

Atmospheric Analysis & Consulting, Inc.

CLIENT : Soil Water Air Protection Enterprise
PROJECT NAME : Bridgeton Sanitary Landfill Air Quality Assessment
AAC PROJECT NO. : 130597
REPORT DATE : 05/21/2013

On May 17, 2013, Atmospheric Analysis & Consulting, Inc. received four (4) Six-Liter Summa Canisters for Volatile Organic Compounds analysis by EPA method TO-15. Upon receipt each sample was assigned a unique Laboratory ID number as follows:

| Client ID | Lab ID | Return Pressure (mmHga) |
|-----------------|--------------|-------------------------|
| U-1 W8-Canister | 130597-63048 | 507.6 |
| U-2 W6-Canister | 130597-63049 | 316.6 |
| D-1 W4-Canister | 130597-63050 | 614.5 |
| D-2 K-Canister | 130597-63051 | 174.8 |

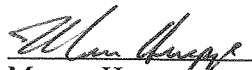
An initial reading of each canister's vacuum was taken and recorded. Subsequently, each canister was brought to positive pressure using UHP-He and the final pressure was recorded.

TO-15 Analysis - Up to a 500 mL aliquot of sample is concentrated, put through a water and CO₂ management system, cryofocused and injected into the GC/MS (full scan mode) for analysis following EPA Method TO-15 as specified in the SOW.

No problems were encountered during receiving, preparation and/ or analysis of these samples. The test results included in this report meet all requirements of the NELAC Standards and/or AAC SOP# TO.15.10.

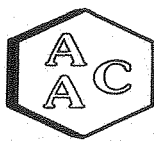
I certify that this data is technically accurate, complete and in compliance with the terms and conditions of the contract. The Laboratory Director or his designee, as verified by the following signature, has authorized the release of the data contained in this hardcopy data package.

If you have any questions or require further explanation of data results, please contact the undersigned.


Marcus Hueppe
Laboratory Director

This report consists of 60 pages.





SAMPLE RECEIPT / LOG-IN REPORT

AAC Project

130597

Received By: J. Zachman

| <u>Sample Receipt Date</u> | <u>Project Desc</u> | <u>Clients ID</u> | <u>Matrix</u> | <u>Sampling Date/Time</u> | <u>Sampled By</u> | <u>Sample #</u> | <u>Analysis Requested</u> |
|----------------------------|--|-------------------|-------------------|---------------------------|-------------------|-----------------|---------------------------|
| 5/17/2013 1045 | Soil Water Air Protection Enterprise (SWAPE) Bridgeton Sanitary Landfill Air Quality Assessment | U-1 W8 Canister | Summa Canister | 5/14/2013 | Client | 63048 | TO15 ASTM D5504 |
| 5/17/2013 1045 | Soil Water Air Protection Enterprise (SWAPE) Bridgeton Sanitary Landfill Air Quality Assessment | U-2 W6 Canister | Summa Canister | 5/14/2013 | Client | 63049 | TO15 ASTM D5504 |
| 5/17/2013 1045 | Soil Water Air Protection Enterprise (SWAPE) Bridgeton Sanitary Landfill Air Quality Assessment | D-1 W4 Canister | Summa Canister | 5/14/2013 | Client | 63050 | TO15 ASTM D5504 |
| 5/17/2013 1045 | Soil Water Air Protection Enterprise (SWAPE) Bridgeton Sanitary Landfill Air Quality Assessment | D-2 K Canister | Summa Canister | 5/14/2013 | Client | 63051 | TO15 ASTM D5504 |

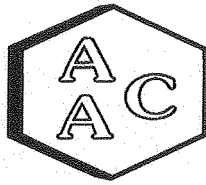
TURN AROUND TIME: Normal (10days)

Lab Due Date: 5/24/2013

Total Samples: 4

REMARKS:

Client returned 4 x Summa canisters + 4 x Flows.



CANISTER PRESSURE LOG

Client: Soil Water Air Protection Ent Project No.: 130597
Date: 5/17/2013

| Canister # | Sample # | Initial Pressure | Final Pressure |
|------------|----------|------------------|----------------|
| 577 | 63048 | 507.6 | 1019.7 |
| 578 | 63049 | 316.6 | 1020.5 |
| 703 | 63050 | 614.5 | 1017.8 |
| 700 | 63051 | 174.8 | 1028.3 |

PH# 30597

CHAIN OF CUSTODY RECORD / ANALYTICAL REQUEST FORM

Bridgeton Sanitary Landfill Air Quality Assessment

Client Name: SOIL / WATER AIR PROTECTION ENTERPRISE
 Project Manager: PAUL ROSENFELD, PH.D.
 Address: 1640 FIFTH STREET, SUITE 204, SANTA MONICA, CA 90401
 Project Name and Location: BRIDGETON SANITARY LANDFILL AIR QUALITY ASSESSMENT
 Telephone No. / Fax No.: (310) 434-0110 / (310) 434-0011
 Date: May 14 2013
 Page 1 of 1

REQUESTED TESTS / ANALYSES

| LAB ID | SAMPLE ID NUMBER | Type | Date | Time | VOCS - EPA TO-15 | Reduced Sulfur Compounds - ASTM D5504 | Carbonyls - EPA TO-11A | Carboxylic Acids - Tube GC-MS | HCL - NIOSH 7903 | Ammonia - OSHA ID-188 | SO2 - OSHA ID-200 | HCN - NIOSH 6010 | Amines - NIOSH 2010M | Fixed Gases - EPA 3C | PAHs / Dioxins EPA TO-13A / 9A | Mercury - NIOSH 6009 | Odor Evaluation | |
|--------|------------------|----------|---------|------|------------------|---------------------------------------|------------------------|-------------------------------|------------------|-----------------------|-------------------|------------------|----------------------|----------------------|--------------------------------|----------------------|-----------------|-----------------|
| 103048 | U-1 W8 | Canister | 5/14/13 | 4 Hr | X | X | | | | | | | | | | | | Canister # 577 |
| 103049 | U-2 W6 | Canister | 5/14/13 | 4 Hr | X | X | | | | | | | | | | | | Canister # 578 |
| 103050 | D-1 W4 | Canister | 5/14/13 | 4 Hr | X | X | | | | | | | | | | | | Canister # 703 |
| 103051 | D-2 K | Canister | 5/14/13 | 4 Hr | X | X | | | | | | | | | | | | Canister # 2155 |
| | | | | | | | | | | | | | | | | | | |
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Special Instructions / Conditions of Receipt

Requested Turnaround Time: Standard turn-around for all analyses. If possible deliver report within 2 weeks.
 OC Requirements: Provide Level IV QC Package for all Analyses.

| | | | | |
|------------------------------------|-------|--------------|-------|---------|
| Relinquished By: John Blank | Date: | Received By: | Date: | Time: |
| | | | | 12 Noon |
| Relinquished By: | Date: | Received By: | Date: | Time: |
| Relinquished By: | Date: | Received By: | Date: | Time: |

- Fed Ex

Atmospheric Analysis and Consulting Inc.

