | 10.0 | -4.6 | 20.0 |
| Isobutyl alcohol | Ave | 0.0099 | 0.0100 | 0.0100 | 252 | 250 | 0.8 | 20.0 |
| Benzene | Ave | 1.216 | 1.187 | 0.5000 | 9.77 | 10.0 | -2.3 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.3544 | 0.3719 | 0.1000 | 10.5 | 10.0 | 4.9 | 20.0 |
| n-Heptane | Ave | 0.2863 | 0.3036 | 0.0100 | 10.6 | 10.0 | 6.0 | 20.0 |
| Trichloroethene | Ave | 0.3059 | 0.2814 | 0.2000 | 9.20 | 10.0 | -8.0 | 20.0 |
| Methylcyclohexane | Ave | 0.4626 | 0.4125 | 0.1000 | 8.92 | 10.0 | -10.8 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Lab Sample ID: CCVIS 180-224919/2 Calibration Date: 10/04/2017 23:29
 Instrument ID: CHHP5 Calib Start Date: 07/27/2017 00:51
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 07/27/2017 04:24
 Lab File ID: 51004D02.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 | 0.2831 | 0.2643 | 0.1000 | 9.34 | 10.0 | -6.6 | 20.0 |
| Dibromomethane | Ave | 0.1659 | 0.1561 | 0.0100 | 9.41 | 10.0 | -5.9 | 20.0 |
| 1,4-Dioxane | Ave | 0.0029 | 0.0031* | 0.0100 | 213 | 200 | 6.7 | 20.0 |
| Bromodichloromethane | Ave | 0.3256 | 0.2918 | 0.2000 | 8.96 | 10.0 | -10.4 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.2037 | 0.1498 | 0.0100 | 14.7 | 20.0 | -26.5* | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3955 | 0.3380 | 0.2000 | 8.55 | 10.0 | -14.5 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 1.282 | 1.257 | 0.1000 | 19.6 | 20.0 | -2.0 | 20.0 |
| Toluene | Ave | 4.986 | 5.491 | 0.4000 | 11.0 | 10.0 | 10.1 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 1.357 | 1.310 | 0.1000 | 9.66 | 10.0 | -3.4 | 20.0 |
| Ethyl methacrylate | Ave | 1.636 | 1.445 | 0.0100 | 8.83 | 10.0 | -11.7 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 1.039 | 1.051 | 0.1000 | 10.1 | 10.0 | 1.2 | 20.0 |
| Tetrachloroethene | Ave | 0.9508 | 1.011 | 0.2000 | 10.6 | 10.0 | 6.3 | 20.0 |
| 1,3-Dichloropropane | Ave | 1.920 | 1.807 | 0.0100 | 9.41 | 10.0 | -5.9 | 20.0 |
| 2-Hexanone | Ave | 0.9836 | 0.8952 | 0.1000 | 18.2 | 20.0 | -9.0 | 20.0 |
| Dibromochloromethane | Ave | 0.8779 | 0.7979 | 0.1000 | 9.09 | 10.0 | -9.1 | 20.0 |
| 1,2-Dibromoethane (EDB) | Ave | 1.065 | 0.997 | 0.1000 | 9.36 | 10.0 | -6.4 | 20.0 |
| 3-Chlorobenzotrifluoride | Ave | 1.718 | 1.970 | 0.0100 | 11.5 | 10.0 | 14.6 | 20.0 |
| Chlorobenzene | Ave | 3.246 | 3.340 | 0.5000 | 10.3 | 10.0 | 2.9 | 20.0 |
| 4-Chlorobenzotrifluoride | Ave | 1.586 | 1.856 | 0.0100 | 11.7 | 10.0 | 17.0 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 1.032 | 1.047 | 0.0100 | 10.1 | 10.0 | 1.4 | 20.0 |
| Ethylbenzene | Ave | 1.812 | 1.802 | 0.1000 | 9.94 | 10.0 | -0.6 | 20.0 |
| m-Xylene & p-Xylene | Ave | 2.214 | 2.210 | 0.1000 | 9.98 | 10.0 | -0.2 | 20.0 |
| o-Xylene | Ave | 2.110 | 2.031 | 0.3000 | 9.63 | 10.0 | -3.7 | 20.0 |
| Styrene | Ave | 3.571 | 3.547 | 0.3000 | 9.93 | 10.0 | -0.7 | 20.0 |
| Bromoform | Ave | 0.5456 | 0.4449 | 0.1000 | 8.15 | 10.0 | -18.5 | 20.0 |
| 2-Chlorobenzotrifluoride | Ave | 1.644 | 1.946 | 0.0100 | 11.8 | 10.0 | 18.4 | 20.0 |
| Isopropylbenzene | Ave | 5.150 | 5.335 | 0.1000 | 10.4 | 10.0 | 3.6 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 1.538 | 1.586 | 0.3000 | 10.3 | 10.0 | 3.1 | 20.0 |
| Bromobenzene | Ave | 0.9704 | 0.8387 | 0.0100 | 8.64 | 10.0 | -13.6 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2926 | 0.3275 | 0.0100 | 11.2 | 10.0 | 11.9 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.4005 | 0.3754 | 0.0100 | 9.37 | 10.0 | -6.3 | 20.0 |
| N-Propylbenzene | Ave | 1.109 | 1.167 | 0.0100 | 10.5 | 10.0 | 5.2 | 20.0 |
| 2-Chlorotoluene | Ave | 0.9585 | 0.9648 | 0.0100 | 10.1 | 10.0 | 0.7 | 20.0 |
| 3-Chlorotoluene | Ave | 1.043 | 1.149 | 0.0100 | 11.0 | 10.0 | 10.2 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.173 | 3.325 | 0.0100 | 10.5 | 10.0 | 4.8 | 20.0 |
| 4-Chlorotoluene | Ave | 1.035 | 1.075 | 0.0100 | 10.4 | 10.0 | 3.8 | 20.0 |
| tert-Butylbenzene | Ave | 2.653 | 2.541 | 0.0100 | 9.58 | 10.0 | -4.2 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.226 | 3.235 | 0.0100 | 10.0 | 10.0 | 0.3 | 20.0 |
| 3,4-Dichlorobenzotrifluoride | Ave | 0.8081 | 0.8717 | 0.0100 | 10.8 | 10.0 | 7.9 | 20.0 |
| sec-Butylbenzene | Ave | 3.701 | 3.708 | 0.0100 | 10.0 | 10.0 | 0.2 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.734 | 1.626 | 0.6000 | 9.38 | 10.0 | -6.2 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Lab Sample ID: CCVIS 180-224919/2 Calibration Date: 10/04/2017 23:29
 Instrument ID: CHHP5 Calib Start Date: 07/27/2017 00:51
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 07/27/2017 04:24
 Lab File ID: 51004D02.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 | 3.083 | 3.092 | 0.0100 | 10.0 | 10.0 | 0.3 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.780 | 1.725 | 0.5000 | 9.69 | 10.0 | -3.1 | 20.0 |
| 2,4-Dichlorobenzotrifluoride | Ave | 0.7524 | 0.7774 | 0.0100 | 10.3 | 10.0 | 3.3 | 20.0 |
| 2,5-Dichlorobenzotrifluoride | Ave | 0.8127 | 0.8442 | 0.0100 | 10.4 | 10.0 | 3.9 | 20.0 |
| n-Butylbenzene | Ave | 2.514 | 2.516 | 0.0100 | 10.0 | 10.0 | 0.0 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.653 | 1.523 | 0.4000 | 9.22 | 10.0 | -7.8 | 20.0 |
| 1,2-Dibromo-3-Chloropropane | Ave | 0.1835 | 0.1456 | 0.0500 | 7.93 | 10.0 | -20.7* | 20.0 |
| 2,4- & 2,5- & 2,6-Dichlorotoluene | Ave | 1.048 | 1.112 | 0.0100 | 31.8 | 30.0 | 6.1 | 20.0 |
| 2,3- & 3,4- Dichlorotoluene | Ave | 1.084 | 1.080 | 0.0100 | 19.9 | 20.0 | -0.4 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.7563 | 0.5647 | 0.2000 | 7.47 | 10.0 | -25.3* | 20.0 |
| Hexachlorobutadiene | Ave | 0.2767 | 0.2431 | 0.0100 | 8.79 | 10.0 | -12.1 | 20.0 |
| Naphthalene | Ave | 2.576 | 1.802 | 0.0100 | 6.99 | 10.0 | -30.1* | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 0.6909 | 0.4797 | 0.0100 | 6.94 | 10.0 | -30.6* | 20.0 |
| 2,4,5-Trichlorotoluene | Ave | 0.3284 | 0.2219 | 0.0100 | 6.76 | 10.0 | -32.4* | 20.0 |
| 2,3,6-Trichlorotoluene | Ave | 0.3055 | 0.2254 | 0.0100 | 7.38 | 10.0 | -26.2* | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2406 | 0.2409 | | 10.0 | 10.0 | 0.1 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.2934 | 0.3259 | | 11.1 | 10.0 | 11.1 | 20.0 |
| Toluene-d8 (Surr) | Ave | 3.979 | 4.657 | | 11.7 | 10.0 | 17.0 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 1.437 | 1.541 | | 10.7 | 10.0 | 7.2 | 20.0 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D02.D
 Lims ID: CCVIS
 Client ID:
 Sample Type: CCVIS
 Inject. Date: 04-Oct-2017 23:29:30 ALS Bottle#: 2 Worklist Smp#: 2
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018725-002
 Misc. Info.: CCVIS
 Operator ID: 034635 Instrument ID: CHHP5
 Sublist: chrom-MSVOA_LL_CHHP5*sub29
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 08-Oct-2017 20:41:20 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK026

First Level Reviewer: bungardf

Date: 05-Oct-2017 00:03:54

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.374 | 4.374 | 0.000 | 0 | 150069 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.337 | 7.337 | 0.000 | 98 | 334551 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.427 | 10.427 | 0.000 | 86 | 71810 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.769 | 12.769 | 0.000 | 93 | 99164 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.613 | 6.613 | 0.000 | 94 | 80595 | 50.0 | 50.1 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.990 | 6.990 | 0.000 | 0 | 109030 | 50.0 | 55.5 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.979 | 8.979 | 0.000 | 93 | 334418 | 50.0 | 58.5 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.613 | 11.613 | 0.000 | 84 | 110668 | 50.0 | 53.6 | |
| 11 Dichlorodifluoromethane | 85 | 1.679 | 1.679 | 0.000 | 99 | 102599 | 50.0 | 52.7 | |
| 12 Chloromethane | 50 | 1.825 | 1.825 | 0.000 | 100 | 143708 | 50.0 | 73.5 | |
| 13 Vinyl chloride | 62 | 1.959 | 1.959 | 0.000 | 98 | 137781 | 50.0 | 69.4 | |
| 14 Butadiene | 39 | 2.008 | 2.008 | 0.000 | 97 | 134074 | 50.0 | 74.4 | |
| 15 Bromomethane | 94 | 2.300 | 2.300 | 0.000 | 89 | 53381 | 50.0 | 56.9 | |
| 16 Chloroethane | 64 | 2.470 | 2.470 | 0.000 | 99 | 66933 | 50.0 | 61.4 | |
| 17 Dichlorofluoromethane | 67 | 2.744 | 2.744 | 0.000 | 97 | 179484 | 50.0 | 65.1 | |
| 18 Trichlorofluoromethane | 101 | 2.768 | 2.768 | 0.000 | 95 | 138344 | 50.0 | 56.7 | |
| 20 Ethyl ether | 59 | 3.121 | 3.121 | 0.000 | 91 | 83345 | 50.0 | 52.6 | |
| 21 Acrolein | 56 | 3.316 | 3.316 | 0.000 | 98 | 65202 | 150.0 | 163.2 | |
| 22 1,1-Dichloroethene | 96 | 3.413 | 3.413 | 0.000 | 98 | 90030 | 50.0 | 55.0 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.504 | 3.504 | 0.000 | 93 | 98418 | 50.0 | 54.8 | |
| 24 Acetone | 43 | 3.529 | 3.529 | 0.000 | 100 | 88991 | 100.0 | 101.7 | |
| 25 Iodomethane | 142 | 3.608 | 3.608 | 0.000 | 97 | 129830 | 50.0 | 50.5 | |
| 26 Carbon disulfide | 76 | 3.699 | 3.699 | 0.000 | 99 | 162025 | 50.0 | 45.1 | |
| 28 3-Chloro-1-propene | 76 | 3.997 | 3.997 | 0.000 | 92 | 50757 | 50.0 | 48.0 | |
| 30 Methyl acetate | 43 | 4.033 | 4.033 | 0.000 | 98 | 185740 | 100.0 | 107.2 | |
| 31 Methylene Chloride | 84 | 4.222 | 4.222 | 0.000 | 92 | 106903 | 50.0 | 52.9 | |
| 32 2-Methyl-2-propanol | 59 | 4.508 | 4.508 | 0.000 | 90 | 88749 | 500.0 | 500.1 | |
| 33 Acrylonitrile | 53 | 4.605 | 4.605 | 0.000 | 100 | 441194 | 500.0 | 523.7 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.642 | 4.642 | 0.000 | 98 | 97694 | 50.0 | 52.3 | |
| 35 Methyl tert-butyl ether | 73 | 4.660 | 4.660 | 0.000 | 96 | 237025 | 50.0 | 47.4 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| 36 Hexane | 57 | 5.049 | 5.049 | 0.000 | 94 | 126582 | 50.0 | 52.8 | |
| 37 1,1-Dichloroethane | 63 | 5.268 | 5.268 | 0.000 | 96 | 161837 | 50.0 | 49.9 | |
| 38 Vinyl acetate | 43 | 5.323 | 5.323 | 0.000 | 97 | 159377 | 50.0 | 48.3 | |
| 44 2,2-Dichloropropane | 97 | 6.004 | 6.004 | 0.000 | 63 | 20981 | 50.0 | 50.8 | |
| 45 cis-1,2-Dichloroethene | 96 | 6.011 | 6.011 | 0.000 | 81 | 103672 | 50.0 | 48.6 | |
| 46 2-Butanone (MEK) | 43 | 6.023 | 6.023 | 0.000 | 100 | 114490 | 100.0 | 91.9 | |
| 49 Chlorobromomethane | 128 | 6.290 | 6.290 | 0.000 | 97 | 45713 | 50.0 | 48.2 | |
| 51 Tetrahydrofuran | 42 | 6.309 | 6.309 | 0.000 | 89 | 65801 | 100.0 | 90.7 | |
| 52 Chloroform | 83 | 6.436 | 6.436 | 0.000 | 93 | 163468 | 50.0 | 50.4 | |
| 53 1,1,1-Trichloroethane | 97 | 6.595 | 6.595 | 0.000 | 99 | 129241 | 50.0 | 52.7 | |
| 54 Cyclohexane | 56 | 6.662 | 6.662 | 0.000 | 91 | 155594 | 50.0 | 51.4 | |
| 56 Carbon tetrachloride | 117 | 6.759 | 6.759 | 0.000 | 97 | 101603 | 50.0 | 49.8 | |
| 55 1,1-Dichloropropene | 75 | 6.783 | 6.783 | 0.000 | 94 | 126417 | 50.0 | 47.7 | |
| 57 Isobutyl alcohol | 41 | 6.990 | 6.990 | 0.000 | 92 | 83904 | 1250.0 | 1260.4 | |
| 58 Benzene | 78 | 6.996 | 6.996 | 0.000 | 97 | 397249 | 50.0 | 48.8 | |
| 59 1,2-Dichloroethane | 62 | 7.069 | 7.069 | 0.000 | 97 | 124403 | 50.0 | 52.5 | |
| 62 n-Heptane | 43 | 7.349 | 7.349 | 0.000 | 88 | 101557 | 50.0 | 53.0 | |
| 64 Trichloroethene | 130 | 7.720 | 7.720 | 0.000 | 98 | 94152 | 50.0 | 46.0 | |
| 66 Methylcyclohexane | 83 | 7.957 | 7.957 | 0.000 | 89 | 137989 | 50.0 | 44.6 | |
| 67 1,2-Dichloropropane | 63 | 7.994 | 7.994 | 0.000 | 93 | 88419 | 50.0 | 46.7 | |
| 68 Dibromomethane | 93 | 8.079 | 8.079 | 0.000 | 96 | 52223 | 50.0 | 47.1 | |
| 70 1,4-Dioxane | 88 | 8.085 | 8.085 | 0.000 | 46 | 20555 | 1000.0 | 1067.2 | |
| 71 Dichlorobromomethane | 83 | 8.280 | 8.280 | 0.000 | 100 | 97634 | 50.0 | 44.8 | |
| 73 2-Chloroethyl vinyl ether | 63 | 8.578 | 8.578 | 0.000 | 94 | 100237 | 100.0 | 73.5 | |
| 74 cis-1,3-Dichloropropene | 75 | 8.718 | 8.718 | 0.000 | 94 | 113092 | 50.0 | 42.7 | |
| 75 4-Methyl-2-pentanone (MIBK) | 43 | 8.876 | 8.876 | 0.000 | 98 | 180464 | 100.0 | 98.0 | |
| 76 Toluene | 91 | 9.046 | 9.046 | 0.000 | 98 | 394306 | 50.0 | 55.1 | |
| 77 trans-1,3-Dichloropropene | 75 | 9.296 | 9.296 | 0.000 | 95 | 94065 | 50.0 | 48.3 | |
| 78 Ethyl methacrylate | 69 | 9.356 | 9.356 | 0.000 | 90 | 103782 | 50.0 | 44.2 | |
| 79 1,1,2-Trichloroethane | 97 | 9.490 | 9.490 | 0.000 | 92 | 75469 | 50.0 | 50.6 | |
| 80 Tetrachloroethene | 164 | 9.563 | 9.563 | 0.000 | 96 | 72606 | 50.0 | 53.2 | |
| 81 1,3-Dichloropropane | 76 | 9.648 | 9.648 | 0.000 | 91 | 129768 | 50.0 | 47.1 | |
| 82 2-Hexanone | 43 | 9.703 | 9.703 | 0.000 | 97 | 128566 | 100.0 | 91.0 | |
| 84 Chlorodibromomethane | 129 | 9.855 | 9.855 | 0.000 | 91 | 57298 | 50.0 | 45.4 | |
| 85 Ethylene Dibromide | 107 | 9.971 | 9.971 | 0.000 | 98 | 71583 | 50.0 | 46.8 | |
| 86 3-Chlorobenzotrifluoride | 180 | 10.433 | 10.433 | 0.000 | 90 | 141442 | 50.0 | 57.3 | |
| 87 Chlorobenzene | 112 | 10.458 | 10.458 | 0.000 | 94 | 239861 | 50.0 | 51.5 | |
| 88 4-Chlorobenzotrifluoride | 180 | 10.518 | 10.518 | 0.000 | 96 | 133245 | 50.0 | 58.5 | |
| 89 1,1,1,2-Tetrachloroethane | 131 | 10.549 | 10.549 | 0.000 | 92 | 75181 | 50.0 | 50.7 | |
| 90 Ethylbenzene | 106 | 10.555 | 10.555 | 0.000 | 99 | 129375 | 50.0 | 49.7 | |
| 91 m-Xylene & p-Xylene | 106 | 10.689 | 10.689 | 0.000 | 0 | 158697 | 50.0 | 49.9 | |
| 92 o-Xylene | 106 | 11.072 | 11.072 | 0.000 | 96 | 145876 | 50.0 | 48.1 | |
| 93 Styrene | 104 | 11.090 | 11.090 | 0.000 | 94 | 254674 | 50.0 | 49.7 | |
| 94 Bromoform | 173 | 11.273 | 11.273 | 0.000 | 95 | 31945 | 50.0 | 40.8 | |
| 96 2-Chlorobenzotrifluoride | 180 | 11.340 | 11.340 | 0.000 | 95 | 139744 | 50.0 | 59.2 | |
| 97 Isopropylbenzene | 105 | 11.437 | 11.437 | 0.000 | 96 | 383113 | 50.0 | 51.8 | |
| 99 1,1,2,2-Tetrachloroethane | 83 | 11.747 | 11.747 | 0.000 | 84 | 113865 | 50.0 | 51.6 | |
| 100 Bromobenzene | 156 | 11.753 | 11.753 | 0.000 | 95 | 83171 | 50.0 | 43.2 | |
| 102 trans-1,4-Dichloro-2-buten | 53 | 11.784 | 11.784 | 0.000 | 62 | 32474 | 50.0 | 56.0 | |
| 101 1,2,3-Trichloropropane | 110 | 11.808 | 11.808 | 0.000 | 86 | 37227 | 50.0 | 46.9 | |
| 103 N-Propylbenzene | 120 | 11.851 | 11.851 | 0.000 | 98 | 115727 | 50.0 | 52.6 | |
| 104 2-Chlorotoluene | 126 | 11.942 | 11.942 | 0.000 | 96 | 95672 | 50.0 | 50.3 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 105 3-Chlorotoluene | 126 | 12.003 | 12.003 | 0.000 | 97 | 113988 | 50.0 | 55.1 | |
| 106 1,3,5-Trimethylbenzene | 105 | 12.033 | 12.033 | 0.000 | 96 | 329687 | 50.0 | 52.4 | |
| 107 4-Chlorotoluene | 126 | 12.064 | 12.064 | 0.000 | 96 | 106560 | 50.0 | 51.9 | |
| 108 tert-Butylbenzene | 119 | 12.350 | 12.350 | 0.000 | 93 | 251952 | 50.0 | 47.9 | |
| 110 1,2,4-Trimethylbenzene | 105 | 12.410 | 12.410 | 0.000 | 97 | 320762 | 50.0 | 50.1 | |
| 111 1,2-dichloro-4-(trifluorom | 214 | 12.453 | 12.453 | 0.000 | 96 | 86440 | 50.0 | 53.9 | |
| 112 sec-Butylbenzene | 105 | 12.575 | 12.575 | 0.000 | 94 | 367696 | 50.0 | 50.1 | |
| 113 1,3-Dichlorobenzene | 146 | 12.690 | 12.690 | 0.000 | 97 | 161271 | 50.0 | 46.9 | |
| 114 4-Isopropyltoluene | 119 | 12.727 | 12.727 | 0.000 | 97 | 306631 | 50.0 | 50.2 | |
| 115 1,4-Dichlorobenzene | 146 | 12.794 | 12.794 | 0.000 | 95 | 171076 | 50.0 | 48.5 | |
| 116 2,4-Dichloro-1-(trifluorom | 214 | 12.818 | 12.818 | 0.000 | 95 | 77085 | 50.0 | 51.7 | |
| 118 2,5-Dichlorobenzotrifluori | 214 | 12.867 | 12.867 | 0.000 | 0 | 83713 | 50.0 | 51.9 | |
| 120 n-Butylbenzene | 91 | 13.140 | 13.140 | 0.000 | 98 | 249535 | 50.0 | 50.0 | |
| 121 1,2-Dichlorobenzene | 146 | 13.153 | 13.153 | 0.000 | 95 | 151030 | 50.0 | 46.1 | |
| 122 1,2-Dibromo-3-Chloropropan | 75 | 13.937 | 13.937 | 0.000 | 79 | 14436 | 50.0 | 39.7 | |
| 123 2,4- & 2,5- & 2,6- Dichlor | 125 | 14.083 | 14.083 | 0.000 | 0 | 330885 | 150.0 | 159.1 | |
| 125 2,3- & 3,4- Dichlorotoluen | 125 | 14.503 | 14.503 | 0.000 | 0 | 214195 | 100.0 | 99.6 | |
| 126 1,2,4-Trichlorobenzene | 180 | 14.765 | 14.765 | 0.000 | 92 | 55998 | 50.0 | 37.3 | |
| 127 Hexachlorobutadiene | 225 | 14.905 | 14.905 | 0.000 | 95 | 24110 | 50.0 | 43.9 | |
| 128 Naphthalene | 128 | 15.026 | 15.026 | 0.000 | 97 | 178669 | 50.0 | 35.0 | |
| 129 1,2,3-Trichlorobenzene | 180 | 15.257 | 15.257 | 0.000 | 93 | 47565 | 50.0 | 34.7 | |
| 131 2,4,5-Trichlorotoluene | 159 | 16.024 | 16.024 | 0.000 | 0 | 22008 | 50.0 | 33.8 | |
| 130 2,3,6-Trichlorotoluene | 159 | 16.121 | 16.121 | 0.000 | 95 | 22350 | 50.0 | 36.9 | |
| 149 3,4-Dichlorotoluene | 1 | | 0.000 | | | | ND | ND | |
| S 133 Xylenes, Total | 106 | | | | 0 | | 100.0 | 98.1 | |
| S 134 1,2-Dichloroethene, Total | 96 | | | | 0 | | 100.0 | 100.9 | |
| S 135 1,3-Dichloropropene, Total | 1 | | | | 0 | | 100.0 | 91.0 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Reagents:

| | | | |
|---------------------|--------------------|-----------|-------------|
| voaWEEmix1stR_00011 | Amount Added: 2.00 | Units: uL | |
| voaWKetmix1st_00006 | Amount Added: 2.00 | Units: uL | |
| voaWVA1stRest_00021 | Amount Added: 2.00 | Units: uL | |
| VOA8260VOA2ND_00267 | Amount Added: 2.00 | Units: uL | |
| VOA2CEVE2ND_00008 | Amount Added: 2.00 | Units: uL | |
| voaWAcro1stRe_00021 | Amount Added: 6.00 | Units: uL | |
| VOA8260INT_00074 | Amount Added: 2.00 | Units: uL | Run Reagent |
| VOA8260SURR_00073 | Amount Added: 2.00 | Units: uL | Run Reagent |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D02.D

Injection Date: 04-Oct-2017 23:29:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: CCVIS

Worklist Smp#: 2

Client ID:

Purge Vol: 5.000 mL

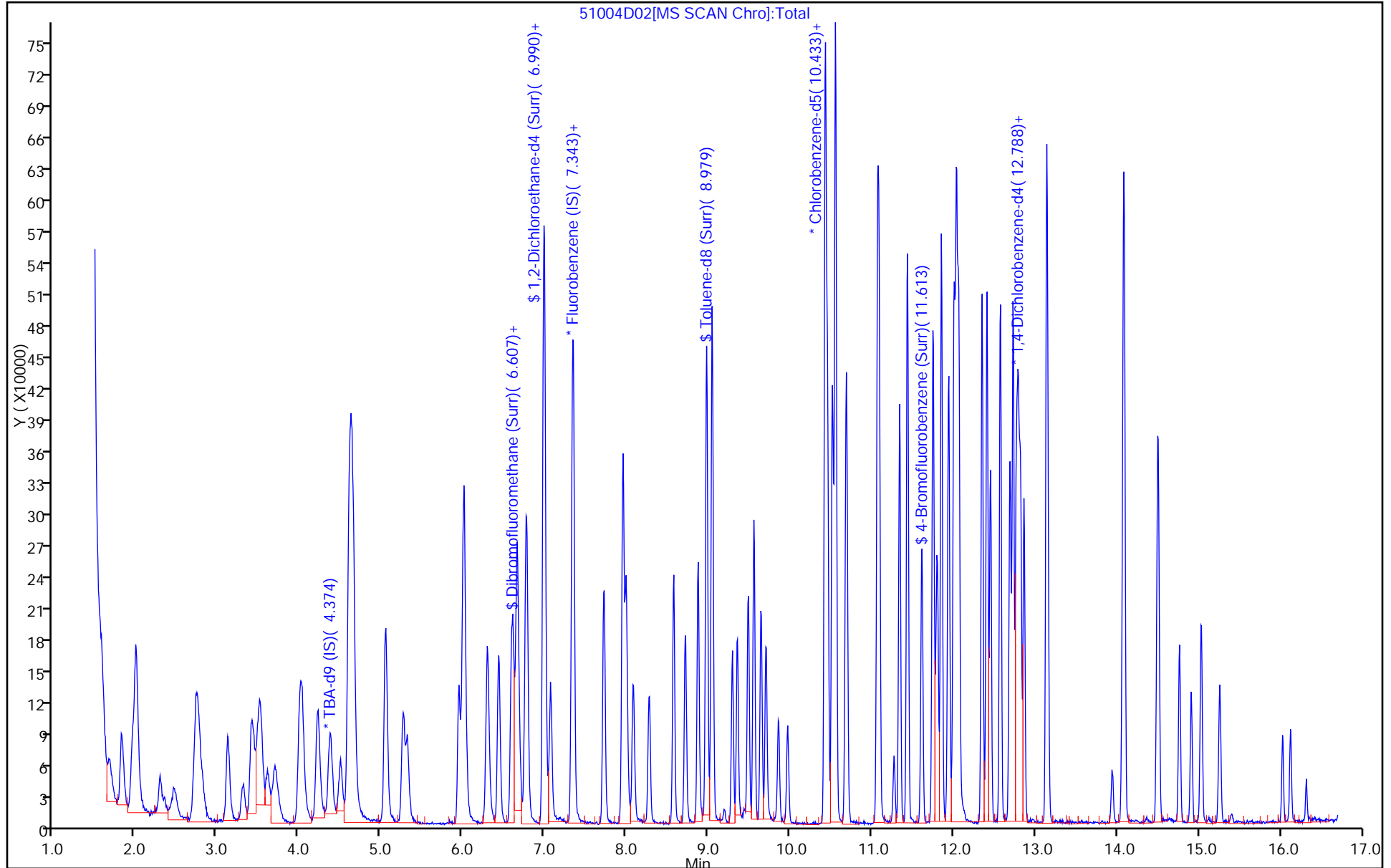
Dil. Factor: 1.0000

ALS Bottle#: 2

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D01.D
 Lims ID: BFB
 Client ID:
 Sample Type: BFB
 Inject. Date: 27-Jul-2017 00:22:30 ALS Bottle#: 1 Worklist Smp#: 1
 Injection Vol: 5.0 mL Dil. Factor: 1.0000
 Sample Info: 180-0017756-001
 Misc. Info.: BFB
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 28-Jul-2017 01:04:43 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK029

First Level Reviewer: bungardf Date: 27-Jul-2017 05:09:11

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

| | | | | | | | | | |
|-----------|----|-------|-------|-------|---|-------|----|----|--|
| \$ 10 BFB | 95 | 8.334 | 8.334 | 0.000 | 0 | 79656 | NR | NR | |
|-----------|----|-------|-------|-------|---|-------|----|----|--|

QC Flag Legend

Processing Flags

NR - Missing Quant Standard

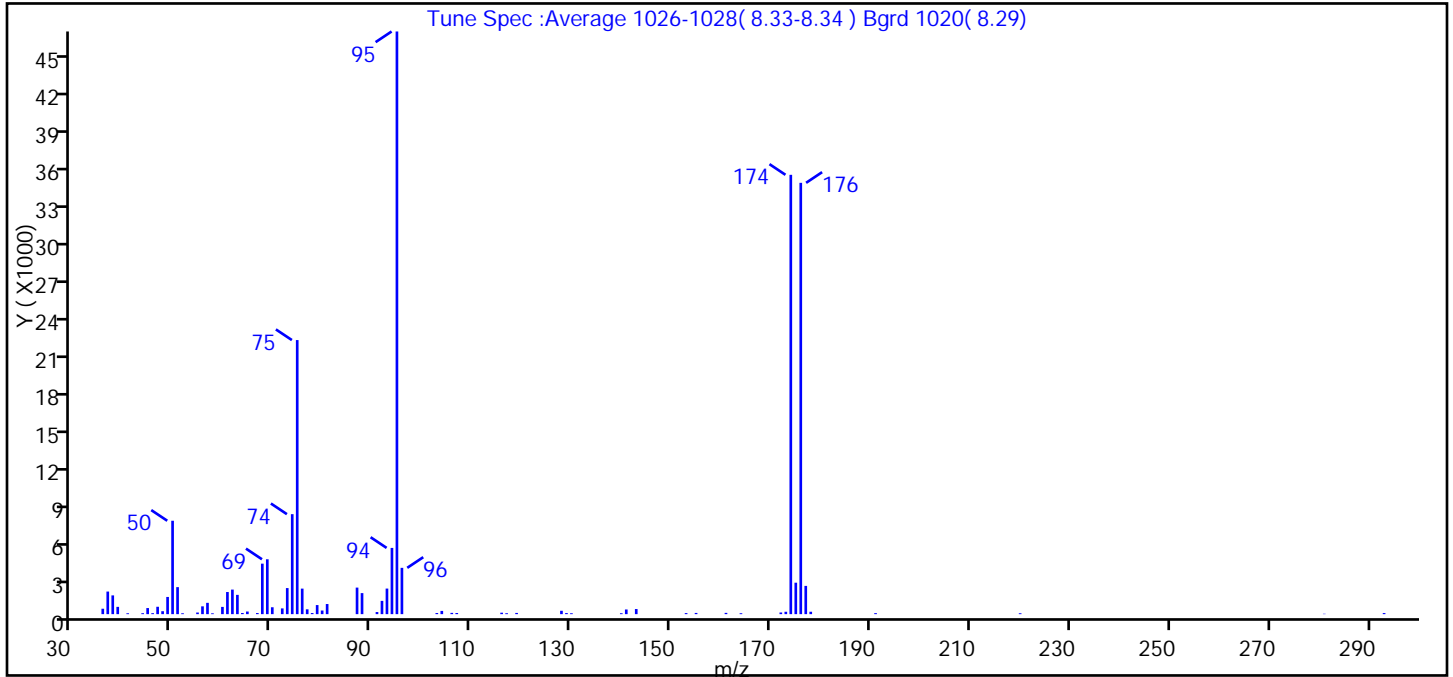
Reagents:

VOABFB25_00090 Amount Added: 1.00 Units: uL

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D01.D
 Injection Date: 27-Jul-2017 00:22:30 Instrument ID: CHHP5
 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_CHHP5 Limit Group: VOA 8260C ICAL
 Tune Method: BFB Method 8260

\$ 10 BFB



| m/z | Ion Abundance Criteria | % Relative Abundance |
|-----|--|----------------------|
| 95 | Base peak, 100% relative abundance | 100.0 |
| 50 | 15 to 40% of m/z 95 | 16.0 |
| 75 | 30 to 60% of m/z 95 | 47.0 |
| 96 | 5 to 9% of m/z 95 | 7.9 |
| 173 | Less than 2% of m/z 174 | 0.4 (0.5) |
| 174 | 50 to 120% of m/z 95 | 75.4 |
| 175 | 5 to 9% of m/z 174 | 5.4 (7.2) |
| 176 | Greater than 95% but less than 101% of m/z 174 | 74.0 (98.2) |
| 177 | 5 to 9% of m/z 176 | 4.8 (6.5) |

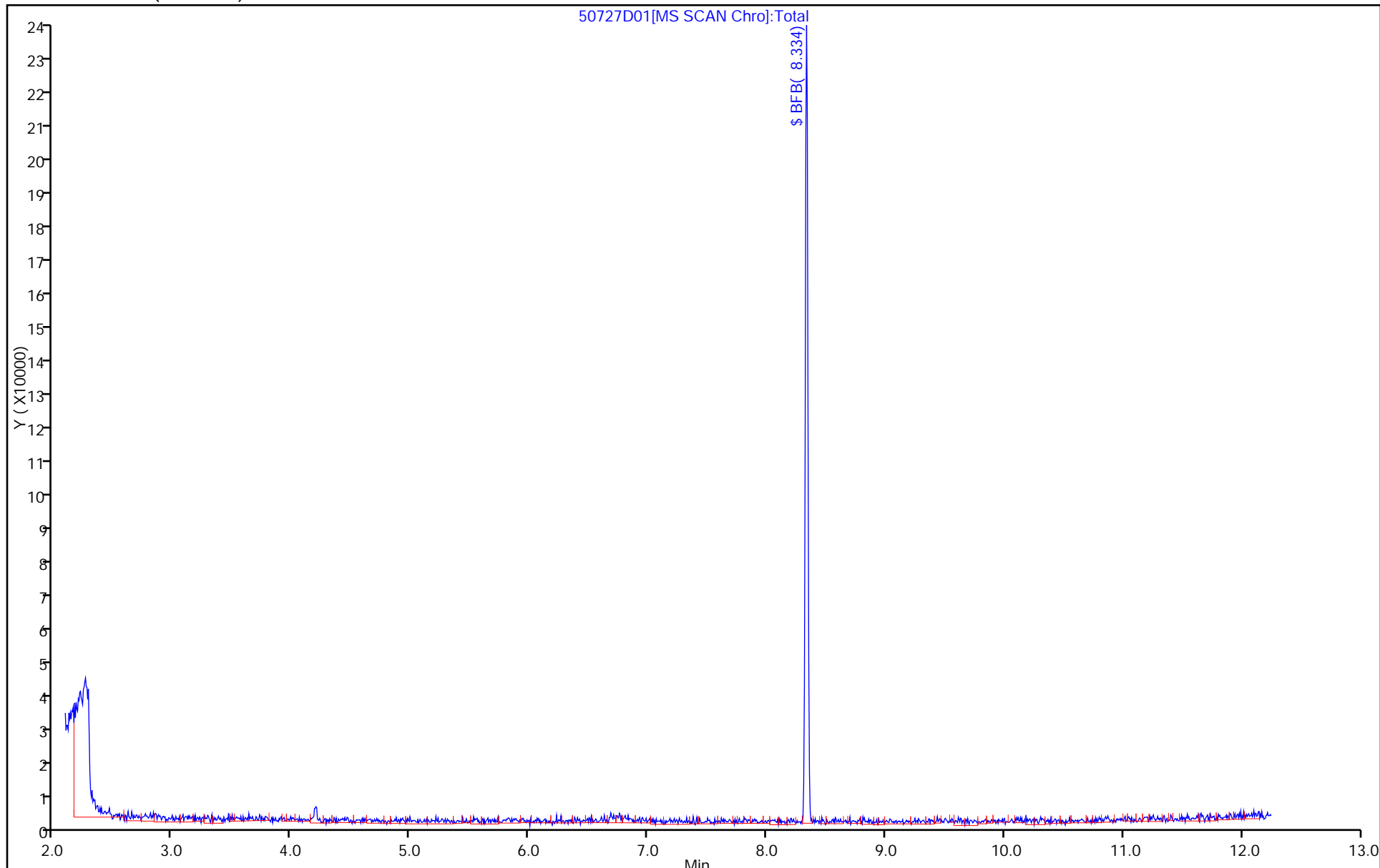
Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D01.D\MSVOA_LL_CHHP5.rsl\spec
 Injection Date: 27-Jul-2017 00:22:30
 Spectrum: Tune Spec :Average 1026-1028(8.33-8.34) Bgrd 1020(8.29)
 Base Peak: 95.00
 Minimum % Base Peak: 0
 Number of Points: 74

| m/z | Y | m/z | Y | m/z | Y | m/z | Y |
|-------|------|-------|-------|--------|-------|--------|-------|
| 36.00 | 433 | 61.00 | 1769 | 87.00 | 2123 | 141.00 | 374 |
| 37.00 | 1806 | 62.00 | 1963 | 88.00 | 1682 | 143.00 | 408 |
| 38.00 | 1500 | 63.00 | 1542 | 91.00 | 169 | 153.00 | 84 |
| 39.00 | 582 | 64.00 | 92 | 92.00 | 1061 | 155.00 | 97 |
| 41.00 | 70 | 65.00 | 209 | 93.00 | 2045 | 161.00 | 102 |
| 44.00 | 76 | 67.00 | 88 | 94.00 | 5297 | 164.00 | 73 |
| 45.00 | 487 | 68.00 | 4038 | 95.00 | 46600 | 172.00 | 132 |
| 46.00 | 79 | 69.00 | 4388 | 96.00 | 3703 | 173.00 | 191 |
| 47.00 | 590 | 70.00 | 551 | 103.00 | 90 | 174.00 | 35136 |
| 48.00 | 235 | 72.00 | 459 | 104.00 | 258 | 175.00 | 2515 |
| 49.00 | 1375 | 73.00 | 2085 | 106.00 | 102 | 176.00 | 34496 |
| 50.00 | 7469 | 74.00 | 7996 | 107.00 | 90 | 177.00 | 2259 |
| 51.00 | 2160 | 75.00 | 21920 | 116.00 | 116 | 178.00 | 192 |
| 52.00 | 70 | 76.00 | 2042 | 117.00 | 73 | 191.00 | 80 |
| 55.00 | 130 | 77.00 | 386 | 119.00 | 97 | 220.00 | 71 |
| 56.00 | 624 | 78.00 | 89 | 128.00 | 269 | 281.00 | 30 |
| 57.00 | 904 | 79.00 | 726 | 129.00 | 86 | 293.00 | 87 |
| 58.00 | 67 | 80.00 | 290 | 130.00 | 72 | | |
| 60.00 | 579 | 81.00 | 809 | 140.00 | 72 | | |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D01.D
Injection Date: 27-Jul-2017 00:22:30 Instrument ID: CHHP5
Lims ID: BFB
Client ID:
Injection Vol: 5.0 mL Dil. Factor: 1.0000
Method: MSVOA_LL_CHHP5 Limit Group: VOA 8260C ICAL
Column: DB-624 (0.18 mm)

Operator ID: 034635
Worklist Smp#: 1
ALS Bottle#: 1



TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D01.D
 Lims ID: BFB
 Client ID:
 Sample Type: BFB
 Inject. Date: 03-Oct-2017 23:49:30 ALS Bottle#: 1 Worklist Smp#: 1
 Injection Vol: 5.0 mL Dil. Factor: 1.0000
 Sample Info: 180-0018710-001
 Misc. Info.: BFB
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 04-Oct-2017 21:10:15 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK005

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|-----------|-----|-----------|---------------|---------------|---|----------|------------|--------------|-------|
| \$ 10 BFB | 95 | 8.334 | 8.334 | 0.000 | 0 | 49616 | NR | NR | |

QC Flag Legend

Processing Flags
 NR - Missing Quant Standard

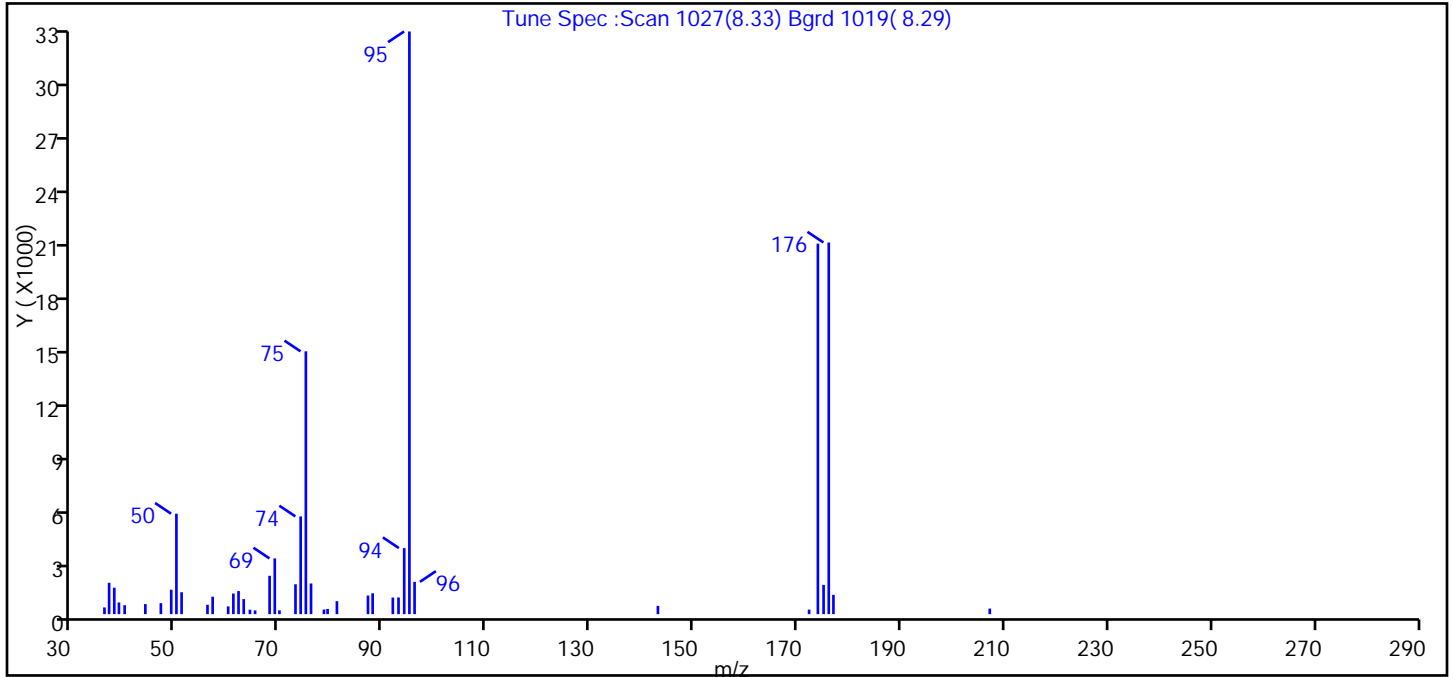
Reagents:

VOABFB25_00093 Amount Added: 1.00 Units: uL

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D01.D
 Injection Date: 03-Oct-2017 23:49:30 Instrument ID: CHHP5
 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_CHHP5 Limit Group: VOA 8260C ICAL
 Tune Method: BFB Method 8260

\$ 10 BFB



| m/z | Ion Abundance Criteria | % Relative Abundance |
|-----|--|----------------------|
| 95 | Base peak, 100% relative abundance | 100.0 |
| 50 | 15 to 40% of m/z 95 | 17.2 |
| 75 | 30 to 60% of m/z 95 | 45.1 |
| 96 | 5 to 9% of m/z 95 | 5.5 |
| 173 | Less than 2% of m/z 174 | 0.0 (0.0) |
| 174 | 50 to 120% of m/z 95 | 63.6 |
| 175 | 5 to 9% of m/z 174 | 5.0 (7.9) |
| 176 | Greater than 95% but less than 101% of m/z 174 | 63.8 (100.3) |
| 177 | 5 to 9% of m/z 176 | 3.3 (5.2) |

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D01.D\MSVOA_LL_CHHP5.rsl\spec
Injection Date: 03-Oct-2017 23:49:30
Spectrum: Tune Spec :Scan 1027(8.33) Bgrd 1019(8.29)
Base Peak: 95.00
Minimum % Base Peak: 0
Number of Points: 43

| m/z | Y | m/z | Y | m/z | Y | m/z | Y |
|-------|------|-------|------|-------|-------|--------|-------|
| 36.10 | 372 | 57.00 | 969 | 74.00 | 5433 | 95.00 | 32384 |
| 37.00 | 1742 | 60.00 | 430 | 75.00 | 14607 | 96.00 | 1795 |
| 38.00 | 1469 | 61.00 | 1147 | 76.00 | 1704 | 143.00 | 460 |
| 38.90 | 649 | 62.00 | 1282 | 78.50 | 256 | 172.20 | 252 |
| 40.00 | 496 | 63.00 | 838 | 79.20 | 289 | 173.90 | 20592 |
| 44.00 | 560 | 64.20 | 246 | 81.00 | 723 | 175.00 | 1627 |
| 47.00 | 612 | 65.20 | 209 | 87.00 | 1035 | 176.00 | 20656 |
| 49.00 | 1357 | 68.00 | 2131 | 87.90 | 1158 | 176.90 | 1071 |
| 50.00 | 5584 | 69.00 | 3098 | 91.80 | 923 | 207.10 | 310 |
| 51.00 | 1217 | 69.90 | 211 | 92.90 | 927 | 281.00 | 1 |
| 56.00 | 521 | 73.00 | 1660 | 94.00 | 3671 | | |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D01.D

Injection Date: 03-Oct-2017 23:49:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: BFB

Worklist Smp#: 1

Client ID:

Injection Vol: 5.0 mL

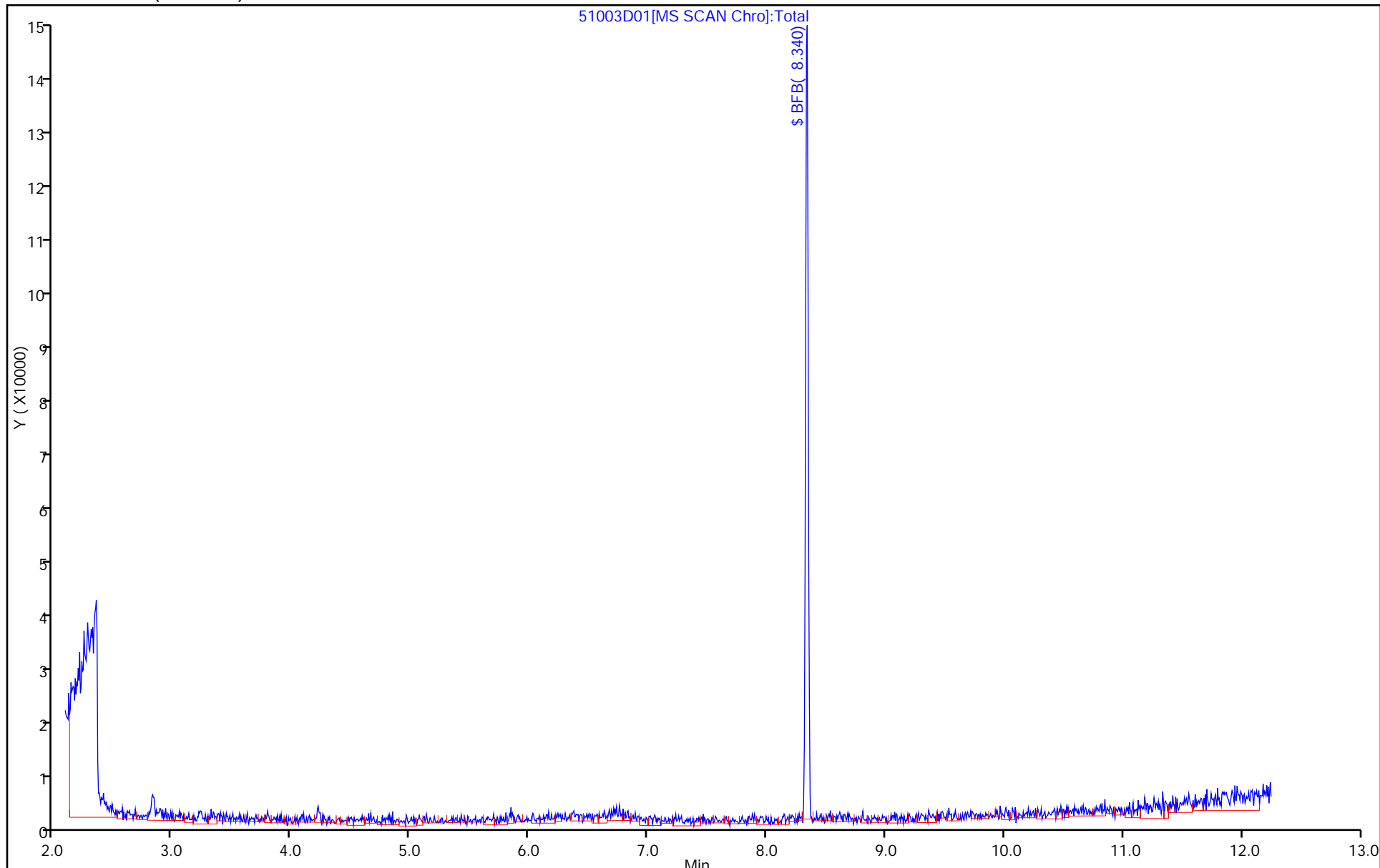
Dil. Factor: 1.0000

ALS Bottle#: 1

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D01.D
 Lims ID: BFB
 Client ID:
 Sample Type: BFB
 Inject. Date: 04-Oct-2017 22:24:30 ALS Bottle#: 1 Worklist Smp#: 1
 Injection Vol: 5.0 mL Dil. Factor: 1.0000
 Sample Info: 180-0018725-001
 Misc. Info.: BFB
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 08-Oct-2017 20:41:18 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK026

First Level Reviewer: bungardf Date: 05-Oct-2017 04:03:09

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

| | | | | | | | | | |
|-----------|----|-------|-------|-------|---|-------|----|----|--|
| \$ 10 BFB | 95 | 8.336 | 8.336 | 0.000 | 0 | 73323 | NR | NR | |
|-----------|----|-------|-------|-------|---|-------|----|----|--|

QC Flag Legend

Processing Flags

NR - Missing Quant Standard

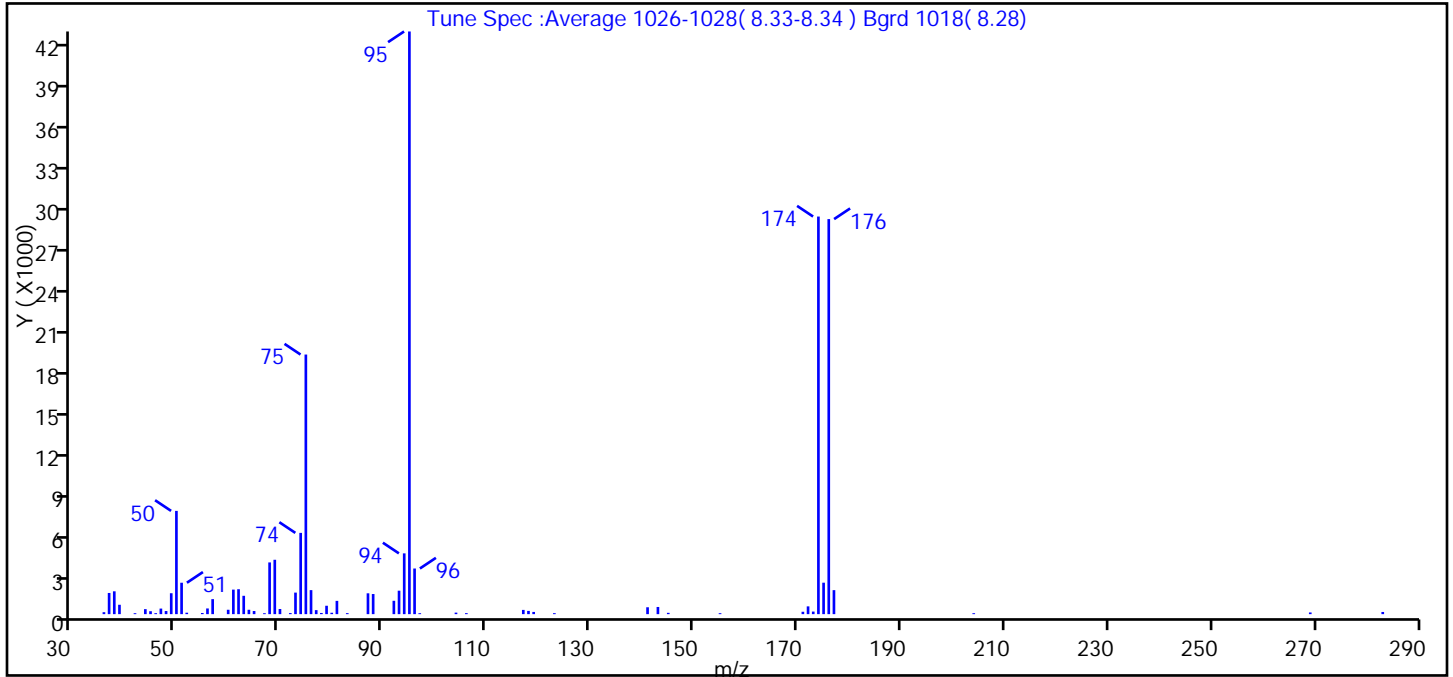
Reagents:

VOABFB25_00093 Amount Added: 1.00 Units: uL

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D01.D
 Injection Date: 04-Oct-2017 22:24:30 Instrument ID: CHHP5
 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_CHHP5 Limit Group: VOA 8260C ICAL
 Tune Method: BFB Method 8260

\$ 10 BFB



| m/z | Ion Abundance Criteria | % Relative Abundance |
|-----|--|----------------------|
| 95 | Base peak, 100% relative abundance | 100.0 |
| 50 | 15 to 40% of m/z 95 | 17.7 |
| 75 | 30 to 60% of m/z 95 | 44.6 |
| 96 | 5 to 9% of m/z 95 | 7.8 |
| 173 | Less than 2% of m/z 174 | 0.4 (0.6) |
| 174 | 50 to 120% of m/z 95 | 68.2 |
| 175 | 5 to 9% of m/z 174 | 5.4 (7.9) |
| 176 | Greater than 95% but less than 101% of m/z 174 | 67.8 (99.4) |
| 177 | 5 to 9% of m/z 176 | 4.1 (6.1) |

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D01.D\MSVOA_LL_CHHP5.rsl\spec
 Injection Date: 04-Oct-2017 22:24:30
 Spectrum: Tune Spec :Average 1026-1028(8.33-8.34) Bgrd 1018(8.28)
 Base Peak: 95.00
 Minimum % Base Peak: 0
 Number of Points: 66

| m/z | Y | m/z | Y | m/z | Y | m/z | Y |
|-------|------|-------|-------|--------|-------|--------|-------|
| 36.00 | 141 | 60.00 | 317 | 79.00 | 606 | 123.00 | 71 |
| 37.00 | 1529 | 61.00 | 1772 | 80.00 | 100 | 141.00 | 497 |
| 38.00 | 1649 | 62.00 | 1800 | 81.00 | 958 | 143.00 | 518 |
| 39.00 | 677 | 63.00 | 1325 | 83.00 | 73 | 145.00 | 95 |
| 42.00 | 68 | 64.00 | 316 | 87.00 | 1505 | 155.00 | 69 |
| 44.00 | 361 | 65.00 | 226 | 88.00 | 1453 | 171.00 | 172 |
| 45.00 | 211 | 67.00 | 67 | 92.00 | 965 | 172.00 | 562 |
| 46.00 | 71 | 68.00 | 3736 | 93.00 | 1694 | 173.00 | 186 |
| 47.00 | 402 | 69.00 | 3927 | 94.00 | 4390 | 174.00 | 28720 |
| 48.00 | 225 | 70.00 | 371 | 95.00 | 42088 | 175.00 | 2272 |
| 49.00 | 1511 | 72.00 | 76 | 96.00 | 3296 | 176.00 | 28536 |
| 50.00 | 7458 | 73.00 | 1556 | 97.00 | 68 | 177.00 | 1733 |
| 51.00 | 2269 | 74.00 | 5870 | 104.00 | 110 | 204.00 | 68 |
| 52.00 | 104 | 75.00 | 18752 | 106.00 | 72 | 269.00 | 122 |
| 55.00 | 83 | 76.00 | 1735 | 117.00 | 306 | 283.00 | 157 |
| 56.00 | 411 | 77.00 | 290 | 118.00 | 233 | | |
| 57.00 | 1084 | 78.00 | 85 | 119.00 | 160 | | |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D01.D

Injection Date: 04-Oct-2017 22:24:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: BFB

Worklist Smp#: 1

Client ID:

Injection Vol: 5.0 mL

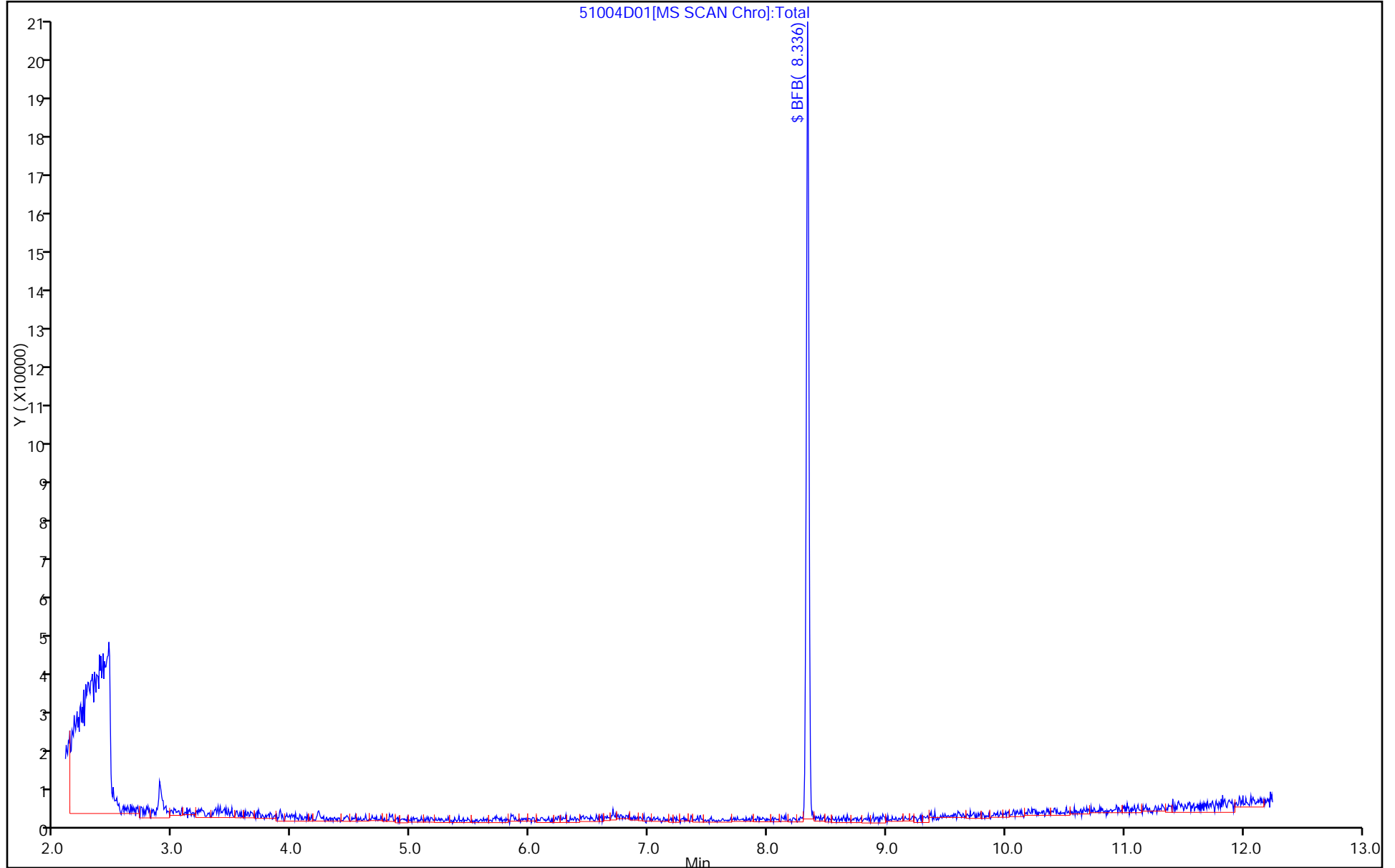
Dil. Factor: 1.0000

ALS Bottle#: 1

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Client Sample ID: _____ Lab Sample ID: MB 180-224792/6
 Matrix: Water Lab File ID: 51003D06.D
 Analysis Method: 8260C Date Collected: _____
 Sample wt/vol: 5 (mL) Date Analyzed: 10/04/2017 02:21
 Soil Aliquot Vol: _____ Dilution Factor: 1
 Soil Extract Vol.: _____ GC Column: DB-624 ID: 0.18 (mm)
 % Moisture: _____ Level: (low/med) Low
 Analysis Batch No.: 224792 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|------------|-----------------------------|--------|---|-----|------|
| 74-87-3 | Chloromethane | 1.0 | U | 1.0 | 0.38 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 1.0 | 0.17 |
| 74-83-9 | Bromomethane | 1.0 | U | 1.0 | 0.59 |
| 75-00-3 | Chloroethane | 1.0 | U | 1.0 | 0.58 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.32 |
| 67-64-1 | Acetone | 5.0 | U | 5.0 | 3.1 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 1.0 | 0.53 |
| 75-09-2 | Methylene Chloride | 1.0 | U | 1.0 | 0.94 |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.20 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.34 |
| 156-59-2 | cis-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.30 |
| 74-97-5 | Bromochloromethane | 1.0 | U | 1.0 | 0.36 |
| 78-93-3 | 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 |
| 67-66-3 | Chloroform | 1.0 | U | 1.0 | 0.27 |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.27 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 1.0 | 0.56 |
| 71-43-2 | Benzene | 1.0 | U | 1.0 | 0.18 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.24 |
| 79-01-6 | Trichloroethene | 1.0 | U | 1.0 | 0.20 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.35 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 1.0 | 0.57 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.32 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 5.0 | U | 5.0 | 2.2 |
| 108-88-3 | Toluene | 1.0 | U | 1.0 | 0.16 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.22 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.31 |
| 127-18-4 | Tetrachloroethene | 1.0 | U | 1.0 | 0.24 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 5.0 | 2.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 1.0 | 0.44 |
| 106-93-4 | 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.51 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 1.0 | 0.15 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.49 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 1.0 | 0.25 |
| 1330-20-7 | Xylenes, Total | 2.0 | U | 2.0 | 0.27 |
| 100-42-5 | Styrene | 1.0 | U | 1.0 | 0.22 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Client Sample ID: _____ Lab Sample ID: MB 180-224792/6
 Matrix: Water Lab File ID: 51003D06.D
 Analysis Method: 8260C Date Collected: _____
 Sample wt/vol: 5 (mL) Date Analyzed: 10/04/2017 02:21
 Soil Aliquot Vol: _____ Dilution Factor: 1
 Soil Extract Vol.: _____ GC Column: DB-624 ID: 0.18 (mm)
 % Moisture: _____ Level: (low/med) Low
 Analysis Batch No.: 224792 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|----------|---------------------------|--------|---|-----|------|
| 75-25-2 | Bromoform | 1.0 | U | 1.0 | 0.76 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.37 |
| 107-13-1 | Acrylonitrile | 20 | U | 20 | 3.3 |
| 123-91-1 | 1,4-Dioxane | 200 | U | 200 | 16 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|------------|------------------------------|------|---|--------|
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 107 | | 65-121 |
| 2037-26-5 | Toluene-d8 (Surr) | 98 | | 73-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 93 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 94 | | 73-120 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D06.D
 Lims ID: MB
 Client ID:
 Sample Type: MB
 Inject. Date: 04-Oct-2017 02:21:30 ALS Bottle#: 6 Worklist Smp#: 6
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018710-006
 Misc. Info.: MB
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 04-Oct-2017 21:10:21 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK005

First Level Reviewer: bungardf

Date: 04-Oct-2017 02:42:56

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.350 | 4.347 | 0.003 | 0 | 193789 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.337 | 7.334 | 0.003 | 99 | 384752 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.427 | 10.431 | -0.004 | 86 | 86410 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.770 | 12.773 | -0.003 | 97 | 122049 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.619 | 6.616 | 0.003 | 91 | 86724 | 50.0 | 46.8 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.984 | 6.981 | 0.003 | 0 | 120939 | 50.0 | 53.6 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.980 | 8.977 | 0.003 | 93 | 336841 | 50.0 | 49.0 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.608 | 11.611 | -0.003 | 86 | 114967 | 50.0 | 46.3 | |
| 11 Dichlorodifluoromethane | 85 | | 1.689 | | | | | ND | |
| 12 Chloromethane | 50 | | 1.823 | | | | | ND | |
| 13 Vinyl chloride | 62 | | 1.963 | | | | | ND | |
| 14 Butadiene | 39 | | 1.993 | | | | | ND | |
| 15 Bromomethane | 94 | | 2.291 | | | | | ND | |
| 16 Chloroethane | 64 | | 2.461 | | | | | ND | |
| 17 Dichlorofluoromethane | 67 | | 2.735 | | | | | ND | |
| 18 Trichlorofluoromethane | 101 | | 2.790 | | | | | ND | |
| 19 Ethanol | 45 | | 2.821 | | | | | ND | |
| 20 Ethyl ether | 59 | | 3.124 | | | | | ND | |
| 21 Acrolein | 56 | | 3.301 | | | | | ND | |
| 22 1,1-Dichloroethene | 96 | | 3.423 | | | | | ND | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | | 3.508 | | | | | ND | |
| 24 Acetone | 43 | | 3.526 | | | | | ND | |
| 25 Iodomethane | 142 | | 3.617 | | | | | ND | |
| 26 Carbon disulfide | 76 | | 3.708 | | | | | ND | |
| 27 Isopropyl alcohol | 45 | | 3.812 | | | | | ND | |
| 29 Acetonitrile | 41 | | 3.964 | | | | | ND | |
| 28 3-Chloro-1-propene | 76 | | 3.994 | | | | | ND | |
| 30 Methyl acetate | 43 | | 4.031 | | | | | ND | |
| 31 Methylene Chloride | 84 | | 4.232 | | | | | ND | |
| 32 2-Methyl-2-propanol | 59 | | 4.493 | | | | | ND | |
| 33 Acrylonitrile | 53 | | 4.609 | | | | | 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 | | 4.633 | | | | | ND | |
| 35 Methyl tert-butyl ether | 73 | | 4.657 | | | | | ND | |
| 36 Hexane | 57 | | 5.053 | | | | | ND | |
| 37 1,1-Dichloroethane | 63 | | 5.272 | | | | | ND | |
| 38 Vinyl acetate | 43 | | 5.315 | | | | | ND | |
| 41 Isopropyl ether | 45 | | 5.363 | | | | | ND | |
| 39 2-Chloro-1,3-butadiene | 53 | | 5.363 | | | | | ND | |
| 40 Isopropyl ether TIC | 45 | | 5.410 | | | | | ND | |
| 42 Tert-butyl ethyl ether | 59 | | 5.832 | | | | | ND | |
| 43 Tert-butyl ethyl ether (TI | 59 | | 5.961 | | | | | ND | |
| 45 cis-1,2-Dichloroethene | 96 | | 6.008 | | | | | ND | |
| 44 2,2-Dichloropropane | 97 | | 6.008 | | | | | ND | |
| 46 2-Butanone (MEK) | 43 | | 6.026 | | | | | ND | |
| 47 Propionitrile | 54 | | 6.093 | | | | | ND | |
| 48 Ethyl acetate | 43 | | 6.099 | | | | | ND | |
| 50 Methacrylonitrile | 41 | | 6.276 | | | | | ND | |
| 49 Chlorobromomethane | 128 | | 6.294 | | | | | ND | |
| 51 Tetrahydrofuran | 42 | | 6.306 | | | | | ND | |
| 52 Chloroform | 83 | | 6.440 | | | | | ND | |
| 53 1,1,1-Trichloroethane | 97 | | 6.592 | | | | | ND | |
| 54 Cyclohexane | 56 | | 6.659 | | | | | ND | |
| 56 Carbon tetrachloride | 117 | | 6.762 | | | | | ND | |
| 55 1,1-Dichloropropene | 75 | | 6.781 | | | | | ND | |
| 57 Isobutyl alcohol | 41 | | 6.981 | | | | | ND | |
| 58 Benzene | 78 | | 6.994 | | | | | ND | |
| 59 1,2-Dichloroethane | 62 | | 7.067 | | | | | ND | |
| 151 Isooctane | 57 | | 7.146 | | | | | ND | |
| 61 Tert-amyl methyl ether | 73 | | 7.170 | | | | | ND | |
| 60 Tert-amyl methyl ether (TI | 73 | | 7.262 | | | | | ND | |
| 62 n-Heptane | 43 | | 7.352 | | | | | ND | |
| 63 n-Butanol | 56 | | 7.687 | | | | | ND | |
| 64 Trichloroethene | 130 | | 7.724 | | | | | ND | |
| 65 Ethyl acrylate | 55 | | 7.845 | | | | | ND | |
| 66 Methylcyclohexane | 83 | | 7.961 | | | | | ND | |
| 67 1,2-Dichloropropane | 63 | | 7.997 | | | | | ND | |
| 69 Methyl methacrylate | 69 | | 8.076 | | | | | ND | |
| 68 Dibromomethane | 93 | | 8.082 | | | | | ND | |
| 70 1,4-Dioxane | 88 | | 8.082 | | | | | ND | |
| 71 Dichlorobromomethane | 83 | | 8.277 | | | | | ND | |
| 73 2-Chloroethyl vinyl ether | 63 | | 8.575 | | | | | ND | |
| 74 cis-1,3-Dichloropropene | 75 | | 8.721 | | | | | ND | |
| 75 4-Methyl-2-pentanone (MIBK | 43 | | 8.873 | | | | | ND | |
| 76 Toluene | 91 | | 9.044 | | | | | ND | |
| 77 trans-1,3-Dichloropropene | 75 | | 9.293 | | | | | ND | |
| 78 Ethyl methacrylate | 69 | | 9.354 | | | | | ND | |
| 79 1,1,2-Trichloroethane | 97 | | 9.488 | | | | | ND | |
| 80 Tetrachloroethene | 164 | | 9.561 | | | | | ND | |
| 81 1,3-Dichloropropane | 76 | | 9.646 | | | | | ND | |
| 82 2-Hexanone | 43 | | 9.707 | | | | | ND | |
| 83 n-Butyl acetate | 43 | | 9.822 | | | | | ND | |
| 84 Chlorodibromomethane | 129 | | 9.853 | | | | | ND | |
| 85 Ethylene Dibromide | 107 | | 9.974 | | | | | ND | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|---|----------|------------|--------------|-------|
| 86 3-Chlorobenzotrifluoride | 180 | | 10.431 | | | | | ND | |
| 87 Chlorobenzene | 112 | | 10.455 | | | | | ND | |
| 88 4-Chlorobenzotrifluoride | 180 | | 10.522 | | | | | ND | |
| 89 1,1,1,2-Tetrachloroethane | 131 | | 10.552 | | | | | ND | |
| 90 Ethylbenzene | 106 | | 10.558 | | | | | ND | |
| 91 m-Xylene & p-Xylene | 106 | | 10.692 | | | | | ND | |
| 92 o-Xylene | 106 | | 11.069 | | | | | ND | |
| 93 Styrene | 104 | | 11.094 | | | | | ND | |
| 94 Bromoform | 173 | | 11.270 | | | | | ND | |
| 95 Cyclohexanol | 57 | | 11.288 | | | | | ND | |
| 96 2-Chlorobenzotrifluoride | 180 | | 11.343 | | | | | ND | |
| 97 Isopropylbenzene | 105 | | 11.441 | | | | | ND | |
| 98 Cyclohexanone | 55 | | 11.526 | | | | | ND | |
| 100 Bromobenzene | 156 | | 11.751 | | | | | ND | |
| 99 1,1,2,2-Tetrachloroethane | 83 | | 11.751 | | | | | ND | |
| 102 trans-1,4-Dichloro-2-buten | 53 | | 11.787 | | | | | ND | |
| 101 1,2,3-Trichloropropane | 110 | | 11.799 | | | | | ND | |
| 103 N-Propylbenzene | 120 | | 11.854 | | | | | ND | |
| 104 2-Chlorotoluene | 126 | | 11.939 | | | | | ND | |
| 105 3-Chlorotoluene | 126 | | 12.006 | | | | | ND | |
| 106 1,3,5-Trimethylbenzene | 105 | | 12.037 | | | | | ND | |
| 107 4-Chlorotoluene | 126 | | 12.061 | | | | | ND | |
| 108 tert-Butylbenzene | 119 | | 12.353 | | | | | ND | |
| 110 1,2,4-Trimethylbenzene | 105 | | 12.408 | | | | | ND | |
| 111 1,2-dichloro-4-(trifluorom | 214 | | 12.450 | | | | | ND | |
| 112 sec-Butylbenzene | 105 | | 12.572 | | | | | ND | |
| 113 1,3-Dichlorobenzene | 146 | | 12.694 | | | | | ND | |
| 114 4-Isopropyltoluene | 119 | | 12.730 | | | | | ND | |
| 115 1,4-Dichlorobenzene | 146 | | 12.797 | | | | | ND | |
| 116 2,4-Dichloro-1-(triflourom | 214 | | 12.822 | | | | | ND | |
| 117 1,2,3-Trimethylbenzene | 105 | | 12.822 | | | | | ND | |
| 118 2,5-Dichlorobenzotrifluori | 214 | | 12.864 | | | | | ND | |
| 119 Benzyl chloride | 91 | | 12.907 | | | | | ND | |
| 120 n-Butylbenzene | 91 | | 13.138 | | | | | ND | |
| 121 1,2-Dichlorobenzene | 146 | | 13.150 | | | | | ND | |
| 122 1,2-Dibromo-3-Chloropropan | 75 | | 13.947 | | | | | ND | |
| 123 2,4- & 2,5- & 2,6- Dichlor | 125 | | 14.081 | | | | | ND | |
| 124 1,3,5-Trichlorobenzene | 180 | | 14.130 | | | | | ND | |
| 125 2,3- & 3,4- Dichlorotoluen | 125 | | 14.501 | | | | | ND | |
| 126 1,2,4-Trichlorobenzene | 180 | | 14.768 | | | | | ND | |
| 127 Hexachlorobutadiene | 225 | | 14.908 | | | | | ND | |
| 128 Naphthalene | 128 | | 15.030 | | | | | ND | |
| 129 1,2,3-Trichlorobenzene | 180 | | 15.255 | | | | | ND | |
| 131 2,4,5-Trichlorotoluene | 159 | | 16.027 | | | | | ND | |
| 130 2,3,6-Trichlorotoluene | 159 | | 16.125 | | | | | ND | |
| 152 Formaldehyde TIC | 1 | | 0.000 | | | | | ND | |
| 149 3,4-Dichlorotoluene | 1 | | 0.000 | | | | | ND | |
| S 133 Xylenes, Total | 106 | | 1.000 | | | | | ND | |
| S 134 1,2-Dichloroethene, Total | 96 | | 1.000 | | | | | ND | |
| S 154 Total BTEX | 106 | | 1.000 | | | | | ND | |
| S 135 1,3-Dichloropropene, Total | 1 | | 0.000 | | | | | ND | |
| T 136 Mesityl oxide TIC | 83 | | 0.000 | | | | | ND | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|---------------|---|----------|------------|--------------|-------|
| T 138 Methyl n-amyl ketone TIC | 43 | | 0.000 | | | | | ND | |
| T 137 Tetrahydrofuran TIC | 42 | | 6.253 | | | | | ND | |
| T 153 1,2 Epoxybutane TIC | 42 | | 6.253 | | | | | ND | |

Reagents:

VOA8260INT_00074

Amount Added: 2.00

Units: uL

Run Reagent

VOA8260SURR_00073

Amount Added: 2.00

Units: uL

Run Reagent

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D06.D

Injection Date: 04-Oct-2017 02:21:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: MB

Worklist Smp#: 6

Client ID:

Purge Vol: 5.000 mL

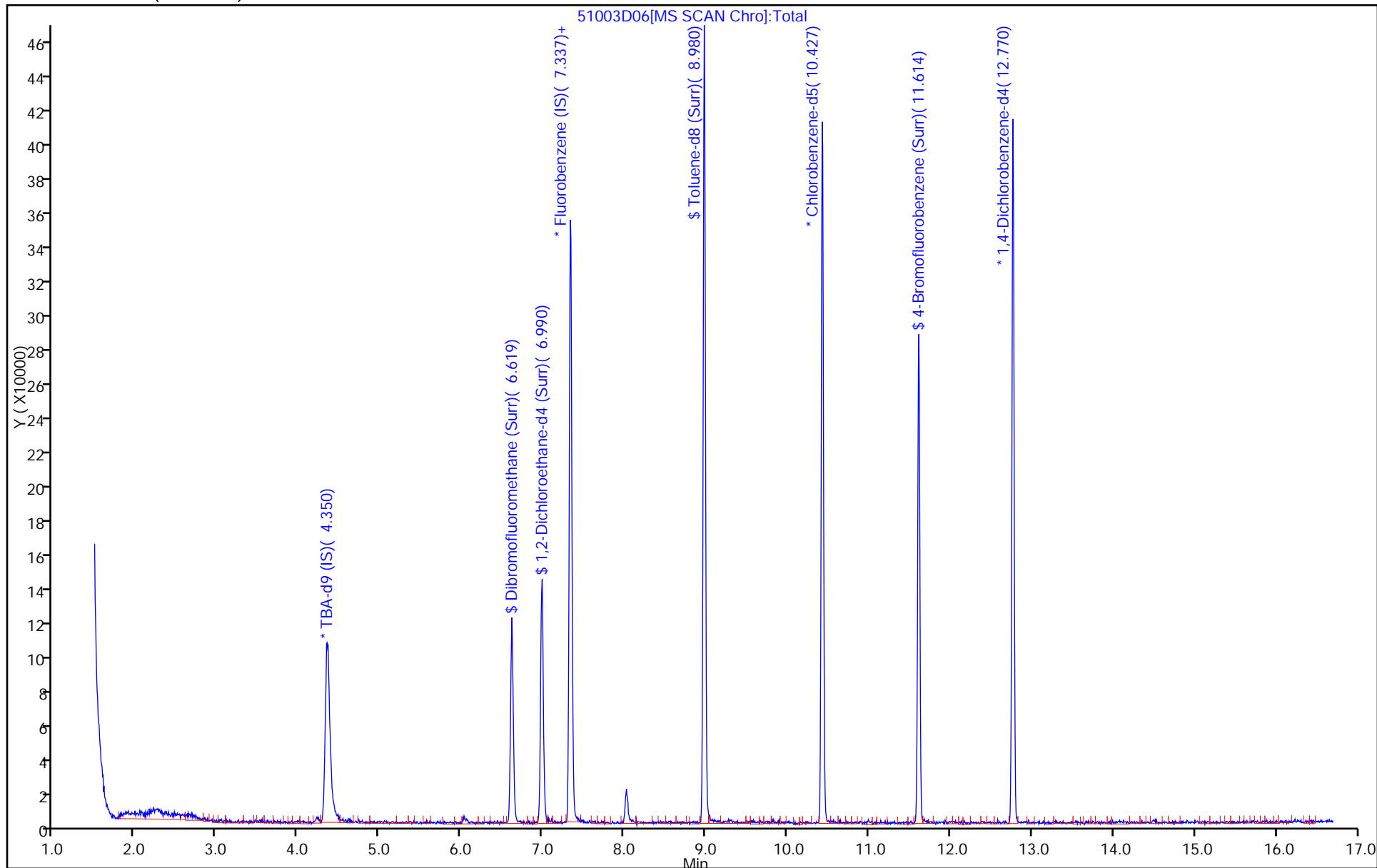
Dil. Factor: 1.0000

ALS Bottle#: 6

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Recovery Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D06.D
 Lims ID: MB
 Client ID:
 Sample Type: MB
 Inject. Date: 04-Oct-2017 02:21:30 ALS Bottle#: 6 Worklist Smp#: 6
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018710-006
 Misc. Info.: MB
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 04-Oct-2017 21:10:21 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK005