Canister Sampling Field Data Sheet

GENERAL INFORMATION

Project Name and/or ID No.: **Bridgeton Sanitary Landfill**

Site Address and/or ID No: **13570 St Charles Rock Rd, Bridgeton, MO 63044**

Sample Name and/or ID No.: ~~U1-WB~~ **U1-WB/577 Canister**

AAC Batch ID: 130597 AAC Sample ID: 63048

SAMPLING INFORMATION

Start Date/Time: May 14 / 10:53 AM

Stop Date/Time: May 14 / 14:53

Start Temp/Pressure*: 27°C 14.7 psi

Stop Temp/Pressure*: 33°C 14.67psi

Initial Can Pressure**: - 29

Final Can Pressure**: - 9

* Ambient Barometric Reading where sample is being taken (C / inHg) ** Flow Controller Gauge Reading (inHg)

Comments: _____

John Blank

May 14, 2013

Sampler Name (Print)

Sampler Signature/Date

Jeff Miller

Jeff Miller 5/14/13

LABORATORY INFORMATION

Canister Size: 6 - Liter

Sampling Period: 4 - Hour

Canister Serial No.: 577

Flow Controller Serial No: 694

Initial Pressure: 1.4

Certified Flow Rate: 18.0

Return Pressure: 507.6

Certified By/Date: JZ 4/10/2013

Final Pressure: 1019.7

Flow Rate upon Return: 19.3

Date Shipped From Lab: 4/25/2013

Shipped By: JZ

Date Returned to Lab: 5/17/2013

Received By: JZ

Flow Controller Certification File ID: MS03/03281321

Canister Certification File ID: MS03/04051315

Certification Type: SIM _____ SCAN NJLL _____ PAMS _____ Other _____

Jeff Miller 5/20/13

Chemist Signature/Date

N/A 5/20/13

Lab Manager Signature/Date

**Sampler is required to fill out all highlighted sections during sampling.
All remaining sections will be completed upon return by the laboratory.**

Atmospheric Analysis and Consulting Inc.

Canister Sampling Field Data Sheet

GENERAL INFORMATION

Project Name and/or ID No.: **Bridgeton Sanitary Landfill**

Site Address and/or ID No: **13570 St Charles Rock Rd, Bridgeton, MO 63044**

Sample Name and/or ID No.: **U2-W/578 Canister**

AAC Batch ID: 130597 AAC Sample ID: 63049

SAMPLING INFORMATION

Start Date/Time: May 14 / 11:09

Stop Date/Time: May 14 / 15:09

Start Temp/Pressure*: 27°C 14.7 psi

Stop Temp/Pressure*: 33°C 14.67 psi

Initial Can Pressure**: - 30

Final Can Pressure**: - 17

* Ambient Barometric Reading where sample is being taken (C/inHg) ** Flow Controller Gauge Reading (inHg)

Comments: _____

John Blank 2013

Samplers Name (Print)

Jeff Miller

LABORATORY INFORMATION

John Blank

May 14, 2013

Samplers Signature/Date

Jeff Miller

5/14/13

Canister Size: 6 - Liter

Sampling Period: 4 - Hour

Canister Serial No.: 578

Flow Controller Serial No.: 710

Initial Pressure: 1.3

Certified Flow Rate: 18.1

Return Pressure: 3/4.6

Certified By/Date: J2 4/5/2013

Final Pressure: 1020.5

Flow Rate upon Return: 13.6

Date Shipped From Lab: 4/25/2013

Shipped By: J2

Date Returned to Lab: 5/17/2013

Received By: J2

Flow Controller Certification File ID: 11603/03281321

Canister Certification File ID: 11603/04021314

Certification Type: SIM SCAN NJLL PAMS Other

Quana Reed

Chemist Signature/Date

John Blank

Lab Manager Signature/Date

**Sampler is required to fill out all highlighted sections during sampling.
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Atmospheric Analysis and Consulting Inc.

Canister Sampling Field Data Sheet

GENERAL INFORMATION

Project Name and/or ID No.: **Bridgeton Sanitary Landfill**

Site Address and/or ID No.: **13570 St Charles Rock Rd, Bridgeton, MO 63044**

Sample Name and/or ID No.: **W/D 1-W4/703 Canister**

AAC Batch ID: 130597 AAC Sample ID: 63050

SAMPLING INFORMATION

Start Date/Time: May 14 / 10:13

Stop Date/Time: May 14 / 14:13

Start Temp/Pressure*: 26 C 14.7 psi

Stop Temp/Pressure*: 33 C 14.67 psi

Initial Can Pressure**: - 29

Final Can Pressure**: - 4

* Ambient Barometric Reading where sample is being taken (C / inHg) ** Flow Controller Gauge Reading (inHg)

Comments: _____

John Blank

Sampler Name (Print)

Jeff Miller

LABORATORY INFORMATION

Phill Blaul

May 14, 2013

Sampler Signature/Date

Jeff Miller 15/14/13

Canister Size: 6 - Liter

Sampling Period: 4 - Hour

Canister Serial No.: 703

Flow Controller Serial No.: 715

Initial Pressure: 1.5

Certified Flow Rate: 18.0

Return Pressure: 614.5

Certified By/Date: J2 4/26/2013

Final Pressure: 1017.8

Flow Rate upon Return: 21.8

Date Shipped From Lab: 4/25/2013

Shipped By: J2

Date Returned to Lab: 5/17/2013

Received By: J2

Flow Controller Certification File ID: 11503/04261305

Canister Certification File ID: 11502/0411321

Certification Type: SIM SCAN NJLL PAMS Other

James R. Osted
Chemist Signature/Date

MM 5/20/13
Lab Manager Signature/Date

Sampler is required to fill out all highlighted sections during sampling.
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Atmospheric Analysis and Consulting Inc.