First Level Reviewer: bungardf

Date: 04-Oct-2017 02:42:56

| Compound | Amount Added | Amount Recovered | % Rec. |
|-----------------------------------|--------------|------------------|--------|
| \$ 5 Dibromofluoromethane (Surr) | 50.0 | 46.8 | 93.69 |
| \$ 6 1,2-Dichloroethane-d4 (Surr) | 50.0 | 53.6 | 107.13 |
| \$ 7 Toluene-d8 (Surr) | 50.0 | 49.0 | 97.96 |
| \$ 8 4-Bromofluorobenzene (Surr) | 50.0 | 46.3 | 92.57 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Client Sample ID: _____ Lab Sample ID: MB 180-224919/7
 Matrix: Water Lab File ID: 51004D07.D
 Analysis Method: 8260C Date Collected: _____
 Sample wt/vol: 5 (mL) Date Analyzed: 10/05/2017 02:41
 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.: 224919 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|------------|-----------------------------|--------|---|-----|------|
| 74-87-3 | Chloromethane | 1.0 | U | 1.0 | 0.38 |
| 75-01-4 | Vinyl chloride | 1.0 | U | 1.0 | 0.17 |
| 74-83-9 | Bromomethane | 1.0 | U | 1.0 | 0.59 |
| 75-00-3 | Chloroethane | 1.0 | U | 1.0 | 0.58 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 | U | 1.0 | 0.32 |
| 67-64-1 | Acetone | 5.0 | U | 5.0 | 3.1 |
| 75-15-0 | Carbon disulfide | 1.0 | U | 1.0 | 0.53 |
| 75-09-2 | Methylene Chloride | 1.0 | U | 1.0 | 0.94 |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 1.0 | U | 1.0 | 0.20 |
| 75-34-3 | 1,1-Dichloroethane | 1.0 | U | 1.0 | 0.34 |
| 156-59-2 | cis-1,2-Dichloroethene | 1.0 | U | 1.0 | 0.30 |
| 74-97-5 | Bromochloromethane | 1.0 | U | 1.0 | 0.36 |
| 78-93-3 | 2-Butanone (MEK) | 5.0 | U | 5.0 | 2.6 |
| 67-66-3 | Chloroform | 1.0 | U | 1.0 | 0.27 |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 | U | 1.0 | 0.27 |
| 56-23-5 | Carbon tetrachloride | 1.0 | U | 1.0 | 0.56 |
| 71-43-2 | Benzene | 1.0 | U | 1.0 | 0.18 |
| 107-06-2 | 1,2-Dichloroethane | 1.0 | U | 1.0 | 0.24 |
| 79-01-6 | Trichloroethene | 1.0 | U | 1.0 | 0.20 |
| 78-87-5 | 1,2-Dichloropropane | 1.0 | U | 1.0 | 0.35 |
| 75-27-4 | Bromodichloromethane | 1.0 | U | 1.0 | 0.57 |
| 10061-01-5 | cis-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.32 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 5.0 | U | 5.0 | 2.2 |
| 108-88-3 | Toluene | 1.0 | U | 1.0 | 0.16 |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 | U | 1.0 | 0.22 |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 | U | 1.0 | 0.31 |
| 127-18-4 | Tetrachloroethene | 1.0 | U | 1.0 | 0.24 |
| 591-78-6 | 2-Hexanone | 5.0 | U | 5.0 | 2.0 |
| 124-48-1 | Dibromochloromethane | 1.0 | U | 1.0 | 0.44 |
| 106-93-4 | 1,2-Dibromoethane (EDB) | 1.0 | U | 1.0 | 0.51 |
| 108-90-7 | Chlorobenzene | 1.0 | U | 1.0 | 0.15 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.49 |
| 100-41-4 | Ethylbenzene | 1.0 | U | 1.0 | 0.25 |
| 1330-20-7 | Xylenes, Total | 2.0 | U | 2.0 | 0.27 |
| 100-42-5 | Styrene | 1.0 | U | 1.0 | 0.22 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Client Sample ID: _____ Lab Sample ID: MB 180-224919/7
 Matrix: Water Lab File ID: 51004D07.D
 Analysis Method: 8260C Date Collected: _____
 Sample wt/vol: 5 (mL) Date Analyzed: 10/05/2017 02:41
 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.: 224919 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|----------|---------------------------|--------|---|-----|------|
| 75-25-2 | Bromoform | 1.0 | U | 1.0 | 0.76 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 | U | 1.0 | 0.37 |
| 107-13-1 | Acrylonitrile | 20 | U | 20 | 3.3 |
| 123-91-1 | 1,4-Dioxane | 200 | U | 200 | 16 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|------------|------------------------------|------|---|--------|
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 111 | | 65-121 |
| 2037-26-5 | Toluene-d8 (Surr) | 101 | | 73-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 95 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 96 | | 73-120 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D07.D
 Lims ID: MB
 Client ID:
 Sample Type: MB
 Inject. Date: 05-Oct-2017 02:41:30 ALS Bottle#: 7 Worklist Smp#: 7
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018725-007
 Misc. Info.: MB
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 08-Oct-2017 20:41:27 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK026

First Level Reviewer: bungardf

Date: 05-Oct-2017 03:04:15

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.359 | 4.352 | 0.007 | 0 | 189568 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.340 | 7.339 | 0.001 | 99 | 369267 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.430 | 10.429 | 0.001 | 85 | 81221 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.766 | 12.771 | -0.005 | 96 | 115186 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.616 | 6.613 | 0.003 | 93 | 85389 | 50.0 | 48.1 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.987 | 6.990 | -0.003 | 0 | 120443 | 50.0 | 55.6 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.983 | 8.979 | 0.004 | 93 | 325966 | 50.0 | 50.4 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.611 | 11.613 | -0.002 | 85 | 111301 | 50.0 | 47.7 | |
| 11 Dichlorodifluoromethane | 85 | | 1.679 | | | | | ND | |
| 12 Chloromethane | 50 | | 1.825 | | | | | ND | |
| 13 Vinyl chloride | 62 | | 1.959 | | | | | ND | |
| 14 Butadiene | 39 | | 2.008 | | | | | ND | |
| 15 Bromomethane | 94 | | 2.300 | | | | | ND | |
| 16 Chloroethane | 64 | | 2.470 | | | | | ND | |
| 17 Dichlorofluoromethane | 67 | | 2.744 | | | | | ND | |
| 18 Trichlorofluoromethane | 101 | | 2.768 | | | | | ND | |
| 19 Ethanol | 45 | | 2.821 | | | | | ND | |
| 20 Ethyl ether | 59 | | 3.121 | | | | | ND | |
| 21 Acrolein | 56 | | 3.316 | | | | | ND | |
| 22 1,1-Dichloroethene | 96 | | 3.413 | | | | | ND | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | | 3.504 | | | | | ND | |
| 24 Acetone | 43 | | 3.529 | | | | | ND | |
| 25 Iodomethane | 142 | | 3.608 | | | | | ND | |
| 26 Carbon disulfide | 76 | | 3.699 | | | | | ND | |
| 27 Isopropyl alcohol | 45 | | 3.811 | | | | | ND | |
| 29 Acetonitrile | 41 | | 3.969 | | | | | ND | |
| 28 3-Chloro-1-propene | 76 | | 3.997 | | | | | ND | |
| 30 Methyl acetate | 43 | | 4.033 | | | | | ND | |
| 31 Methylene Chloride | 84 | | 4.222 | | | | | ND | |
| 32 2-Methyl-2-propanol | 59 | | 4.508 | | | | | ND | |
| 33 Acrylonitrile | 53 | | 4.605 | | | | | 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 | | 4.642 | | | | | ND | |
| 35 Methyl tert-butyl ether | 73 | | 4.660 | | | | | ND | |
| 36 Hexane | 57 | | 5.049 | | | | | ND | |
| 37 1,1-Dichloroethane | 63 | | 5.268 | | | | | ND | |
| 38 Vinyl acetate | 43 | | 5.323 | | | | | ND | |
| 39 2-Chloro-1,3-butadiene | 53 | | 5.356 | | | | | ND | |
| 41 Isopropyl ether | 45 | | 5.368 | | | | | ND | |
| 40 Isopropyl ether TIC | 45 | | 5.410 | | | | | ND | |
| 42 Tert-butyl ethyl ether | 59 | | 5.836 | | | | | ND | |
| 43 Tert-butyl ethyl ether (TI | 59 | | 5.961 | | | | | ND | |
| 44 2,2-Dichloropropane | 97 | | 6.004 | | | | | ND | |
| 45 cis-1,2-Dichloroethene | 96 | | 6.011 | | | | | ND | |
| 46 2-Butanone (MEK) | 43 | | 6.023 | | | | | ND | |
| 48 Ethyl acetate | 43 | | 6.092 | | | | | ND | |
| 47 Propionitrile | 54 | | 6.098 | | | | | ND | |
| 50 Methacrylonitrile | 41 | | 6.274 | | | | | ND | |
| 49 Chlorobromomethane | 128 | | 6.290 | | | | | ND | |
| 51 Tetrahydrofuran | 42 | | 6.309 | | | | | ND | |
| 52 Chloroform | 83 | | 6.436 | | | | | ND | |
| 53 1,1,1-Trichloroethane | 97 | | 6.595 | | | | | ND | |
| 54 Cyclohexane | 56 | | 6.662 | | | | | ND | |
| 56 Carbon tetrachloride | 117 | | 6.759 | | | | | ND | |
| 55 1,1-Dichloropropene | 75 | | 6.783 | | | | | ND | |
| 57 Isobutyl alcohol | 41 | | 6.990 | | | | | ND | |
| 58 Benzene | 78 | | 6.996 | | | | | ND | |
| 59 1,2-Dichloroethane | 62 | | 7.069 | | | | | ND | |
| 151 Isooctane | 57 | | 7.144 | | | | | ND | |
| 61 Tert-amyl methyl ether | 73 | | 7.175 | | | | | ND | |
| 60 Tert-amyl methyl ether (TI | 73 | | 7.262 | | | | | ND | |
| 62 n-Heptane | 43 | | 7.349 | | | | | ND | |
| 63 n-Butanol | 56 | | 7.686 | | | | | ND | |
| 64 Trichloroethene | 130 | | 7.720 | | | | | ND | |
| 65 Ethyl acrylate | 55 | | 7.844 | | | | | ND | |
| 66 Methylcyclohexane | 83 | | 7.957 | | | | | ND | |
| 67 1,2-Dichloropropane | 63 | | 7.994 | | | | | ND | |
| 68 Dibromomethane | 93 | | 8.079 | | | | | ND | |
| 69 Methyl methacrylate | 69 | | 8.081 | | | | | ND | |
| 70 1,4-Dioxane | 88 | | 8.085 | | | | | ND | |
| 71 Dichlorobromomethane | 83 | | 8.280 | | | | | ND | |
| 73 2-Chloroethyl vinyl ether | 63 | | 8.578 | | | | | ND | |
| 74 cis-1,3-Dichloropropene | 75 | | 8.718 | | | | | ND | |
| 75 4-Methyl-2-pentanone (MIBK | 43 | | 8.876 | | | | | ND | |
| 76 Toluene | 91 | | 9.046 | | | | | ND | |
| 77 trans-1,3-Dichloropropene | 75 | | 9.296 | | | | | ND | |
| 78 Ethyl methacrylate | 69 | | 9.356 | | | | | ND | |
| 79 1,1,2-Trichloroethane | 97 | | 9.490 | | | | | ND | |
| 80 Tetrachloroethene | 164 | | 9.563 | | | | | ND | |
| 81 1,3-Dichloropropane | 76 | | 9.648 | | | | | ND | |
| 82 2-Hexanone | 43 | | 9.703 | | | | | ND | |
| 83 n-Butyl acetate | 43 | | 9.827 | | | | | ND | |
| 84 Chlorodibromomethane | 129 | | 9.855 | | | | | ND | |
| 85 Ethylene Dibromide | 107 | | 9.971 | | | | | ND | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|---|----------|------------|--------------|-------|
| 86 3-Chlorobenzotrifluoride | 180 | | 10.433 | | | | | ND | |
| 87 Chlorobenzene | 112 | | 10.458 | | | | | ND | |
| 88 4-Chlorobenzotrifluoride | 180 | | 10.518 | | | | | ND | |
| 89 1,1,1,2-Tetrachloroethane | 131 | | 10.549 | | | | | ND | |
| 90 Ethylbenzene | 106 | | 10.555 | | | | | ND | |
| 91 m-Xylene & p-Xylene | 106 | | 10.689 | | | | | ND | |
| 92 o-Xylene | 106 | | 11.072 | | | | | ND | |
| 93 Styrene | 104 | | 11.090 | | | | | ND | |
| 94 Bromoform | 173 | | 11.273 | | | | | ND | |
| 95 Cyclohexanol | 57 | | 11.288 | | | | | ND | |
| 96 2-Chlorobenzotrifluoride | 180 | | 11.340 | | | | | ND | |
| 97 Isopropylbenzene | 105 | | 11.437 | | | | | ND | |
| 98 Cyclohexanone | 55 | | 11.530 | | | | | ND | |
| 99 1,1,2,2-Tetrachloroethane | 83 | | 11.747 | | | | | ND | |
| 100 Bromobenzene | 156 | | 11.753 | | | | | ND | |
| 102 trans-1,4-Dichloro-2-buten | 53 | | 11.784 | | | | | ND | |
| 101 1,2,3-Trichloropropane | 110 | | 11.808 | | | | | ND | |
| 103 N-Propylbenzene | 120 | | 11.851 | | | | | ND | |
| 104 2-Chlorotoluene | 126 | | 11.942 | | | | | ND | |
| 105 3-Chlorotoluene | 126 | | 12.003 | | | | | ND | |
| 106 1,3,5-Trimethylbenzene | 105 | | 12.033 | | | | | ND | |
| 107 4-Chlorotoluene | 126 | | 12.064 | | | | | ND | |
| 108 tert-Butylbenzene | 119 | | 12.350 | | | | | ND | |
| 110 1,2,4-Trimethylbenzene | 105 | | 12.410 | | | | | ND | |
| 111 1,2-dichloro-4-(trifluorom | 214 | | 12.453 | | | | | ND | |
| 112 sec-Butylbenzene | 105 | | 12.575 | | | | | ND | |
| 113 1,3-Dichlorobenzene | 146 | | 12.690 | | | | | ND | |
| 114 4-Isopropyltoluene | 119 | | 12.727 | | | | | ND | |
| 115 1,4-Dichlorobenzene | 146 | | 12.794 | | | | | ND | |
| 116 2,4-Dichloro-1-(triflourom | 214 | | 12.818 | | | | | ND | |
| 117 1,2,3-Trimethylbenzene | 105 | | 12.820 | | | | | ND | |
| 118 2,5-Dichlorobenzotrifluori | 214 | | 12.867 | | | | | ND | |
| 119 Benzyl chloride | 91 | | 12.911 | | | | | ND | |
| 120 n-Butylbenzene | 91 | | 13.140 | | | | | ND | |
| 121 1,2-Dichlorobenzene | 146 | | 13.153 | | | | | ND | |
| 122 1,2-Dibromo-3-Chloropropan | 75 | | 13.937 | | | | | ND | |
| 123 2,4- & 2,5- & 2,6- Dichlor | 125 | | 14.083 | | | | | ND | |
| 124 1,3,5-Trichlorobenzene | 180 | 14.129 | 14.128 | 0.001 | 1 | 280 | | 0.1174 | |
| 125 2,3- & 3,4- Dichlorotoluen | 125 | | 14.503 | | | | | ND | |
| 126 1,2,4-Trichlorobenzene | 180 | | 14.765 | | | | | ND | |
| 127 Hexachlorobutadiene | 225 | | 14.905 | | | | | ND | |
| 128 Naphthalene | 128 | | 15.026 | | | | | ND | |
| 129 1,2,3-Trichlorobenzene | 180 | | 15.257 | | | | | ND | |
| 131 2,4,5-Trichlorotoluene | 159 | | 16.024 | | | | | ND | |
| 130 2,3,6-Trichlorotoluene | 159 | | 16.121 | | | | | ND | |
| 152 Formaldehyde TIC | 1 | | 0.000 | | | | | ND | |
| 149 3,4-Dichlorotoluene | 1 | | 0.000 | | | | | ND | |
| S 133 Xylenes, Total | 106 | | 1.000 | | | | | ND | |
| S 134 1,2-Dichloroethene, Total | 96 | | 1.000 | | | | | ND | |
| S 154 Total BTEX | 106 | | 1.000 | | | | | ND | |
| S 135 1,3-Dichloropropene, Total | 1 | | 0.000 | | | | | ND | |
| T 136 Mesityl oxide TIC | 83 | | 0.000 | | | | | ND | |

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

| | | | | | | | | | |
|--------------------------------|----|--|-------|--|--|--|--|--|----|
| T 138 Methyl n-amyl ketone TIC | 43 | | 0.000 | | | | | | ND |
| T 137 Tetrahydrofuran TIC | 42 | | 6.253 | | | | | | ND |
| T 153 1,2 Epoxybutane TIC | 42 | | 6.253 | | | | | | ND |

Reagents:

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

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D07.D

Injection Date: 05-Oct-2017 02:41:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: MB

Worklist Smp#: 7

Client ID:

Purge Vol: 5.000 mL

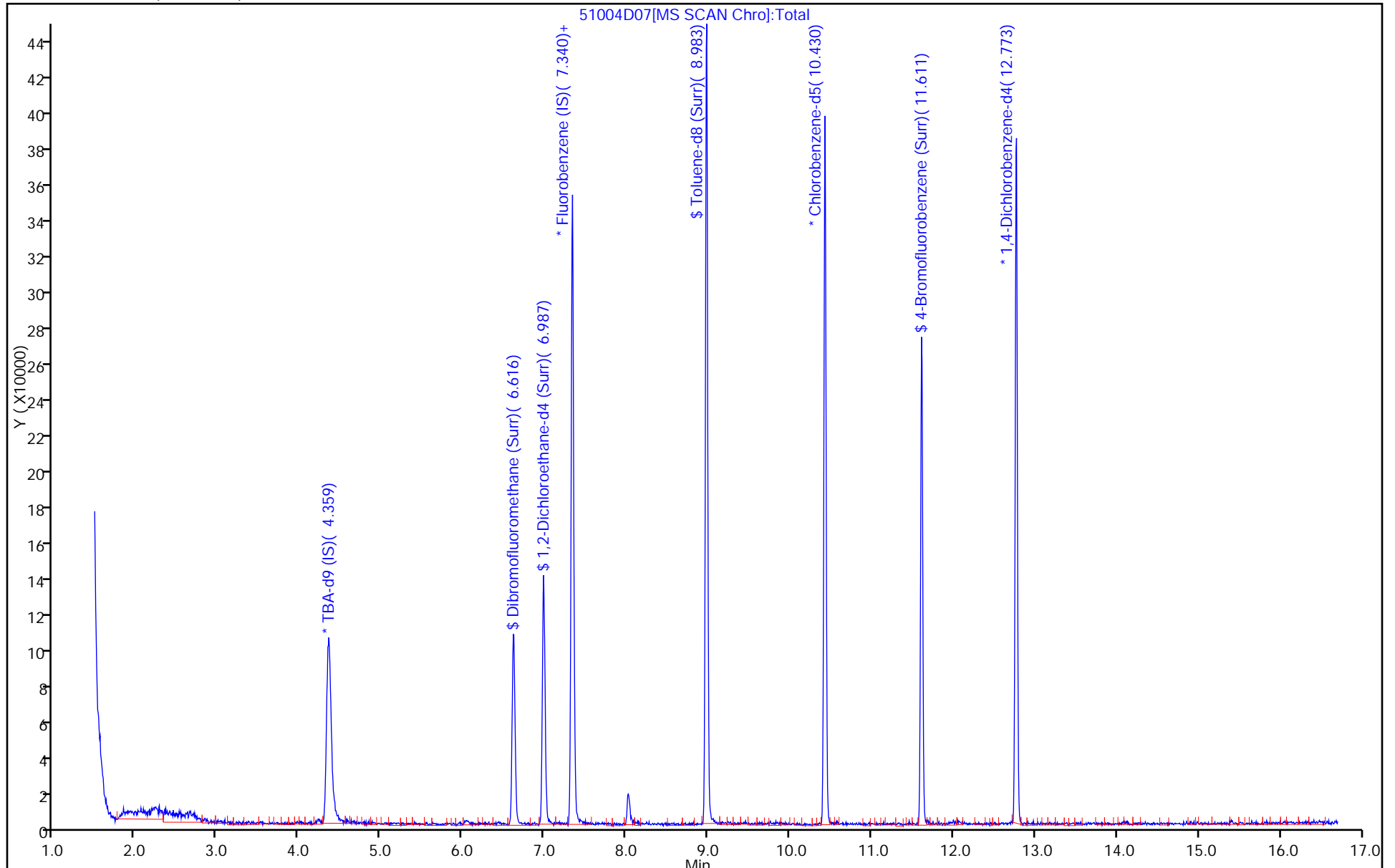
Dil. Factor: 1.0000

ALS Bottle#: 7

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Recovery Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D07.D
 Lims ID: MB
 Client ID:
 Sample Type: MB
 Inject. Date: 05-Oct-2017 02:41:30 ALS Bottle#: 7 Worklist Smp#: 7
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018725-007
 Misc. Info.: MB
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 08-Oct-2017 20:41:27 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK026