Canister Sampling Field Data Sheet

GENERAL INFORMATION

Project Name and/or ID No.: **Bridgeton Sanitary Landfill**

Site Address and/or ID No: **13570 St Charles Rock Rd, Bridgeton, MO 63044**

Sample Name and/or ID No.: **WD 2-K / 2155 Canister**

AAC Batch ID: 130597 AAC Sample ID: 700 CG 05/17/13 63051

SAMPLING INFORMATION

Start Date/Time: May 14 / 10:36

Stop Date/Time: May 14 / 14:36

Start Temp/Pressure*: 26 C 14.7 psi

Stop Temp/Pressure*: 33 C 14.67 psi

Initial Can Pressure**: - 30

Final Can Pressure**: - 22.5

* Ambient Barometric Reading where sample is being taken (C / inHg) ** Flow Controller Gauge Reading (inHg)

Comments: _____

John Blank
Sampler Name (Print)

Jeff Miller

John Blank

May 14, 2013

Sampler Signature/Date

Jeff Miller 5/14/13

LABORATORY INFORMATION

Canister Size: 6 - Liter

Sampling Period: 4 - Hour

Canister Serial No.: 2155 700

Flow Controller Serial No: 717

Initial Pressure: 2.3 CG 05/17/13

Certified Flow Rate: 18.0

Return Pressure: 174.8

Certified By/Date: JZ 4/5/2013

Final Pressure: 1028.3

Flow Rate upon Return: 11.0

Date Shipped From Lab: 4/25/2013

Shipped By: JZ

Date Returned to Lab: 5/17/2013

Received By: JZ

Flow Controller Certification File ID: MS03/03260320

Canister Certification File ID: MS03/03071316

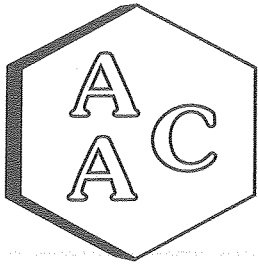
Certification Type: SIM _____ SCAN NJLL _____ PAMS _____ Other _____

James Ben Postol
Chemist Signature/Date

MM 5/20/13
Lab Manager Signature/Date

Sampler is required to fill out all highlighted sections during sampling.
All remaining sections will be completed upon return by the laboratory.

TO-15 REPORTS



Atmospheric Analysis & Consulting, Inc.

Laboratory Analysis Report

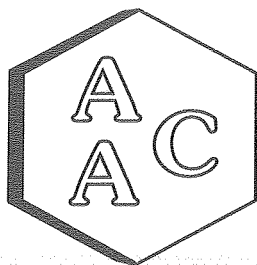
CLIENT : Soil Water Air Protection Enterprise
PROJECT NO : 130597
MATRIX : AIR
UNITS : PPB (v/v)

DATE RECEIVED : 05/17/2013
DATE REPORTED : 05/21/2013

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

| Client ID AAC ID | U-1 W8-Canister 130597-63048 | | | Sample Reporting Limit (SRL) (MRLxDF's) | U-2 W6-Canister 130597-63049 | | | Sample Reporting Limit (SRL) (MRLxDF's) | Method Reporting Limit (MRL) |
|--------------------------------|---------------------------------|---------------|---------------------|---|---------------------------------|---------------|---------------------|---|---------------------------------|
| | Date Sampled | Date Analyzed | Can Dilution Factor | | Date Sampled | Date Analyzed | Can Dilution Factor | | |
| | Result | Qualifier | Analysis DF | | Result | Qualifier | Analysis DF | | |
| Chlorodifluoromethane | 0.32 | J | 1.0 | 1.00 | 0.35 | J | 1.0 | 1.61 | 0.5 |
| Propene | <SRL | U | 1.0 | 2.01 | <SRL | U | 1.0 | 3.22 | 1.0 |
| Dichlorodifluoromethane | 0.58 | J | 1.0 | 1.00 | 0.68 | J | 1.0 | 1.61 | 0.5 |
| Chloromethane | 0.42 | J | 1.0 | 1.00 | 0.58 | J | 1.0 | 1.61 | 0.5 |
| Dichlorotetrafluoroethane | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Vinyl Chloride | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Methanol | 17.2 | | 1.0 | 10.0 | 19.1 | | 1.0 | 16.1 | 5.0 |
| 1,3-Butadiene | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Bromomethane | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Chloroethane | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Dichlorofluoromethane | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Ethanol | 5.54 | | 1.0 | 4.02 | 8.06 | | 1.0 | 6.45 | 2.0 |
| Vinyl Bromide | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Acetone | 8.24 | | 1.0 | 4.02 | 8.25 | | 1.0 | 6.45 | 2.0 |
| Trichlorofluoromethane | 0.28 | J | 1.0 | 1.00 | 0.26 | J | 1.0 | 1.61 | 0.5 |
| 2-Propanol (IPA) | 0.48 | J | 1.0 | 4.02 | 1.74 | J | 1.0 | 6.45 | 2.0 |
| Acrylonitrile | <SRL | U | 1.0 | 2.01 | <SRL | U | 1.0 | 3.22 | 1.0 |
| 1,1-Dichloroethene | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Methylene Chloride (DCM) | <SRL | U | 1.0 | 2.01 | <SRL | U | 1.0 | 3.22 | 1.0 |
| Allyl Chloride | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Carbon Disulfide | NR | U | 1.0 | 1.00 | NR | U | 1.0 | 1.61 | 0.5 |
| Trichlorotrifluoroethane | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| trans-1,2-Dichloroethene | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| 1,1-Dichloroethane | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Methyl Tert Butyl Ether (MTBE) | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Vinyl Acetate | <SRL | U | 1.0 | 2.01 | <SRL | U | 1.0 | 3.22 | 1.0 |
| 2-Butanone (MEK) | 0.98 | J | 1.0 | 2.01 | 0.90 | J | 1.0 | 3.22 | 1.0 |
| cis-1,2-Dichloroethene | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Hexane | 1.02 | | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Chloroform | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Ethyl Acetate | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Tetrahydrofuran | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| 1,2-Dichloroethane | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| 1,1,1-Trichloroethane | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |





Atmospheric Analysis & Consulting, Inc.

Laboratory Analysis Report

CLIENT : Soil Water Air Protection Enterprise
PROJECT NO : 130597
MATRIX : AIR
UNITS : PPB (v/v)

DATE RECEIVED : 05/17/2013
DATE REPORTED : 05/21/2013

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

| Client ID AAC ID | U-1 W8-Canister | | | Sample Reporting Limit (SRL) (MRLxDF's) | U-2 W6-Canister | | | Sample Reporting Limit (SRL) (MRLxDF's) | Method Reporting Limit (MRL) |
|-----------------------------------|-----------------|-----------|-------------|--|-----------------|-----------|-------------|--|------------------------------|
| | 130597-63048 | | | | 130597-63049 | | | | |
| Date Sampled | 05/14/2013 | | | | 05/14/2013 | | | | |
| Date Analyzed | 05/20/2013 | | | | 05/20/2013 | | | | |
| Can Dilution Factor | 2.01 | | | | 3.22 | | | | |
| | Result | Qualifier | Analysis DF | | Result | Qualifier | Analysis DF | | |
| Benzene | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Carbon Tetrachloride | 0.10 | J | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Cyclohexane | 0.14 | J | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| 1,2-Dichloropropane | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Bromodichloromethane | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| 1,4-Dioxane | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Trichloroethene (TCE) | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| 2,2,4-Trimethylpentane | 0.76 | J | 1.0 | 1.00 | 0.81 | J | 1.0 | 1.61 | 0.5 |
| Heptane | 0.44 | J | 1.0 | 1.00 | 0.42 | J | 1.0 | 1.61 | 0.5 |
| cis-1,3-Dichloropropene | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| 4-Methyl-2-pentanone (MIBK) | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| trans-1,3-Dichloropropene | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| 1,1,2-Trichloroethane | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Toluene | 1.17 | J | 1.0 | 1.00 | 1.74 | J | 1.0 | 1.61 | 0.5 |
| 2-Hexanone (MBK) | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Dibromochloromethane | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| 1,2-Dibromoethane | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Tetrachloroethene (PCE) | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Chlorobenzene | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Ethylbenzene | 0.30 | J | 1.0 | 1.00 | 0.52 | J | 1.0 | 1.61 | 0.5 |
| m & p-Xylenes | 1.02 | J | 1.0 | 2.01 | 1.68 | J | 1.0 | 3.22 | 1.0 |
| Bromoform | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Styrene | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| 1,1,2,2-Tetrachloroethane | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| o-Xylene | 0.44 | J | 1.0 | 1.00 | 0.74 | J | 1.0 | 1.61 | 0.5 |
| 4-Ethyltoluene | 0.20 | J | 1.0 | 1.00 | 0.29 | J | 1.0 | 1.61 | 0.5 |
| 1,3,5-Trimethylbenzene | 0.20 | J | 1.0 | 1.00 | 0.32 | J | 1.0 | 1.61 | 0.5 |
| 1,2,4-Trimethylbenzene | 0.70 | J | 1.0 | 1.00 | 1.13 | J | 1.0 | 1.61 | 0.5 |
| Benzyl Chloride (a-Chlorotoluene) | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| 1,3-Dichlorobenzene | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| 1,4-Dichlorobenzene | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| 1,2-Dichlorobenzene | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| 1,2,4-Trichlorobenzene | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| Hexachlorobutadiene | <SRL | U | 1.0 | 1.00 | <SRL | U | 1.0 | 1.61 | 0.5 |
| BFB-Surrogate Std. % Recovery | 102% | | | | 101% | | | | 70-130% |

U - Compound was analyzed for, but was not detected at or above the SRL.

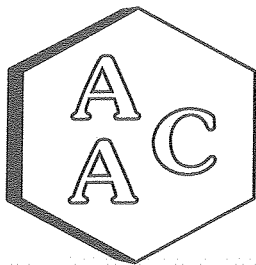
J - Analyte was detected. However the analyte concentration is an estimated value, which is between the Sample Reporting Limit (SRL) and the Sample Quantitation Limit (SQL).

SQL - Sample Quantitation Limit is the Limit Of Quantitation (LOQ) x Dilution Factors.

NR - Not Reported on these analysis.


 Marcus Hueppe
 Laboratory Director





Atmospheric Analysis & Consulting, Inc.

Laboratory Analysis Report

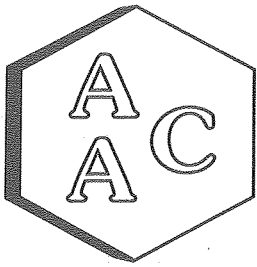
CLIENT : Soil Water Air Protection Enterprise
PROJECT NO : 130597
MATRIX : AIR
UNITS : ug/m3