First Level Reviewer: bungardf

Date: 05-Oct-2017 03:04:15

| Compound | Amount Added | Amount Recovered | % Rec. |
|-----------------------------------|--------------|------------------|--------|
| \$ 5 Dibromofluoromethane (Surr) | 50.0 | 48.1 | 96.12 |
| \$ 6 1,2-Dichloroethane-d4 (Surr) | 50.0 | 55.6 | 111.16 |
| \$ 7 Toluene-d8 (Surr) | 50.0 | 50.4 | 100.85 |
| \$ 8 4-Bromofluorobenzene (Surr) | 50.0 | 47.7 | 95.35 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Client Sample ID: _____ Lab Sample ID: LCS 180-224792/4
 Matrix: Water Lab File ID: 51003D04.D
 Analysis Method: 8260C Date Collected: _____
 Sample wt/vol: 5 (mL) Date Analyzed: 10/04/2017 01:24
 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.: 224792 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|------------|-----------------------------|--------|---|-----|------|
| 74-87-3 | Chloromethane | 11.6 | | 1.0 | 0.38 |
| 75-01-4 | Vinyl chloride | 11.8 | | 1.0 | 0.17 |
| 74-83-9 | Bromomethane | 10.1 | | 1.0 | 0.59 |
| 75-00-3 | Chloroethane | 13.5 | | 1.0 | 0.58 |
| 75-35-4 | 1,1-Dichloroethene | 10.3 | | 1.0 | 0.32 |
| 67-64-1 | Acetone | 22.4 | | 5.0 | 3.1 |
| 75-15-0 | Carbon disulfide | 9.37 | | 1.0 | 0.53 |
| 75-09-2 | Methylene Chloride | 9.77 | | 1.0 | 0.94 |
| 156-60-5 | trans-1,2-Dichloroethene | 10.0 | | 1.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 9.62 | | 1.0 | 0.20 |
| 75-34-3 | 1,1-Dichloroethane | 9.85 | | 1.0 | 0.34 |
| 156-59-2 | cis-1,2-Dichloroethene | 9.81 | | 1.0 | 0.30 |
| 74-97-5 | Bromochloromethane | 9.53 | | 1.0 | 0.36 |
| 78-93-3 | 2-Butanone (MEK) | 21.0 | | 5.0 | 2.6 |
| 67-66-3 | Chloroform | 9.83 | | 1.0 | 0.27 |
| 71-55-6 | 1,1,1-Trichloroethane | 10.4 | | 1.0 | 0.27 |
| 56-23-5 | Carbon tetrachloride | 9.98 | | 1.0 | 0.56 |
| 71-43-2 | Benzene | 9.68 | | 1.0 | 0.18 |
| 107-06-2 | 1,2-Dichloroethane | 10.6 | | 1.0 | 0.24 |
| 79-01-6 | Trichloroethene | 9.28 | | 1.0 | 0.20 |
| 78-87-5 | 1,2-Dichloropropane | 9.20 | | 1.0 | 0.35 |
| 75-27-4 | Bromodichloromethane | 9.11 | | 1.0 | 0.57 |
| 10061-01-5 | cis-1,3-Dichloropropene | 8.97 | | 1.0 | 0.32 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 21.2 | | 5.0 | 2.2 |
| 108-88-3 | Toluene | 10.9 | | 1.0 | 0.16 |
| 10061-02-6 | trans-1,3-Dichloropropene | 9.70 | | 1.0 | 0.22 |
| 79-00-5 | 1,1,2-Trichloroethane | 10.8 | | 1.0 | 0.31 |
| 127-18-4 | Tetrachloroethene | 10.5 | | 1.0 | 0.24 |
| 591-78-6 | 2-Hexanone | 21.2 | | 5.0 | 2.0 |
| 124-48-1 | Dibromochloromethane | 9.60 | | 1.0 | 0.44 |
| 106-93-4 | 1,2-Dibromoethane (EDB) | 9.86 | | 1.0 | 0.51 |
| 108-90-7 | Chlorobenzene | 10.3 | | 1.0 | 0.15 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 10.6 | | 1.0 | 0.49 |
| 100-41-4 | Ethylbenzene | 9.83 | | 1.0 | 0.25 |
| 1330-20-7 | Xylenes, Total | 20.2 | | 2.0 | 0.27 |
| 100-42-5 | Styrene | 10.0 | | 1.0 | 0.22 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Client Sample ID: _____ Lab Sample ID: LCS 180-224792/4
 Matrix: Water Lab File ID: 51003D04.D
 Analysis Method: 8260C Date Collected: _____
 Sample wt/vol: 5 (mL) Date Analyzed: 10/04/2017 01:24
 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.: 224792 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|----------|---------------------------|--------|---|-----|------|
| 75-25-2 | Bromoform | 8.98 | | 1.0 | 0.76 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 10.4 | | 1.0 | 0.37 |
| 107-13-1 | Acrylonitrile | 108 | | 20 | 3.3 |
| 123-91-1 | 1,4-Dioxane | 271 | | 200 | 16 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|------------|------------------------------|------|---|--------|
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 106 | | 65-121 |
| 2037-26-5 | Toluene-d8 (Surr) | 113 | | 73-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 107 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 98 | | 73-120 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D04.D
 Lims ID: LCS
 Client ID:
 Sample Type: LCS
 Inject. Date: 04-Oct-2017 01:24:30 ALS Bottle#: 4 Worklist Smp#: 4
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018710-004
 Misc. Info.: LCS
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 04-Oct-2017 21:10:21 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK005

First Level Reviewer: bungardf

Date: 04-Oct-2017 02:09:21

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.355 | 4.347 | 0.008 | 0 | 183806 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.342 | 7.334 | 0.008 | 97 | 350927 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.426 | 10.431 | -0.005 | 85 | 75706 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.768 | 12.773 | -0.005 | 95 | 111504 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.618 | 6.616 | 0.002 | 93 | 82762 | 50.0 | 49.0 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.983 | 6.981 | 0.002 | 0 | 109269 | 50.0 | 53.1 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.978 | 8.977 | 0.001 | 93 | 341582 | 50.0 | 56.7 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.612 | 11.611 | 0.001 | 86 | 116849 | 50.0 | 53.7 | |
| 11 Dichlorodifluoromethane | 85 | 1.678 | 1.689 | -0.011 | 99 | 103253 | 50.0 | 50.6 | |
| 12 Chloromethane | 50 | 1.824 | 1.823 | 0.001 | 98 | 118538 | 50.0 | 57.8 | |
| 13 Vinyl chloride | 62 | 1.952 | 1.963 | -0.011 | 97 | 122919 | 50.0 | 59.1 | |
| 14 Butadiene | 39 | 1.994 | 1.993 | 0.001 | 98 | 121229 | 50.0 | 64.1 | |
| 15 Bromomethane | 94 | 2.305 | 2.291 | 0.014 | 87 | 49551 | 50.0 | 50.4 | |
| 16 Chloroethane | 64 | 2.451 | 2.461 | -0.010 | 98 | 76988 | 50.0 | 67.3 | |
| 17 Dichlorofluoromethane | 67 | 2.743 | 2.735 | 0.008 | 97 | 174905 | 50.0 | 60.4 | |
| 18 Trichlorofluoromethane | 101 | 2.743 | 2.790 | -0.047 | 94 | 147056 | 50.0 | 57.5 | |
| 20 Ethyl ether | 59 | 3.120 | 3.124 | -0.004 | 92 | 90317 | 50.0 | 54.3 | |
| 21 Acrolein | 56 | 3.308 | 3.301 | 0.007 | 99 | 72376 | 150.0 | 172.7 | |
| 22 1,1-Dichloroethene | 96 | 3.430 | 3.423 | 0.007 | 97 | 88643 | 50.0 | 51.6 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.503 | 3.508 | -0.005 | 90 | 102362 | 50.0 | 54.3 | |
| 24 Acetone | 43 | 3.521 | 3.526 | -0.005 | 99 | 102980 | 100.0 | 112.2 | |
| 25 Iodomethane | 142 | 3.619 | 3.617 | 0.002 | 97 | 136016 | 50.0 | 50.4 | |
| 26 Carbon disulfide | 76 | 3.710 | 3.708 | 0.002 | 99 | 176720 | 50.0 | 46.9 | |
| 28 3-Chloro-1-propene | 76 | 4.008 | 3.994 | 0.014 | 95 | 52288 | 50.0 | 47.1 | |
| 30 Methyl acetate | 43 | 4.032 | 4.031 | 0.001 | 96 | 190761 | 100.0 | 105.0 | |
| 31 Methylene Chloride | 84 | 4.221 | 4.232 | -0.011 | 93 | 104105 | 50.0 | 48.8 | |
| 32 2-Methyl-2-propanol | 59 | 4.483 | 4.493 | -0.010 | 94 | 108171 | 500.0 | 497.6 | |
| 33 Acrylonitrile | 53 | 4.610 | 4.609 | 0.001 | 99 | 478450 | 500.0 | 541.5 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.641 | 4.633 | 0.008 | 99 | 98310 | 50.0 | 50.2 | |
| 35 Methyl tert-butyl ether | 73 | 4.659 | 4.657 | 0.002 | 97 | 252538 | 50.0 | 48.1 | |
| 36 Hexane | 57 | 5.054 | 5.053 | 0.001 | 94 | 129294 | 50.0 | 51.5 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 37 1,1-Dichloroethane | 63 | 5.267 | 5.272 | -0.005 | 97 | 167599 | 50.0 | 49.2 | |
| 38 Vinyl acetate | 43 | 5.322 | 5.315 | 0.008 | 97 | 181173 | 50.0 | 52.3 | |
| 45 cis-1,2-Dichloroethene | 96 | 6.016 | 6.008 | 0.008 | 80 | 109826 | 50.0 | 49.1 | |
| 44 2,2-Dichloropropane | 97 | 5.997 | 6.008 | -0.011 | 62 | 23638 | 50.0 | 54.5 | |
| 46 2-Butanone (MEK) | 43 | 6.022 | 6.026 | -0.004 | 99 | 136976 | 100.0 | 104.9 | |
| 49 Chlorobromomethane | 128 | 6.295 | 6.294 | 0.001 | 97 | 47438 | 50.0 | 47.7 | |
| 51 Tetrahydrofuran | 42 | 6.308 | 6.306 | 0.002 | 93 | 71574 | 100.0 | 94.1 | |
| 52 Chloroform | 83 | 6.435 | 6.440 | -0.005 | 94 | 166997 | 50.0 | 49.1 | |
| 53 1,1,1-Trichloroethane | 97 | 6.594 | 6.592 | 0.002 | 98 | 133821 | 50.0 | 52.0 | |
| 54 Cyclohexane | 56 | 6.667 | 6.659 | 0.008 | 91 | 163595 | 50.0 | 51.5 | |
| 56 Carbon tetrachloride | 117 | 6.764 | 6.762 | 0.002 | 97 | 106851 | 50.0 | 49.9 | |
| 55 1,1-Dichloropropene | 75 | 6.776 | 6.781 | -0.005 | 97 | 135425 | 50.0 | 48.7 | |
| 57 Isobutyl alcohol | 41 | 6.977 | 6.981 | -0.004 | 88 | 101254 | 1250.0 | 1450.1 | |
| 58 Benzene | 78 | 6.995 | 6.994 | 0.001 | 97 | 413043 | 50.0 | 48.4 | |
| 59 1,2-Dichloroethane | 62 | 7.074 | 7.067 | 0.007 | 98 | 131706 | 50.0 | 53.0 | |
| 62 n-Heptane | 43 | 7.354 | 7.352 | 0.002 | 89 | 103197 | 50.0 | 51.4 | |
| 64 Trichloroethene | 130 | 7.725 | 7.724 | 0.001 | 98 | 99645 | 50.0 | 46.4 | |
| 66 Methylcyclohexane | 83 | 7.962 | 7.961 | 0.001 | 88 | 150041 | 50.0 | 46.2 | |
| 67 1,2-Dichloropropane | 63 | 7.993 | 7.997 | -0.004 | 93 | 91445 | 50.0 | 46.0 | |
| 68 Dibromomethane | 93 | 8.078 | 8.082 | -0.004 | 96 | 53748 | 50.0 | 46.2 | |
| 70 1,4-Dioxane | 88 | 8.078 | 8.082 | -0.004 | 51 | 27346 | 1000.0 | 1353.5 | |
| 71 Dichlorobromomethane | 83 | 8.273 | 8.277 | -0.004 | 99 | 104118 | 50.0 | 45.6 | |
| 73 2-Chloroethyl vinyl ether | 63 | 8.577 | 8.575 | 0.002 | 93 | 104178 | 100.0 | 72.9 | |
| 74 cis-1,3-Dichloropropene | 75 | 8.723 | 8.721 | 0.002 | 94 | 124521 | 50.0 | 44.9 | |
| 75 4-Methyl-2-pentanone (MIBK) | 43 | 8.875 | 8.873 | 0.002 | 96 | 205423 | 100.0 | 105.8 | |
| 76 Toluene | 91 | 9.051 | 9.044 | 0.007 | 99 | 411022 | 50.0 | 54.4 | |
| 77 trans-1,3-Dichloropropene | 75 | 9.295 | 9.293 | 0.002 | 95 | 99666 | 50.0 | 48.5 | |
| 78 Ethyl methacrylate | 69 | 9.355 | 9.354 | 0.001 | 89 | 111013 | 50.0 | 44.8 | |
| 79 1,1,2-Trichloroethane | 97 | 9.489 | 9.488 | 0.001 | 89 | 84987 | 50.0 | 54.0 | |
| 80 Tetrachloroethene | 164 | 9.562 | 9.561 | 0.001 | 96 | 75842 | 50.0 | 52.7 | |
| 81 1,3-Dichloropropane | 76 | 9.647 | 9.646 | 0.001 | 90 | 146002 | 50.0 | 50.2 | |
| 82 2-Hexanone | 43 | 9.708 | 9.707 | 0.001 | 99 | 157606 | 100.0 | 105.8 | |
| 84 Chlorodibromomethane | 129 | 9.860 | 9.853 | 0.007 | 91 | 63786 | 50.0 | 48.0 | |
| 85 Ethylene Dibromide | 107 | 9.970 | 9.974 | -0.004 | 99 | 79479 | 50.0 | 49.3 | |
| 86 3-Chlorobenzotrifluoride | 180 | 10.432 | 10.431 | 0.001 | 89 | 157842 | 50.0 | 60.7 | |
| 87 Chlorobenzene | 112 | 10.457 | 10.455 | 0.002 | 95 | 252385 | 50.0 | 51.4 | |
| 88 4-Chlorobenzotrifluoride | 180 | 10.517 | 10.522 | -0.005 | 95 | 150026 | 50.0 | 62.5 | |
| 89 1,1,1,2-Tetrachloroethane | 131 | 10.548 | 10.552 | -0.004 | 91 | 83133 | 50.0 | 53.2 | |
| 90 Ethylbenzene | 106 | 10.560 | 10.558 | 0.002 | 99 | 134813 | 50.0 | 49.1 | |
| 91 m-Xylene & p-Xylene | 106 | 10.688 | 10.692 | -0.004 | 0 | 172133 | 50.0 | 51.3 | |
| 92 o-Xylene | 106 | 11.071 | 11.069 | 0.002 | 96 | 158520 | 50.0 | 49.6 | |
| 93 Styrene | 104 | 11.089 | 11.094 | -0.005 | 95 | 271030 | 50.0 | 50.1 | |
| 94 Bromoform | 173 | 11.272 | 11.270 | 0.002 | 94 | 37077 | 50.0 | 44.9 | |
| 96 2-Chlorobenzotrifluoride | 180 | 11.345 | 11.343 | 0.002 | 96 | 152607 | 50.0 | 61.3 | |
| 97 Isopropylbenzene | 105 | 11.436 | 11.441 | -0.005 | 96 | 410317 | 50.0 | 52.6 | |
| 100 Bromobenzene | 156 | 11.752 | 11.751 | 0.001 | 95 | 95235 | 50.0 | 44.0 | |
| 99 1,1,2,2-Tetrachloroethane | 83 | 11.752 | 11.751 | 0.001 | 86 | 121341 | 50.0 | 52.1 | |
| 102 trans-1,4-Dichloro-2-buten | 53 | 11.789 | 11.787 | 0.002 | 72 | 33910 | 50.0 | 52.0 | |
| 101 1,2,3-Trichloropropane | 110 | 11.807 | 11.799 | 0.008 | 87 | 42509 | 50.0 | 47.6 | |
| 103 N-Propylbenzene | 120 | 11.856 | 11.854 | 0.002 | 99 | 116124 | 50.0 | 47.0 | |
| 104 2-Chlorotoluene | 126 | 11.941 | 11.939 | 0.002 | 96 | 99698 | 50.0 | 46.6 | |
| 105 3-Chlorotoluene | 126 | 12.008 | 12.006 | 0.002 | 97 | 126202 | 50.0 | 54.3 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 106 1,3,5-Trimethylbenzene | 105 | 12.038 | 12.037 | 0.001 | 95 | 346189 | 50.0 | 48.9 | |
| 107 4-Chlorotoluene | 126 | 12.063 | 12.061 | 0.002 | 98 | 109975 | 50.0 | 47.6 | |
| 108 tert-Butylbenzene | 119 | 12.349 | 12.353 | -0.004 | 94 | 267680 | 50.0 | 45.2 | |
| 110 1,2,4-Trimethylbenzene | 105 | 12.409 | 12.408 | 0.001 | 97 | 336477 | 50.0 | 46.8 | |
| 111 1,2-dichloro-4-(trifluorom | 214 | 12.452 | 12.450 | 0.002 | 97 | 95957 | 50.0 | 53.2 | |
| 112 sec-Butylbenzene | 105 | 12.574 | 12.572 | 0.002 | 94 | 392992 | 50.0 | 47.6 | |
| 113 1,3-Dichlorobenzene | 146 | 12.689 | 12.694 | -0.005 | 97 | 181959 | 50.0 | 47.1 | |
| 114 4-Isopropyltoluene | 119 | 12.732 | 12.730 | 0.002 | 97 | 326742 | 50.0 | 47.5 | |
| 115 1,4-Dichlorobenzene | 146 | 12.793 | 12.797 | -0.004 | 95 | 186990 | 50.0 | 47.1 | |
| 116 2,4-Dichloro-1-(trifluorom | 214 | 12.823 | 12.822 | 0.001 | 95 | 86163 | 50.0 | 51.4 | |
| 118 2,5-Dichlorobenzotrifluori | 214 | 12.866 | 12.864 | 0.002 | 0 | 97154 | 50.0 | 53.6 | |
| 120 n-Butylbenzene | 91 | 13.139 | 13.138 | 0.001 | 98 | 255354 | 50.0 | 45.5 | |
| 121 1,2-Dichlorobenzene | 146 | 13.152 | 13.150 | 0.002 | 96 | 174412 | 50.0 | 47.3 | |
| 122 1,2-Dibromo-3-Chloropropan | 75 | 13.942 | 13.947 | -0.005 | 77 | 17569 | 50.0 | 42.9 | |
| 123 2,4- & 2,5- & 2,6- Dichlor | 125 | 14.082 | 14.081 | 0.001 | 0 | 361209 | 150.0 | 154.5 | |
| 125 2,3- & 3,4- Dichlorotoluen | 125 | 14.502 | 14.501 | 0.001 | 0 | 234784 | 100.0 | 97.1 | |
| 126 1,2,4-Trichlorobenzene | 180 | 14.764 | 14.768 | -0.004 | 94 | 68840 | 50.0 | 40.8 | |
| 127 Hexachlorobutadiene | 225 | 14.910 | 14.908 | 0.002 | 97 | 27914 | 50.0 | 45.2 | |
| 128 Naphthalene | 128 | 15.031 | 15.030 | 0.001 | 97 | 212015 | 50.0 | 36.9 | |
| 129 1,2,3-Trichlorobenzene | 180 | 15.256 | 15.255 | 0.001 | 96 | 60881 | 50.0 | 39.5 | |
| 131 2,4,5-Trichlorotoluene | 159 | 16.023 | 16.027 | -0.004 | 0 | 24923 | 50.0 | 34.0 | |
| 130 2,3,6-Trichlorotoluene | 159 | 16.120 | 16.125 | -0.005 | 98 | 23556 | 50.0 | 34.6 | |
| 149 3,4-Dichlorotoluene | 1 | | 0.000 | | | | ND | ND | |
| S 133 Xylenes, Total | 106 | | | | 0 | | 100.0 | 101.0 | |
| S 134 1,2-Dichloroethene, Total | 96 | | | | 0 | | 100.0 | 99.3 | |
| S 135 1,3-Dichloropropene, Total | 1 | | | | 0 | | 100.0 | 93.4 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Reagents:

| | | | |
|---------------------|--------------------|-----------|-------------|
| voaWEEmix1stR_00011 | Amount Added: 2.00 | Units: uL | |
| voaWKetmix1st_00006 | Amount Added: 2.00 | Units: uL | |
| voaWVA1stRest_00021 | Amount Added: 2.00 | Units: uL | |
| VOA8260VOA2ND_00267 | Amount Added: 2.00 | Units: uL | |
| VOA2CEVE2ND_00008 | Amount Added: 2.00 | Units: uL | |
| voaWAcro1stRe_00021 | Amount Added: 6.00 | Units: uL | |
| VOA8260INT_00074 | Amount Added: 2.00 | Units: uL | Run Reagent |
| VOA8260SURR_00073 | Amount Added: 2.00 | Units: uL | Run Reagent |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D04.D

Injection Date: 04-Oct-2017 01:24:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: LCS

Worklist Smp#: 4

Client ID:

Purge Vol: 5.000 mL

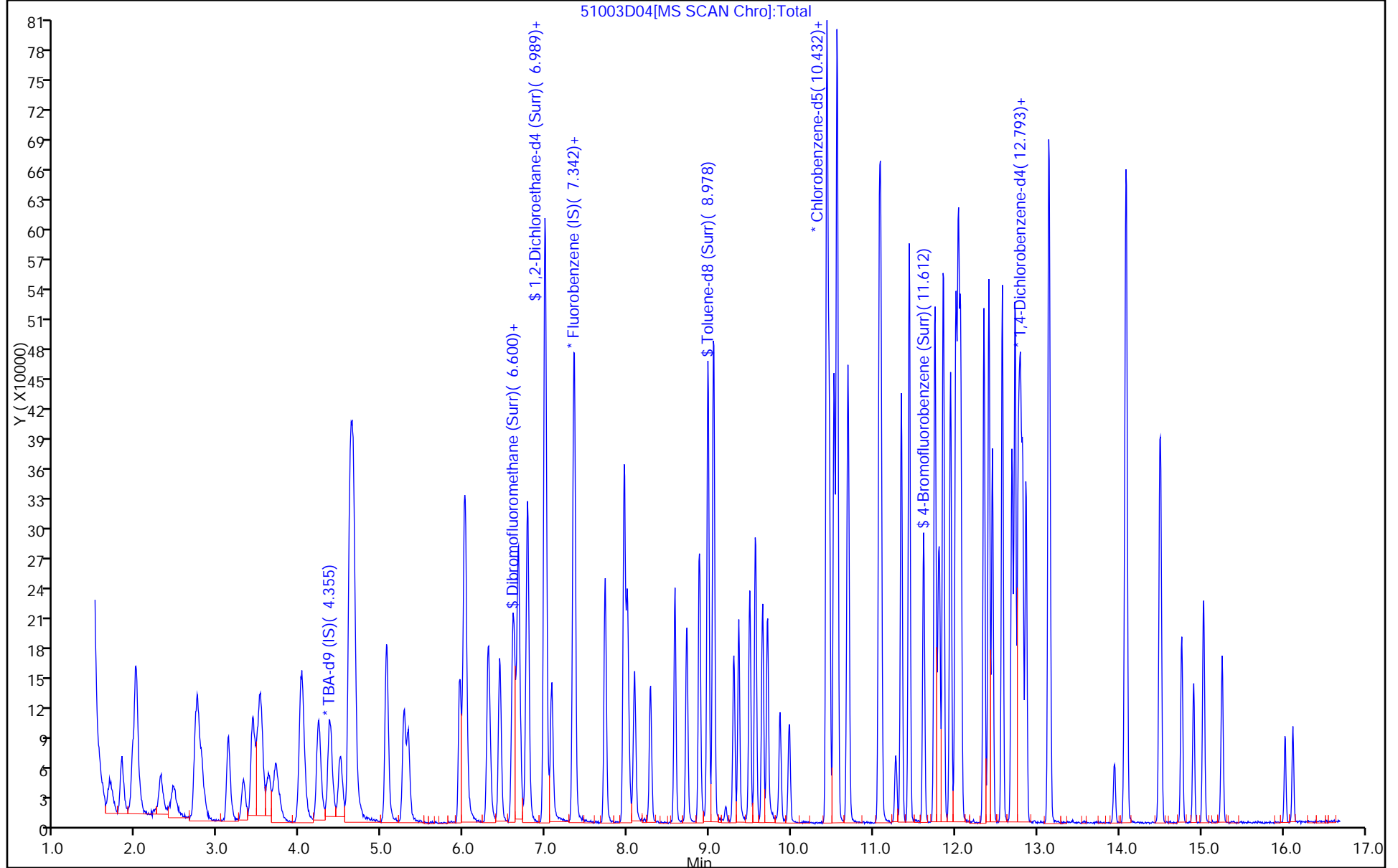
Dil. Factor: 1.0000

ALS Bottle#: 4

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Recovery Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\51003D04.D
 Lims ID: LCS
 Client ID:
 Sample Type: LCS
 Inject. Date: 04-Oct-2017 01:24:30 ALS Bottle#: 4 Worklist Smp#: 4
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018710-004
 Misc. Info.: LCS
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171003-18710.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 04-Oct-2017 21:10:21 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK005

First Level Reviewer: bungardf Date: 04-Oct-2017 02:09:21

| Compound | Amount Added | Amount Recovered | % Rec. |
|-----------------------------------|--------------|------------------|--------|
| \$ 5 Dibromofluoromethane (Surr) | 50.0 | 49.0 | 98.03 |
| \$ 6 1,2-Dichloroethane-d4 (Surr) | 50.0 | 53.1 | 106.12 |
| \$ 7 Toluene-d8 (Surr) | 50.0 | 56.7 | 113.38 |
| \$ 8 4-Bromofluorobenzene (Surr) | 50.0 | 53.7 | 107.39 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Client Sample ID: _____ Lab Sample ID: LCS 180-224919/4
 Matrix: Water Lab File ID: 51004D04.D
 Analysis Method: 8260C Date Collected: _____
 Sample wt/vol: 5 (mL) Date Analyzed: 10/05/2017 01:09
 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.: 224919 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|------------|-----------------------------|--------|---|-----|------|
| 74-87-3 | Chloromethane | 12.0 | | 1.0 | 0.38 |
| 75-01-4 | Vinyl chloride | 12.6 | | 1.0 | 0.17 |
| 74-83-9 | Bromomethane | 12.4 | | 1.0 | 0.59 |
| 75-00-3 | Chloroethane | 12.5 | | 1.0 | 0.58 |
| 75-35-4 | 1,1-Dichloroethene | 9.95 | | 1.0 | 0.32 |
| 67-64-1 | Acetone | 21.7 | | 5.0 | 3.1 |
| 75-15-0 | Carbon disulfide | 9.04 | | 1.0 | 0.53 |
| 75-09-2 | Methylene Chloride | 8.90 | | 1.0 | 0.94 |
| 156-60-5 | trans-1,2-Dichloroethene | 9.52 | | 1.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 9.23 | | 1.0 | 0.20 |
| 75-34-3 | 1,1-Dichloroethane | 9.51 | | 1.0 | 0.34 |
| 156-59-2 | cis-1,2-Dichloroethene | 8.65 | | 1.0 | 0.30 |
| 74-97-5 | Bromochloromethane | 8.75 | | 1.0 | 0.36 |
| 78-93-3 | 2-Butanone (MEK) | 18.3 | | 5.0 | 2.6 |
| 67-66-3 | Chloroform | 9.05 | | 1.0 | 0.27 |
| 71-55-6 | 1,1,1-Trichloroethane | 9.46 | | 1.0 | 0.27 |
| 56-23-5 | Carbon tetrachloride | 9.46 | | 1.0 | 0.56 |
| 71-43-2 | Benzene | 8.85 | | 1.0 | 0.18 |
| 107-06-2 | 1,2-Dichloroethane | 9.68 | | 1.0 | 0.24 |
| 79-01-6 | Trichloroethene | 8.20 | | 1.0 | 0.20 |
| 78-87-5 | 1,2-Dichloropropane | 8.44 | | 1.0 | 0.35 |
| 75-27-4 | Bromodichloromethane | 8.21 | | 1.0 | 0.57 |
| 10061-01-5 | cis-1,3-Dichloropropene | 8.18 | | 1.0 | 0.32 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 20.4 | | 5.0 | 2.2 |
| 108-88-3 | Toluene | 10.4 | | 1.0 | 0.16 |
| 10061-02-6 | trans-1,3-Dichloropropene | 9.78 | | 1.0 | 0.22 |
| 79-00-5 | 1,1,2-Trichloroethane | 10.6 | | 1.0 | 0.31 |
| 127-18-4 | Tetrachloroethene | 9.77 | | 1.0 | 0.24 |
| 591-78-6 | 2-Hexanone | 19.4 | | 5.0 | 2.0 |
| 124-48-1 | Dibromochloromethane | 9.57 | | 1.0 | 0.44 |
| 106-93-4 | 1,2-Dibromoethane (EDB) | 9.26 | | 1.0 | 0.51 |
| 108-90-7 | Chlorobenzene | 9.72 | | 1.0 | 0.15 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 9.84 | | 1.0 | 0.49 |
| 100-41-4 | Ethylbenzene | 9.31 | | 1.0 | 0.25 |
| 1330-20-7 | Xylenes, Total | 18.9 | | 2.0 | 0.27 |
| 100-42-5 | Styrene | 9.11 | | 1.0 | 0.22 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1
 SDG No.: _____
 Client Sample ID: _____ Lab Sample ID: LCS 180-224919/4
 Matrix: Water Lab File ID: 51004D04.D
 Analysis Method: 8260C Date Collected: _____
 Sample wt/vol: 5 (mL) Date Analyzed: 10/05/2017 01:09
 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.: 224919 Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|----------|---------------------------|--------|---|-----|------|
| 75-25-2 | Bromoform | 8.75 | | 1.0 | 0.76 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 9.67 | | 1.0 | 0.37 |
| 107-13-1 | Acrylonitrile | 102 | | 20 | 3.3 |
| 123-91-1 | 1,4-Dioxane | 189 | J | 200 | 16 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|------------|------------------------------|------|---|--------|
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 102 | | 65-121 |
| 2037-26-5 | Toluene-d8 (Surr) | 110 | | 73-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 102 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 90 | | 73-120 |

TestAmerica Pittsburgh
Target Compound Quantitation Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D04.D
 Lims ID: LCS
 Client ID:
 Sample Type: LCS
 Inject. Date: 05-Oct-2017 01:09:30 ALS Bottle#: 4 Worklist Smp#: 4
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018725-004
 Misc. Info.: LCS
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 08-Oct-2017 20:41:20 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK026

First Level Reviewer: bungardf

Date: 05-Oct-2017 01:33:14

| Compound | Sig | RT (min.) | Exp RT (min.) | Diff RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|---------------------------------|-----|-----------|---------------|----------------|----|----------|------------|--------------|-------|
| * 1 TBA-d9 (IS) | 65 | 4.365 | 4.374 | -0.009 | 0 | 181022 | 1000.0 | 1000.0 | |
| * 2 Fluorobenzene (IS) | 96 | 7.340 | 7.337 | 0.003 | 97 | 356874 | 50.0 | 50.0 | |
| * 3 Chlorobenzene-d5 | 119 | 10.430 | 10.427 | 0.003 | 86 | 73297 | 50.0 | 50.0 | |
| * 4 1,4-Dichlorobenzene-d4 | 152 | 12.766 | 12.769 | -0.003 | 93 | 101073 | 50.0 | 50.0 | |
| \$ 5 Dibromofluoromethane (Surr | 113 | 6.616 | 6.613 | 0.003 | 92 | 76986 | 50.0 | 44.8 | |
| \$ 6 1,2-Dichloroethane-d4 (Sur | 65 | 6.987 | 6.990 | -0.003 | 0 | 107095 | 50.0 | 51.1 | |
| \$ 7 Toluene-d8 (Surr) | 98 | 8.982 | 8.979 | 0.003 | 93 | 321517 | 50.0 | 55.1 | |
| \$ 8 4-Bromofluorobenzene (Surr | 95 | 11.610 | 11.613 | -0.003 | 85 | 107404 | 50.0 | 51.0 | |
| 11 Dichlorodifluoromethane | 85 | 1.676 | 1.679 | -0.003 | 99 | 96918 | 50.0 | 46.7 | |
| 12 Chloromethane | 50 | 1.822 | 1.825 | -0.003 | 99 | 124731 | 50.0 | 59.8 | |
| 13 Vinyl chloride | 62 | 1.962 | 1.959 | 0.003 | 97 | 133058 | 50.0 | 62.9 | |
| 14 Butadiene | 39 | 1.998 | 2.008 | -0.010 | 96 | 129922 | 50.0 | 67.6 | |
| 15 Bromomethane | 94 | 2.309 | 2.300 | 0.009 | 90 | 62099 | 50.0 | 62.1 | |
| 16 Chloroethane | 64 | 2.455 | 2.470 | -0.015 | 99 | 72511 | 50.0 | 62.3 | |
| 17 Dichlorofluoromethane | 67 | 2.747 | 2.744 | 0.003 | 97 | 176398 | 50.0 | 59.9 | |
| 18 Trichlorofluoromethane | 101 | 2.765 | 2.768 | -0.003 | 92 | 133128 | 50.0 | 51.2 | |
| 20 Ethyl ether | 59 | 3.124 | 3.121 | 0.003 | 92 | 82872 | 50.0 | 49.0 | |
| 21 Acrolein | 56 | 3.319 | 3.316 | 0.003 | 97 | 67213 | 150.0 | 157.7 | |
| 22 1,1-Dichloroethene | 96 | 3.416 | 3.413 | 0.003 | 96 | 86923 | 50.0 | 49.8 | |
| 23 1,1,2-Trichloro-1,2,2-trif | 101 | 3.501 | 3.504 | -0.003 | 95 | 88802 | 50.0 | 46.3 | |
| 24 Acetone | 43 | 3.532 | 3.529 | 0.003 | 99 | 101281 | 100.0 | 108.5 | |
| 25 Iodomethane | 142 | 3.623 | 3.608 | 0.015 | 96 | 124276 | 50.0 | 45.3 | |
| 26 Carbon disulfide | 76 | 3.714 | 3.699 | 0.015 | 99 | 173289 | 50.0 | 45.2 | |
| 28 3-Chloro-1-propene | 76 | 4.012 | 3.997 | 0.015 | 97 | 50237 | 50.0 | 44.5 | |
| 30 Methyl acetate | 43 | 4.030 | 4.033 | -0.003 | 98 | 190730 | 100.0 | 103.2 | |
| 31 Methylene Chloride | 84 | 4.219 | 4.222 | -0.003 | 92 | 97064 | 50.0 | 44.5 | |
| 32 2-Methyl-2-propanol | 59 | 4.493 | 4.508 | -0.015 | 96 | 99506 | 500.0 | 464.8 | |
| 33 Acrylonitrile | 53 | 4.608 | 4.605 | 0.003 | 99 | 456048 | 500.0 | 507.5 | |
| 34 trans-1,2-Dichloroethene | 96 | 4.639 | 4.642 | -0.003 | 79 | 94726 | 50.0 | 47.6 | |
| 35 Methyl tert-butyl ether | 73 | 4.651 | 4.660 | -0.009 | 97 | 246267 | 50.0 | 46.1 | |
| 36 Hexane | 57 | 5.065 | 5.049 | 0.016 | 95 | 112742 | 50.0 | 44.1 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|--------------------------------|-----|-----------|---------------|---------------|-----|----------|------------|--------------|-------|
| 37 1,1-Dichloroethane | 63 | 5.271 | 5.268 | 0.003 | 96 | 164517 | 50.0 | 47.5 | |
| 38 Vinyl acetate | 43 | 5.320 | 5.323 | -0.003 | 97 | 170116 | 50.0 | 48.3 | |
| 44 2,2-Dichloropropane | 97 | 6.014 | 6.004 | 0.010 | 63 | 20953 | 50.0 | 47.5 | |
| 45 cis-1,2-Dichloroethene | 96 | 6.014 | 6.011 | 0.003 | 82 | 98469 | 50.0 | 43.2 | |
| 46 2-Butanone (MEK) | 43 | 6.026 | 6.023 | 0.003 | 100 | 121500 | 100.0 | 91.5 | |
| 49 Chlorobromomethane | 128 | 6.299 | 6.290 | 0.009 | 98 | 44281 | 50.0 | 43.8 | |
| 51 Tetrahydrofuran | 42 | 6.312 | 6.309 | 0.003 | 92 | 65444 | 100.0 | 84.6 | |
| 52 Chloroform | 83 | 6.439 | 6.436 | 0.003 | 93 | 156442 | 50.0 | 45.3 | |
| 53 1,1,1-Trichloroethane | 97 | 6.591 | 6.595 | -0.004 | 97 | 123765 | 50.0 | 47.3 | |
| 54 Cyclohexane | 56 | 6.664 | 6.662 | 0.002 | 92 | 147819 | 50.0 | 45.8 | |
| 56 Carbon tetrachloride | 117 | 6.768 | 6.759 | 0.009 | 94 | 103011 | 50.0 | 47.3 | |
| 55 1,1-Dichloropropene | 75 | 6.780 | 6.783 | -0.003 | 96 | 123461 | 50.0 | 43.7 | |
| 57 Isobutyl alcohol | 41 | 6.981 | 6.990 | -0.009 | 63 | 90778 | 1250.0 | 1278.4 | |
| 58 Benzene | 78 | 6.999 | 6.996 | 0.003 | 97 | 383948 | 50.0 | 44.2 | |
| 59 1,2-Dichloroethane | 62 | 7.066 | 7.069 | -0.003 | 98 | 122378 | 50.0 | 48.4 | |
| 62 n-Heptane | 43 | 7.352 | 7.349 | 0.003 | 87 | 93326 | 50.0 | 45.7 | |
| 64 Trichloroethene | 130 | 7.723 | 7.720 | 0.003 | 98 | 89505 | 50.0 | 41.0 | |
| 66 Methylcyclohexane | 83 | 7.954 | 7.957 | -0.003 | 91 | 132103 | 50.0 | 40.0 | |
| 67 1,2-Dichloropropane | 63 | 7.997 | 7.994 | 0.003 | 95 | 85303 | 50.0 | 42.2 | |
| 68 Dibromomethane | 93 | 8.088 | 8.079 | 0.009 | 96 | 49915 | 50.0 | 42.2 | |
| 70 1,4-Dioxane | 88 | 8.088 | 8.085 | 0.003 | 46 | 19395 | 1000.0 | 944.0 | |
| 71 Dichlorobromomethane | 83 | 8.277 | 8.280 | -0.003 | 99 | 95397 | 50.0 | 41.0 | |
| 73 2-Chloroethyl vinyl ether | 63 | 8.575 | 8.578 | -0.003 | 94 | 90727 | 100.0 | 62.4 | |
| 74 cis-1,3-Dichloropropene | 75 | 8.721 | 8.718 | 0.003 | 94 | 115387 | 50.0 | 40.9 | |
| 75 4-Methyl-2-pentanone (MIBK) | 43 | 8.873 | 8.876 | -0.003 | 98 | 191631 | 100.0 | 101.9 | |
| 76 Toluene | 91 | 9.049 | 9.046 | 0.003 | 98 | 378741 | 50.0 | 51.8 | |
| 77 trans-1,3-Dichloropropene | 75 | 9.293 | 9.296 | -0.003 | 93 | 97217 | 50.0 | 48.9 | |
| 78 Ethyl methacrylate | 69 | 9.353 | 9.356 | -0.003 | 90 | 98259 | 50.0 | 41.0 | |
| 79 1,1,2-Trichloroethane | 97 | 9.487 | 9.490 | -0.003 | 89 | 81055 | 50.0 | 53.2 | |
| 80 Tetrachloroethene | 164 | 9.554 | 9.563 | -0.009 | 96 | 68116 | 50.0 | 48.9 | |
| 81 1,3-Dichloropropane | 76 | 9.645 | 9.648 | -0.003 | 92 | 131293 | 50.0 | 46.6 | |
| 82 2-Hexanone | 43 | 9.706 | 9.703 | 0.003 | 99 | 139888 | 100.0 | 97.0 | |
| 84 Chlorodibromomethane | 129 | 9.858 | 9.855 | 0.003 | 88 | 61567 | 50.0 | 47.8 | |
| 85 Ethylene Dibromide | 107 | 9.968 | 9.971 | -0.003 | 98 | 72271 | 50.0 | 46.3 | |
| 86 3-Chlorobenzotrifluoride | 180 | 10.430 | 10.433 | -0.003 | 91 | 128526 | 50.0 | 51.0 | |
| 87 Chlorobenzene | 112 | 10.461 | 10.458 | 0.003 | 94 | 231138 | 50.0 | 48.6 | |
| 88 4-Chlorobenzotrifluoride | 180 | 10.515 | 10.518 | -0.003 | 95 | 120347 | 50.0 | 51.8 | |
| 89 1,1,1,2-Tetrachloroethane | 131 | 10.552 | 10.549 | 0.003 | 92 | 74457 | 50.0 | 49.2 | |
| 90 Ethylbenzene | 106 | 10.558 | 10.555 | 0.003 | 98 | 123640 | 50.0 | 46.5 | |
| 91 m-Xylene & p-Xylene | 106 | 10.692 | 10.689 | 0.003 | 0 | 152835 | 50.0 | 47.1 | |
| 92 o-Xylene | 106 | 11.075 | 11.072 | 0.003 | 96 | 146737 | 50.0 | 47.4 | |
| 93 Styrene | 104 | 11.087 | 11.090 | -0.003 | 94 | 238407 | 50.0 | 45.5 | |
| 94 Bromoform | 173 | 11.270 | 11.273 | -0.003 | 94 | 35008 | 50.0 | 43.8 | |
| 96 2-Chlorobenzotrifluoride | 180 | 11.343 | 11.340 | 0.003 | 96 | 124136 | 50.0 | 51.5 | |
| 97 Isopropylbenzene | 105 | 11.434 | 11.437 | -0.003 | 96 | 374831 | 50.0 | 49.7 | |
| 99 1,1,2,2-Tetrachloroethane | 83 | 11.750 | 11.747 | 0.003 | 85 | 109038 | 50.0 | 48.4 | |
| 100 Bromobenzene | 156 | 11.750 | 11.753 | -0.003 | 95 | 84119 | 50.0 | 42.9 | |
| 102 trans-1,4-Dichloro-2-buten | 53 | 11.787 | 11.784 | 0.003 | 71 | 31722 | 50.0 | 53.6 | |
| 101 1,2,3-Trichloropropane | 110 | 11.805 | 11.808 | -0.003 | 85 | 37075 | 50.0 | 45.8 | |
| 103 N-Propylbenzene | 120 | 11.854 | 11.851 | 0.003 | 99 | 101012 | 50.0 | 45.1 | |
| 104 2-Chlorotoluene | 126 | 11.945 | 11.942 | 0.003 | 96 | 87718 | 50.0 | 45.3 | |
| 105 3-Chlorotoluene | 126 | 12.006 | 12.003 | 0.003 | 97 | 105384 | 50.0 | 50.0 | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | Cal Amt ng | OnCol Amt ng | Flags |
|----------------------------------|-----|-----------|---------------|---------------|----|----------|------------|--------------|-------|
| 106 1,3,5-Trimethylbenzene | 105 | 12.036 | 12.033 | 0.003 | 94 | 305642 | 50.0 | 47.7 | |
| 107 4-Chlorotoluene | 126 | 12.067 | 12.064 | 0.003 | 96 | 97339 | 50.0 | 46.5 | |
| 108 tert-Butylbenzene | 119 | 12.353 | 12.350 | 0.003 | 93 | 231203 | 50.0 | 43.1 | |
| 110 1,2,4-Trimethylbenzene | 105 | 12.407 | 12.410 | -0.003 | 97 | 290686 | 50.0 | 44.6 | |
| 111 1,2-dichloro-4-(trifluorom | 214 | 12.450 | 12.453 | -0.003 | 96 | 73226 | 50.0 | 44.8 | |
| 112 sec-Butylbenzene | 105 | 12.572 | 12.575 | -0.003 | 94 | 342058 | 50.0 | 45.7 | |
| 113 1,3-Dichlorobenzene | 146 | 12.693 | 12.690 | 0.003 | 97 | 156593 | 50.0 | 44.7 | |
| 114 4-Isopropyltoluene | 119 | 12.730 | 12.727 | 0.003 | 96 | 286076 | 50.0 | 45.9 | |
| 115 1,4-Dichlorobenzene | 146 | 12.797 | 12.794 | 0.003 | 95 | 163897 | 50.0 | 45.5 | |
| 116 2,4-Dichloro-1-(trifluorom | 214 | 12.821 | 12.818 | 0.003 | 93 | 71502 | 50.0 | 47.0 | |
| 118 2,5-Dichlorobenzotrifluori | 214 | 12.864 | 12.867 | -0.003 | 0 | 72679 | 50.0 | 44.2 | |
| 120 n-Butylbenzene | 91 | 13.137 | 13.140 | -0.003 | 98 | 229550 | 50.0 | 45.2 | |
| 121 1,2-Dichlorobenzene | 146 | 13.149 | 13.153 | -0.004 | 95 | 148873 | 50.0 | 44.6 | |
| 122 1,2-Dibromo-3-Chloropropan | 75 | 13.940 | 13.937 | 0.003 | 77 | 16109 | 50.0 | 43.4 | |
| 123 2,4- & 2,5- & 2,6- Dichlor | 125 | 14.080 | 14.083 | -0.003 | 0 | 300006 | 150.0 | 141.6 | |
| 125 2,3- & 3,4- Dichlorotoluen | 125 | 14.500 | 14.503 | -0.003 | 0 | 201265 | 100.0 | 91.8 | |
| 126 1,2,4-Trichlorobenzene | 180 | 14.768 | 14.765 | 0.003 | 93 | 59098 | 50.0 | 38.7 | |
| 127 Hexachlorobutadiene | 225 | 14.914 | 14.905 | 0.009 | 95 | 24081 | 50.0 | 43.0 | |
| 128 Naphthalene | 128 | 15.029 | 15.026 | 0.003 | 97 | 195929 | 50.0 | 37.6 | |
| 129 1,2,3-Trichlorobenzene | 180 | 15.260 | 15.257 | 0.003 | 96 | 48438 | 50.0 | 34.7 | |
| 131 2,4,5-Trichlorotoluene | 159 | 16.027 | 16.024 | 0.003 | 0 | 23903 | 50.0 | 36.0 | |
| 130 2,3,6-Trichlorotoluene | 159 | 16.118 | 16.121 | -0.003 | 97 | 23680 | 50.0 | 38.3 | |
| 149 3,4-Dichlorotoluene | 1 | | 0.000 | | | | ND | ND | |
| S 133 Xylenes, Total | 106 | | | | 0 | | 100.0 | 94.5 | |
| S 134 1,2-Dichloroethene, Total | 96 | | | | 0 | | 100.0 | 90.8 | |
| S 135 1,3-Dichloropropene, Total | 1 | | | | 0 | | 100.0 | 89.8 | |