DATE RECEIVED : 05/17/2013
DATE REPORTED : 05/21/2013

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

| Client ID | U-1 W8-Canister | | | Sample Reporting Limit (SRL) (MRLxDF's) | U-2 W6-Canister | | | Sample Reporting Limit (SRL) (MRLxDF's) | Method Reporting Limit (MRL) |
|--------------------------------|-----------------|--------|-----------|---|-----------------|--------|-----------|---|------------------------------|
| | AAC ID | Result | Qualifier | | Analysis DF | Result | Qualifier | | |
| | 130597-63048 | | | | | | | | |
| Date Sampled | 05/14/2013 | | | | 05/14/2013 | | | | |
| Date Analyzed | 05/20/2013 | | | | 05/20/2013 | | | | |
| Can Dilution Factor | 2.01 | | | | 3.22 | | | | |
| Chlorodifluoromethane | 1.1 | J | 1.0 | 3.6 | 1.3 | J | 1.0 | 5.7 | 1.8 |
| Propene | <SRL | U | 1.0 | 3.5 | <SRL | U | 1.0 | 5.5 | 1.7 |
| Dichlorodifluoromethane | 2.9 | J | 1.0 | 5.0 | 3.4 | J | 1.0 | 8.0 | 2.5 |
| Chloromethane | 0.9 | J | 1.0 | 2.1 | 1.2 | J | 1.0 | 3.3 | 1.0 |
| Dichlorotetrafluoroethane | <SRL | U | 1.0 | 7.0 | <SRL | U | 1.0 | 11.3 | 3.5 |
| Vinyl Chloride | <SRL | U | 1.0 | 2.6 | <SRL | U | 1.0 | 4.1 | 1.3 |
| Methanol | 22.5 | | 1.0 | 13.2 | 25.0 | | 1.0 | 21.1 | 6.6 |
| 1,3-Butadiene | <SRL | U | 1.0 | 2.2 | <SRL | U | 1.0 | 3.6 | 1.1 |
| Bromomethane | <SRL | U | 1.0 | 3.9 | <SRL | U | 1.0 | 6.3 | 1.9 |
| Chloroethane | <SRL | U | 1.0 | 2.6 | <SRL | U | 1.0 | 4.3 | 1.3 |
| Dichlorofluoromethane | <SRL | U | 1.0 | 4.2 | <SRL | U | 1.0 | 6.8 | 2.1 |
| Ethanol | 10.4 | | 1.0 | 7.6 | 15.2 | | 1.0 | 12.1 | 3.8 |
| Vinyl Bromide | <SRL | U | 1.0 | 4.4 | <SRL | U | 1.0 | 7.1 | 2.2 |
| Acetone | 19.6 | | 1.0 | 9.5 | 19.6 | | 1.0 | 15.3 | 4.8 |
| Trichlorofluoromethane | 1.6 | J | 1.0 | 5.6 | 1.5 | J | 1.0 | 9.1 | 2.8 |
| 2-Propanol (IPA) | 1.2 | J | 1.0 | 9.9 | 4.3 | J | 1.0 | 15.8 | 4.9 |
| Acrylonitrile | <SRL | U | 1.0 | 4.4 | <SRL | U | 1.0 | 7.0 | 2.2 |
| 1,1-Dichloroethene | <SRL | U | 1.0 | 4.0 | <SRL | U | 1.0 | 6.4 | 2.0 |
| Methylene Chloride (DCM) | <SRL | U | 1.0 | 7.0 | <SRL | U | 1.0 | 11.2 | 3.5 |
| Allyl Chloride | <SRL | U | 1.0 | 3.1 | <SRL | U | 1.0 | 5.0 | 1.6 |
| Carbon Disulfide | NR | U | 1.0 | 3.1 | NR | U | 1.0 | 5.0 | 1.6 |
| Trichlorotrifluoroethane | <SRL | U | 1.0 | 7.7 | <SRL | U | 1.0 | 12.4 | 3.8 |
| trans-1,2-Dichloroethene | <SRL | U | 1.0 | 4.0 | <SRL | U | 1.0 | 6.4 | 2.0 |
| 1,1-Dichloroethane | <SRL | U | 1.0 | 4.1 | <SRL | U | 1.0 | 6.5 | 2.0 |
| Methyl Tert Butyl Ether (MTBE) | <SRL | U | 1.0 | 3.6 | <SRL | U | 1.0 | 5.8 | 1.8 |
| Vinyl Acetate | <SRL | U | 1.0 | 7.1 | <SRL | U | 1.0 | 11.3 | 3.5 |
| 2-Butanone (MEK) | 2.9 | J | 1.0 | 5.9 | 2.7 | J | 1.0 | 9.5 | 2.9 |
| cis-1,2-Dichloroethene | <SRL | U | 1.0 | 4.0 | <SRL | U | 1.0 | 6.4 | 2.0 |
| Hexane | 3.6 | | 1.0 | 3.5 | <SRL | U | 1.0 | 5.7 | 1.8 |
| Chloroform | <SRL | U | 1.0 | 4.9 | <SRL | U | 1.0 | 7.9 | 2.4 |
| Ethyl Acetate | <SRL | U | 1.0 | 3.6 | <SRL | U | 1.0 | 5.8 | 1.8 |
| Tetrahydrofuran | <SRL | U | 1.0 | 3.0 | <SRL | U | 1.0 | 4.8 | 1.5 |
| 1,2-Dichloroethane | <SRL | U | 1.0 | 4.1 | <SRL | U | 1.0 | 6.5 | 2.0 |
| 1,1,1-Trichloroethane | <SRL | U | 1.0 | 5.5 | <SRL | U | 1.0 | 8.8 | 2.7 |





Atmospheric Analysis & Consulting, Inc.

Laboratory Analysis Report

CLIENT : Soil Water Air Protection Enterprise
PROJECT NO : 130597
MATRIX : AIR
UNITS : ug/m3

DATE RECEIVED : 05/17/2013
DATE REPORTED : 05/21/2013

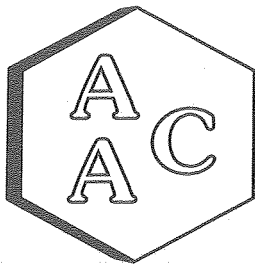
VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

| Client ID | U-1 W8-Canister | | | Sample Reporting Limit (SRL) (MRLxDF's) | U-2 W6-Canister | | | Sample Reporting Limit (SRL) (MRLxDF's) | Method Reporting Limit (MRL) |
|-----------------------------------|-----------------|--------|-----------|---|-----------------|--------|--------|---|------------------------------|
| | AAC ID | Result | Qualifier | | Analysis DF | AAC ID | Result | | |
| Date Sampled | 130597-63048 | | | | 130597-63049 | | | | |
| Date Analyzed | 05/14/2013 | | | | 05/14/2013 | | | | |
| Can Dilution Factor | 05/20/2013 | | | | 05/20/2013 | | | | |
| | 2.01 | | | | 3.22 | | | | |
| Benzene | <SRL | U | 1.0 | 3.2 | <SRL | U | 1.0 | 5.1 | 1.6 |
| Carbon Tetrachloride | 0.6 | J | 1.0 | 6.3 | <SRL | U | 1.0 | 10.1 | 3.1 |
| Cyclohexane | 0.5 | J | 1.0 | 3.5 | <SRL | U | 1.0 | 5.5 | 1.7 |
| 1,2-Dichloropropane | <SRL | U | 1.0 | 4.6 | <SRL | U | 1.0 | 7.4 | 2.3 |
| Bromodichloromethane | <SRL | U | 1.0 | 6.7 | <SRL | U | 1.0 | 10.8 | 3.4 |
| 1,4-Dioxane | <SRL | U | 1.0 | 3.6 | <SRL | U | 1.0 | 5.8 | 1.8 |
| Trichloroethene (TCE) | <SRL | U | 1.0 | 5.4 | <SRL | U | 1.0 | 8.7 | 2.7 |
| 2,2,4-Trimethylpentane | 3.6 | J | 1.0 | 4.7 | 3.8 | J | 1.0 | 7.5 | 2.3 |
| Heptane | 1.8 | J | 1.0 | 4.1 | 1.7 | J | 1.0 | 6.6 | 2.0 |
| cis-1,3-Dichloropropene | <SRL | U | 1.0 | 4.6 | <SRL | U | 1.0 | 7.3 | 2.3 |
| 4-Methyl-2-pentanone (MiBK) | <SRL | U | 1.0 | 4.1 | <SRL | U | 1.0 | 6.6 | 2.0 |
| trans-1,3-Dichloropropene | <SRL | U | 1.0 | 4.6 | <SRL | U | 1.0 | 7.3 | 2.3 |
| 1,1,2-Trichloroethane | <SRL | U | 1.0 | 5.5 | <SRL | U | 1.0 | 8.8 | 2.7 |
| Toluene | 4.4 | J | 1.0 | 3.8 | 6.6 | J | 1.0 | 6.1 | 1.9 |
| 2-Hexanone (MBK) | <SRL | U | 1.0 | 4.1 | <SRL | U | 1.0 | 6.6 | 2.0 |
| Dibromochloromethane | <SRL | U | 1.0 | 8.6 | <SRL | U | 1.0 | 13.7 | 4.3 |
| 1,2-Dibromoethane | <SRL | U | 1.0 | 7.7 | <SRL | U | 1.0 | 12.4 | 3.8 |
| Tetrachloroethene (PCE) | <SRL | U | 1.0 | 6.8 | <SRL | U | 1.0 | 10.9 | 3.4 |
| Chlorobenzene | <SRL | U | 1.0 | 4.6 | <SRL | U | 1.0 | 7.4 | 2.3 |
| Ethylbenzene | 1.3 | J | 1.0 | 4.4 | 2.2 | J | 1.0 | 7.0 | 2.2 |
| m & p-Xylenes | 4.5 | J | 1.0 | 8.7 | 7.3 | J | 1.0 | 14.0 | 4.3 |
| Bromoform | <SRL | U | 1.0 | 10.4 | <SRL | U | 1.0 | 16.7 | 5.2 |
| Styrene | <SRL | U | 1.0 | 4.3 | <SRL | U | 1.0 | 6.9 | 2.1 |
| 1,1,2,2-Tetrachloroethane | <SRL | U | 1.0 | 6.9 | <SRL | U | 1.0 | 11.1 | 3.4 |
| o-Xylene | 1.9 | J | 1.0 | 4.4 | 3.2 | J | 1.0 | 7.0 | 2.2 |
| 4-Ethyltoluene | 1.0 | J | 1.0 | 4.9 | 1.4 | J | 1.0 | 7.9 | 2.5 |
| 1,3,5-Trimethylbenzene | 1.0 | J | 1.0 | 4.9 | 1.6 | J | 1.0 | 7.9 | 2.5 |
| 1,2,4-Trimethylbenzene | 3.5 | J | 1.0 | 4.9 | 5.6 | J | 1.0 | 7.9 | 2.5 |
| Benzyl Chloride (a-Chlorotoluene) | <SRL | U | 1.0 | 5.2 | <SRL | U | 1.0 | 8.3 | 2.6 |
| 1,3-Dichlorobenzene | <SRL | U | 1.0 | 6.0 | <SRL | U | 1.0 | 9.7 | 3.0 |
| 1,4-Dichlorobenzene | <SRL | U | 1.0 | 6.0 | <SRL | U | 1.0 | 9.7 | 3.0 |
| 1,2-Dichlorobenzene | <SRL | U | 1.0 | 6.0 | <SRL | U | 1.0 | 9.7 | 3.0 |
| 1,2,4-Trichlorobenzene | <SRL | U | 1.0 | 7.5 | <SRL | U | 1.0 | 12.0 | 3.7 |
| Hexachlorobutadiene | <SRL | U | 1.0 | 10.7 | <SRL | U | 1.0 | 17.2 | 5.3 |
| BFB-Surrogate Std. % Recovery | 102% | | | | 101% | | | | 70-130% |

U - Compound was analyzed for, but was not detected at or above the SRL.
 J - Analyte was detected. However the analyte concentration is an estimated value, which is between the Sample Reporting Limit (SRL) and the Sample Quantitation Limit (SQL).
 SQL - Sample Quantitation Limit is the Limit Of Quantitation (LOQ) x Dilution Factors.
 NR - Not Reported on these analysis.


 Marcus Hueppe
 Laboratory Director





Atmospheric Analysis & Consulting, Inc.

Laboratory Analysis Report

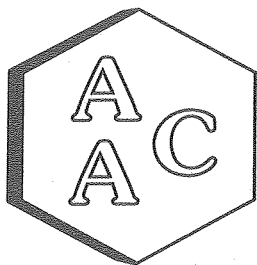
CLIENT : Soil Water Air Protection Enterprise
PROJECT NO : 130597
MATRIX : AIR
UNITS : PPB (v/v)