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Reagents:

| | | | |
|---------------------|--------------------|-----------|-------------|
| voaWEEmix1stR_00011 | Amount Added: 2.00 | Units: uL | |
| voaWKetmix1st_00006 | Amount Added: 2.00 | Units: uL | |
| voaWVA1stRest_00021 | Amount Added: 2.00 | Units: uL | |
| VOA8260VOA2ND_00267 | Amount Added: 2.00 | Units: uL | |
| VOA2CEVE2ND_00008 | Amount Added: 2.00 | Units: uL | |
| voaWAcro1stRe_00021 | Amount Added: 6.00 | Units: uL | |
| VOA8260INT_00074 | Amount Added: 2.00 | Units: uL | Run Reagent |
| VOA8260SURR_00073 | Amount Added: 2.00 | Units: uL | Run Reagent |

TestAmerica Pittsburgh

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D04.D

Injection Date: 05-Oct-2017 01:09:30

Instrument ID: CHHP5

Operator ID: 034635

Lims ID: LCS

Worklist Smp#: 4

Client ID:

Purge Vol: 5.000 mL

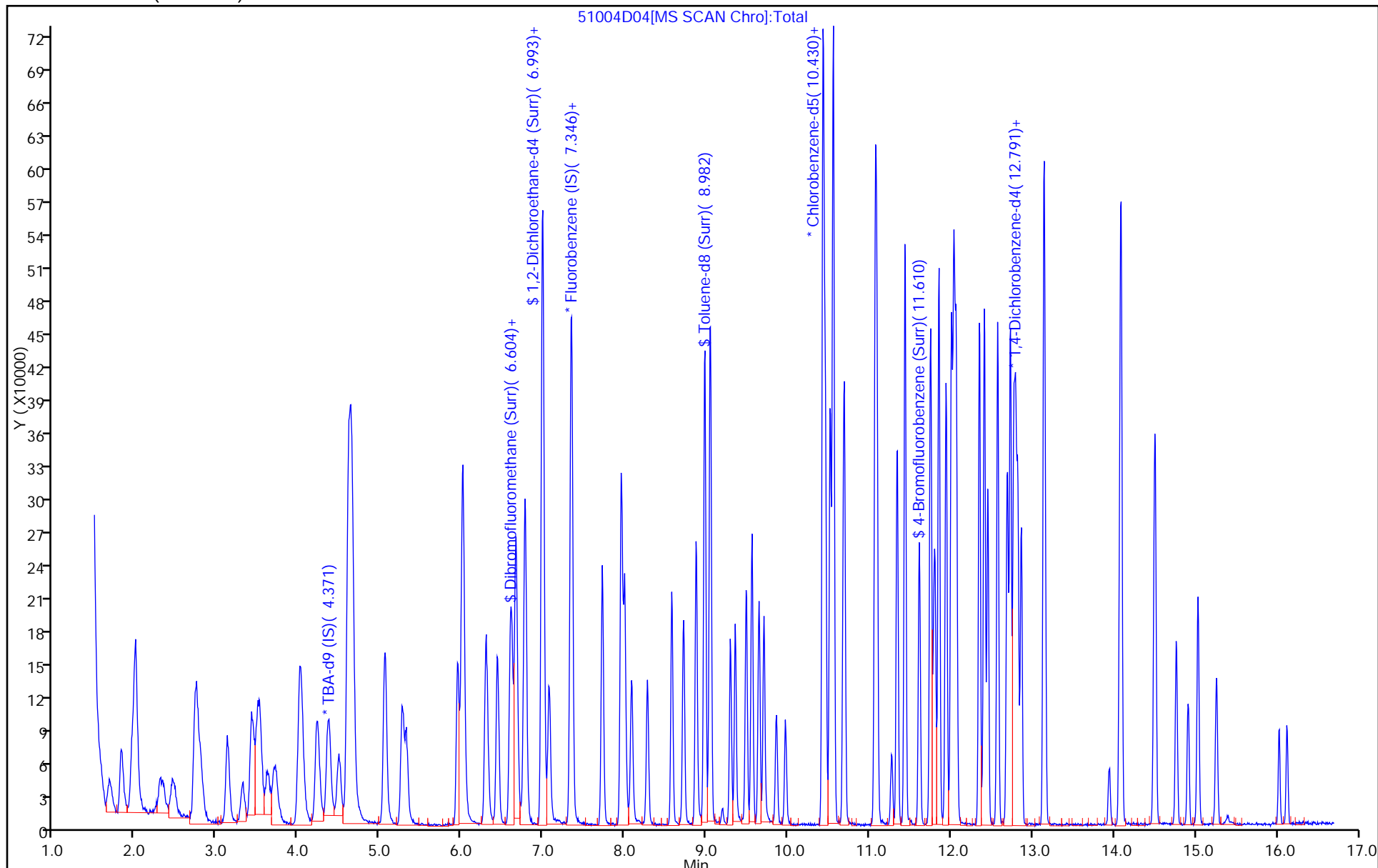
Dil. Factor: 1.0000

ALS Bottle#: 4

Method: MSVOA_LL_CHHP5

Limit Group: VOA 8260C ICAL

Column: DB-624 (0.18 mm)



TestAmerica Pittsburgh
Recovery Report

Data File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\51004D04.D
 Lims ID: LCS
 Client ID:
 Sample Type: LCS
 Inject. Date: 05-Oct-2017 01:09:30 ALS Bottle#: 4 Worklist Smp#: 4
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 180-0018725-004
 Misc. Info.: LCS
 Operator ID: 034635 Instrument ID: CHHP5
 Method: \\ChromNA\Pittsburgh\ChromData\CHHP5\20171004-18725.b\MSVOA_LL_CHHP5.m
 Limit Group: VOA 8260C ICAL
 Last Update: 08-Oct-2017 20:41:20 Calib Date: 27-Jul-2017 04:24:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Pittsburgh\ChromData\CHHP5\20170726-17756.b\50727D11.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK026

First Level Reviewer: bungardf

Date: 05-Oct-2017 01:33:14

| Compound | Amount Added | Amount Recovered | % Rec. |
|-----------------------------------|--------------|------------------|--------|
| \$ 5 Dibromofluoromethane (Surr) | 50.0 | 44.8 | 89.67 |
| \$ 6 1,2-Dichloroethane-d4 (Surr) | 50.0 | 51.1 | 102.27 |
| \$ 7 Toluene-d8 (Surr) | 50.0 | 55.1 | 110.23 |
| \$ 8 4-Bromofluorobenzene (Surr) | 50.0 | 51.0 | 101.96 |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1

SDG No.: _____

Instrument ID: CHHP5 Start Date: 07/27/2017 00:22

Analysis Batch Number: 218218 End Date: 07/27/2017 05:50

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|-------------------|------------------|------------------|-----------------|-------------|------------------|
| BFB 180-218218/1 | | 07/27/2017 00:22 | 1 | 50727D01.D | DB-624 0.18 (mm) |
| IC 180-218218/2 | | 07/27/2017 00:51 | 1 | 50727D02.D | DB-624 0.18 (mm) |
| IC 180-218218/3 | | 07/27/2017 01:15 | 1 | 50727D03.D | DB-624 0.18 (mm) |
| ICIS 180-218218/4 | | 07/27/2017 01:39 | 1 | 50727D04.D | DB-624 0.18 (mm) |
| ZZZZZ | | 07/27/2017 01:39 | 1 | | DB-624 0.18 (mm) |
| IC 180-218218/5 | | 07/27/2017 02:02 | 1 | 50727D05.D | DB-624 0.18 (mm) |
| IC 180-218218/6 | | 07/27/2017 02:26 | 1 | 50727D06.D | DB-624 0.18 (mm) |
| IC 180-218218/8 | | 07/27/2017 03:13 | 1 | 50727D08.D | DB-624 0.18 (mm) |
| IC 180-218218/10 | | 07/27/2017 04:00 | 1 | 50727D10.D | DB-624 0.18 (mm) |
| IC 180-218218/11 | | 07/27/2017 04:24 | 1 | 50727D11.D | DB-624 0.18 (mm) |
| ICV 180-218218/12 | | 07/27/2017 05:03 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 07/27/2017 05:50 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 07/27/2017 05:50 | 1 | | DB-624 0.18 (mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica PittsburghJob No.: 180-70873-1

SDG No.: _____

Instrument ID: CHHP5Start Date: 10/03/2017 23:49Analysis Batch Number: 224792End Date: 10/04/2017 11:24

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|--------------------|------------------|------------------|-----------------|-------------|------------------|
| BFB 180-224792/1 | | 10/03/2017 23:49 | 1 | 51003D01.D | DB-624 0.18 (mm) |
| CCVIS 180-224792/2 | | 10/04/2017 00:22 | 1 | 51003D02.D | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 00:22 | 1 | | DB-624 0.18 (mm) |
| CCV 180-224792/3 | | 10/04/2017 01:00 | 1 | | DB-624 0.18 (mm) |
| LCS 180-224792/4 | | 10/04/2017 01:24 | 1 | 51003D04.D | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 01:57 | 1 | | DB-624 0.18 (mm) |
| MB 180-224792/6 | | 10/04/2017 02:21 | 1 | 51003D06.D | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 02:54 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 03:23 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 03:47 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 04:35 | 1 | | DB-624 0.18 (mm) |
| 180-70873-2 | | 10/04/2017 04:59 | 1 | 51003D12.D | DB-624 0.18 (mm) |
| 180-70873-1 | | 10/04/2017 05:23 | 1 | 51003D13.D | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 06:58 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 07:46 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 08:11 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 08:34 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 08:59 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 09:47 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 10:11 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 10:35 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 10:59 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 11:24 | 1 | | DB-624 0.18 (mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Pittsburgh Job No.: 180-70873-1

SDG No.: _____

Instrument ID: CHHP5 Start Date: 10/04/2017 22:24

Analysis Batch Number: 224919 End Date: 10/05/2017 06:31

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|--------------------|------------------|------------------|-----------------|-------------|------------------|
| BFB 180-224919/1 | | 10/04/2017 22:24 | 1 | 51004D01.D | DB-624 0.18 (mm) |
| CCVIS 180-224919/2 | | 10/04/2017 23:29 | 1 | 51004D02.D | DB-624 0.18 (mm) |
| ZZZZZ | | 10/04/2017 23:29 | 1 | | DB-624 0.18 (mm) |
| LCS 180-224919/4 | | 10/05/2017 01:09 | 1 | 51004D04.D | DB-624 0.18 (mm) |
| CCV 180-224919/5 | | 10/05/2017 01:43 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/05/2017 02:17 | 1 | | DB-624 0.18 (mm) |
| MB 180-224919/7 | | 10/05/2017 02:41 | 1 | 51004D07.D | DB-624 0.18 (mm) |
| ZZZZZ | | 10/05/2017 03:15 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/05/2017 03:41 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/05/2017 04:55 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/05/2017 05:19 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/05/2017 05:43 | 1 | | DB-624 0.18 (mm) |
| ZZZZZ | | 10/05/2017 06:07 | 1 | | DB-624 0.18 (mm) |
| 180-70873-1 DL | | 10/05/2017 06:31 | 12.5 | 51004D16.D | DB-624 0.18 (mm) |

Shipping and Receiving Documents

TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

777 New Durham Road
Edison, New Jersey 08817
Phone: (732) 549-3900 Fax: (732) 549-3679

CHAIN OF CUSTODY / ANALYSIS REQUEST

Page 1 of 1

| | | | |
|---|---------|---|-------|
| Name (for report and invoice) Chris O'Neil / Carrie Gamber | | Site/Project Identification Hurley - Davidson | |
| Company Groundwaterservices Cwp | | State (Location of site): NJ: <input type="checkbox"/> NY: <input type="checkbox"/> Other: PA | |
| Address 2601 market place SE310 | | Regulatory Program: DKQP: <input type="checkbox"/> | |
| City Harrisburg | | LAB USE ONLY Project No: | |
| Phone 717 652 6632 | | Job No: | |
| State PA 17110 | | Sample Numbers (72 hours sample) Trip Blank | |
| Sample Identification | | ANALYSIS REQUESTED (ENTER X BELOW TO INDICATE REQUEST) | |
| Date | | No. of | |
| HD-SPBA-CW-22-0/1-0 | 9/29/17 | Matrix | Cont. |
| HD-QCH-0/1/2 | 9/29/17 | W | 3 |
| | | W | 2 |
| Preservation Used: 1 = ICE, 2 = HCl, 3 = H ₂ SO ₄ , 4 = HNO ₃ , 5 = NaOH, 6 = Other, 7 = Other | | Soil: <input checked="" type="checkbox"/> Water: <input type="checkbox"/> | |

| | | | |
|---------------------------------------|----------------|---------------------------------|--------------------------------------|
| Special Instructions | | Water Metals Filtered (Yes/No)? | |
| Relinquished by <i>[Signature]</i> | Company GSC | Date / Time 9/17/17 1200 | Received by 1) <i>[Signature]</i> |
| Relinquished by | Company | Date / Time | Received by |
| 2) | | | 2) |
| Relinquished by | Company | Date / Time | Received by |
| 3) | | | 3) |
| Relinquished by | Company | Date / Time | Received by |
| 4) | | | 4) |

Laboratory Certifications: New Jersey (12028), New York (11452), Pennsylvania (68-522), Connecticut (PH-0200), Rhode Island (132), Massachusetts (M-NJ312), North Carolina (No. 578) TAL-0016 (0715)

FedEx Express Package US Airbill

FedEx Tracking Number **81116 7637 1538**

From **10/01/13** Date

Sender's Name **Kirsten Fleming** Phone **717 652-6832**

Company **Greenwater Sciences Corp.**

Address **2601 market place ste 310**

City **Harrisburg** State **PA** ZIP **17110**

2 Your Internal Billing Reference

To Recipients Name **Sampe Reeling** Phone **412 963-7058**

Company **Test America Pittsburgh**

Address **301 Alpha Drive**

Address **0A**

Uncorrected temp **2.4** °C
Thermometer ID **13**

CF **0** Initials **TJ**

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

Form ID No. **0200**

4 Express Package Service * To most locations.

Next Business Day

FedEx First Overnight
Earliest next business morning delivery to select locations. Friday shipments will be delivered on Monday unless Saturday Delivery is selected.

FedEx Priority Overnight
Next business morning. * Friday shipments will be delivered on Monday unless Saturday Delivery is selected.

FedEx Standard Overnight
Next business afternoon. * Saturday Delivery NOT available.

Packaging * Declare value June 2008

FedEx Envelope* FedEx Pak* FedEx Box Other FedEx Tube

2 or 3 Business Days

FedEx 2Day AM
Second business morning. * Saturday Delivery NOT available.

FedEx 2Day
Second business afternoon. * Thursday shipments will be delivered on Monday unless Saturday Delivery is selected.

FedEx Express Saver
Third business day. * Saturday Delivery NOT available.

4 Express Package Service * To most locations.

Special Handling and Delivery Signature Options Fees may apply. See the FedEx Service Guide.

Saturday Delivery
NOT available for FedEx Standard Overnight, FedEx 2Day AM, or FedEx Express Saver.

No Signature Required
Package may be left without obtaining a signature for delivery.

Direct Signature
Someone at recipient's address may sign for delivery.

Indirect Signature
Someone at recipient's address may sign for delivery for residential deliveries only.

Does this shipment contain dangerous goods?
One box must be checked.
 NO YES
As per attached Shipper's Declaration, Shipper's Declaration not required.

Restrictions apply for dangerous goods — see the current FedEx Service Guide.

Payment Bill to:

Sender Recipient Third Party Credit Card

Total Packages **Total Weight** **lb.**

Enter FedEx Acct. No. or Credit Card No. below.

Obtain recip. Acct. No. **Check**

Credit Card Auth.

This liability is limited to USD\$500 unless you declare a higher value. See the current FedEx Service Guide for details.



180-70873 Waybill

fedex.com 1.800.GoFedEx 1.800.463.3339

10:30 A 15:38 40:03

R197
FZ B02

fedex.com 1.800.GoFedEx 1.800.463.3339

Login Sample Receipt Checklist

Client: Groundwater Sciences Corporation

Job Number: 180-70873-1

Login Number: 70873
List Number: 1
Creator: Say, Thomas C

List Source: TestAmerica Pittsburgh

| Question | Answer | Comment |
|--|--------|---------|
| Radioactivity wasn't checked or is \leq 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 $<6\text{mm}$ (1/4"). | True | |
| Multiphasic samples are not present. | True | |
| Samples do not require splitting or compositing. | True | |
| Residual Chlorine Checked. | N/A | |