DATE RECEIVED : 05/17/2013
DATE REPORTED : 05/21/2013

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

| Client ID | D-1 W4-Canister | | | Sample Reporting Limit (SRL) (MRLxDF's) | D-2 K-Canister | | | Sample Reporting Limit (SRL) (MRLxDF's) | Method Reporting Limit (MRL) |
|--------------------------------|-----------------|--------|-----------|---|----------------|--------------|--------|---|------------------------------|
| | AAC ID | Result | Qualifier | | Analysis DF | 130597-63051 | Result | | |
| Date Sampled | 05/14/2013 | | | | 05/14/2013 | | | | |
| Date Analyzed | 05/20/2013 | | | | 05/20/2013 | | | | |
| Can Dilution Factor | 1.66 | | | | 5.88 | | | | |
| Chlorodifluoromethane | 0.31 | J | 1.0 | 0.83 | 0.35 | J | 1.0 | 2.94 | 0.5 |
| Propene | <SRL | U | 1.0 | 1.66 | 5.06 | J | 1.0 | 5.88 | 1.0 |
| Dichlorodifluoromethane | 0.60 | J | 1.0 | 0.83 | 0.71 | J | 1.0 | 2.94 | 0.5 |
| Chloromethane | 0.46 | J | 1.0 | 0.83 | 0.71 | J | 1.0 | 2.94 | 0.5 |
| Dichlorotetrafluoroethane | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Vinyl Chloride | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Methanol | 37.3 | | 1.0 | 8.28 | 62.3 | | 1.0 | 29.4 | 5.0 |
| 1,3-Butadiene | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Bromomethane | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Chloroethane | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Dichlorofluoromethane | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Ethanol | 6.26 | | 1.0 | 3.31 | 32.5 | | 1.0 | 11.8 | 2.0 |
| Vinyl Bromide | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Acetone | 8.55 | | 1.0 | 3.31 | 24.9 | | 1.0 | 11.8 | 2.0 |
| Trichlorofluoromethane | 0.27 | J | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| 2-Propanol (IPA) | 0.80 | J | 1.0 | 3.31 | 31.2 | | 1.0 | 11.8 | 2.0 |
| Acrylonitrile | <SRL | U | 1.0 | 1.66 | <SRL | U | 1.0 | 5.88 | 1.0 |
| 1,1-Dichloroethene | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Methylene Chloride (DCM) | <SRL | U | 1.0 | 1.66 | <SRL | U | 1.0 | 5.88 | 1.0 |
| Allyl Chloride | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Carbon Disulfide | NR | U | 1.0 | 0.83 | NR | U | 1.0 | 2.94 | 0.5 |
| Trichlorotrifluoroethane | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| trans-1,2-Dichloroethene | <SRL | U | 1.0 | 0.83 | 0.71 | J | 1.0 | 2.94 | 0.5 |
| 1,1-Dichloroethane | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Methyl Tert Butyl Ether (MTBE) | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Vinyl Acetate | <SRL | U | 1.0 | 1.66 | <SRL | U | 1.0 | 5.88 | 1.0 |
| 2-Butanone (MEK) | 1.86 | | 1.0 | 1.66 | 2.65 | J | 1.0 | 5.88 | 1.0 |
| cis-1,2-Dichloroethene | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Hexane | 0.91 | | 1.0 | 0.83 | 4.35 | | 1.0 | 2.94 | 0.5 |
| Chloroform | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Ethyl Acetate | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Tetrahydrofuran | 0.94 | | 1.0 | 0.83 | 1.06 | J | 1.0 | 2.94 | 0.5 |
| 1,2-Dichloroethane | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| 1,1,1-Trichloroethane | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |





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Laboratory Analysis Report

CLIENT : Soil Water Air Protection Enterprise
PROJECT NO : 130597
MATRIX : AIR
UNITS : PPB (v/v)

DATE RECEIVED : 05/17/2013
DATE REPORTED : 05/21/2013

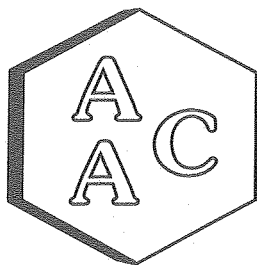
VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

| Client ID AAC ID Date Sampled Date Analyzed Can Dilution Factor | D-1 W4-Canister 130597-63050 | | | Sample Reporting Limit (SRL) (MRLxDF's) | D-2 K-Canister 130597-63051 | | | Sample Reporting Limit (SRL) (MRLxDF's) | Method Reporting Limit (MRL) |
|---|---------------------------------|-----------|-------------|--|--------------------------------|-----------|-------------|--|------------------------------|
| | Result | Qualifier | Analysis DF | | Result | Qualifier | Analysis DF | | |
| | | | 1.66 | | | | 5.88 | | |
| Benzene | 2.34 | | 1.0 | 0.83 | 3.12 | | 1.0 | 2.94 | 0.5 |
| Carbon Tetrachloride | 0.10 | J | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Cyclohexane | 0.13 | J | 1.0 | 0.83 | 0.76 | J | 1.0 | 2.94 | 0.5 |
| 1,2-Dichloropropane | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Bromodichloromethane | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| 1,4-Dioxane | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Trichloroethene (TCE) | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| 2,2,4-Trimethylpentane | 0.65 | J | 1.0 | 0.83 | 3.88 | | 1.0 | 2.94 | 0.5 |
| Heptane | 0.46 | J | 1.0 | 0.83 | 2.47 | J | 1.0 | 2.94 | 0.5 |
| cis-1,3-Dichloropropene | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| 4-Methyl-2-pentanone (MiBK) | 0.10 | J | 1.0 | 0.83 | 0.24 | J | 1.0 | 2.94 | 0.5 |
| trans-1,3-Dichloropropene | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| 1,1,2-Trichloroethane | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Toluene | 1.39 | | 1.0 | 0.83 | 8.53 | | 1.0 | 2.94 | 0.5 |
| 2-Hexanone (MBK) | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Dibromochloromethane | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| 1,2-Dibromoethane | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Tetrachloroethene (PCE) | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Chlorobenzene | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Ethylbenzene | 0.38 | J | 1.0 | 0.83 | 1.53 | J | 1.0 | 2.94 | 0.5 |
| m & p-Xylenes | 1.11 | J | 1.0 | 1.66 | 4.71 | J | 1.0 | 5.88 | 1.0 |
| Bromoform | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Styrene | <SRL | U | 1.0 | 0.83 | 0.88 | J | 1.0 | 2.94 | 0.5 |
| 1,1,2,2-Tetrachloroethane | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| o-Xylene | 0.45 | J | 1.0 | 0.83 | 2.24 | J | 1.0 | 2.94 | 0.5 |
| 4-Ethyltoluene | 0.18 | J | 1.0 | 0.83 | 0.82 | J | 1.0 | 2.94 | 0.5 |
| 1,3,5-Trimethylbenzene | 0.18 | J | 1.0 | 0.83 | 1.00 | J | 1.0 | 2.94 | 0.5 |
| 1,2,4-Trimethylbenzene | 0.65 | J | 1.0 | 0.83 | 3.18 | | 1.0 | 2.94 | 0.5 |
| Benzyl Chloride (a-Chlorotoluene) | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| 1,3-Dichlorobenzene | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| 1,4-Dichlorobenzene | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| 1,2-Dichlorobenzene | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| 1,2,4-Trichlorobenzene | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| Hexachlorobutadiene | <SRL | U | 1.0 | 0.83 | <SRL | U | 1.0 | 2.94 | 0.5 |
| BFB-Surrogate Std. % Recovery | 104% | | | | 103% | | | | 70-130% |

U - Compound was analyzed for, but was not detected at or above the SRL.
 J - Analyte was detected. However the analyte concentration is an estimated value, which is between the Sample Reporting Limit (SRL) and the Sample Quantitation Limit (SQL).
 SQL - Sample Quantitation Limit is the Limit Of Quantitation (LOQ) x Dilution Factors.
 NR - Not Reported on these analysis.


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 Laboratory Director





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Laboratory Analysis Report

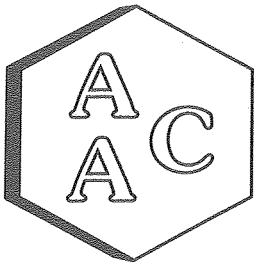
CLIENT : Soil Water Air Protection Enterprise
PROJECT NO : 130597
MATRIX : AIR
UNITS : ug/m3

DATE RECEIVED : 05/17/2013
DATE REPORTED : 05/21/2013

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

| Client ID AAC ID Date Sampled Date Analyzed Can Dilution Factor | D-1 W4-Canister | | | Sample Reporting Limit (SRL) (MRLxDF's) | D-2 K-Canister | | | Sample Reporting Limit (SRL) (MRLxDF's) | Method Reporting Limit (MRL) |
|---|-----------------|-----------|-------------|---|----------------|-------------|-----|---|------------------------------|
| | 130597-63050 | | | | 130597-63051 | | | | |
| | 05/14/2013 | | | | 05/14/2013 | | | | |
| | 05/20/2013 | | | | 05/20/2013 | | | | |
| | 1.66 | | | 5.88 | | | | | |
| | Result | Qualifier | Analysis DF | Result | Qualifier | Analysis DF | | | |
| Chlorodifluoromethane | 1.1 | J | 1.0 | 2.9 | 1.3 | J | 1.0 | 10.4 | 1.8 |
| Propene | <SRL | U | 1.0 | 2.9 | 8.7 | J | 1.0 | 10.1 | 1.7 |
| Dichlorodifluoromethane | 3.0 | J | 1.0 | 4.1 | 3.5 | J | 1.0 | 14.5 | 2.5 |
| Chloromethane | 1.0 | J | 1.0 | 1.7 | 1.5 | J | 1.0 | 6.1 | 1.0 |
| Dichlorotetrafluoroethane | <SRL | U | 1.0 | 5.8 | <SRL | U | 1.0 | 20.6 | 3.5 |
| Vinyl Chloride | <SRL | U | 1.0 | 2.1 | <SRL | U | 1.0 | 7.5 | 1.3 |
| Methanol | 48.9 | | 1.0 | 10.9 | 81.6 | | 1.0 | 38.5 | 6.6 |
| 1,3-Butadiene | <SRL | U | 1.0 | 1.8 | <SRL | U | 1.0 | 6.5 | 1.1 |
| Bromomethane | <SRL | U | 1.0 | 3.2 | <SRL | U | 1.0 | 11.4 | 1.9 |
| Chloroethane | <SRL | U | 1.0 | 2.2 | <SRL | U | 1.0 | 7.8 | 1.3 |
| Dichlorofluoromethane | <SRL | U | 1.0 | 3.5 | <SRL | U | 1.0 | 12.4 | 2.1 |
| Ethanol | 11.8 | | 1.0 | 6.2 | 61.3 | | 1.0 | 22.2 | 3.8 |
| Vinyl Bromide | <SRL | U | 1.0 | 3.6 | <SRL | U | 1.0 | 12.9 | 2.2 |
| Acetone | 20.3 | | 1.0 | 7.9 | 59.1 | | 1.0 | 27.9 | 4.8 |
| Trichlorofluoromethane | 1.5 | J | 1.0 | 4.7 | <SRL | U | 1.0 | 16.5 | 2.8 |
| 2-Propanol (IPA) | 2.0 | J | 1.0 | 8.1 | 76.6 | | 1.0 | 28.9 | 4.9 |
| Acrylonitrile | <SRL | U | 1.0 | 3.6 | <SRL | U | 1.0 | 12.8 | 2.2 |
| 1,1-Dichloroethene | <SRL | U | 1.0 | 3.3 | <SRL | U | 1.0 | 11.7 | 2.0 |
| Methylene Chloride (DCM) | <SRL | U | 1.0 | 5.8 | <SRL | U | 1.0 | 20.4 | 3.5 |
| Allyl Chloride | <SRL | U | 1.0 | 2.6 | <SRL | U | 1.0 | 9.2 | 1.6 |
| Carbon Disulfide | NR | U | 1.0 | 2.6 | NR | U | 1.0 | 9.2 | 1.6 |
| Trichlorotrifluoroethane | <SRL | U | 1.0 | 6.3 | <SRL | U | 1.0 | 22.5 | 3.8 |
| trans-1,2-Dichloroethene | <SRL | U | 1.0 | 3.3 | 2.8 | J | 1.0 | 11.7 | 2.0 |
| 1,1-Dichloroethane | <SRL | U | 1.0 | 3.4 | <SRL | U | 1.0 | 11.9 | 2.0 |
| Methyl Tert Butyl Ether (MTBE) | <SRL | U | 1.0 | 3.0 | <SRL | U | 1.0 | 10.6 | 1.8 |
| Vinyl Acetate | <SRL | U | 1.0 | 5.8 | <SRL | U | 1.0 | 20.7 | 3.5 |
| 2-Butanone (MEK) | 5.5 | | 1.0 | 4.9 | 7.8 | J | 1.0 | 17.3 | 2.9 |
| cis-1,2-Dichloroethene | <SRL | U | 1.0 | 3.3 | <SRL | U | 1.0 | 11.7 | 2.0 |
| Hexane | 3.2 | | 1.0 | 2.9 | 15.3 | | 1.0 | 10.4 | 1.8 |
| Chloroform | <SRL | U | 1.0 | 4.0 | <SRL | U | 1.0 | 14.4 | 2.4 |
| Ethyl Acetate | <SRL | U | 1.0 | 3.0 | <SRL | U | 1.0 | 10.6 | 1.8 |
| Tetrahydrofuran | 2.8 | | 1.0 | 2.4 | 3.1 | J | 1.0 | 8.7 | 1.5 |
| 1,2-Dichloroethane | <SRL | U | 1.0 | 3.4 | <SRL | U | 1.0 | 11.9 | 2.0 |
| 1,1,1-Trichloroethane | <SRL | U | 1.0 | 4.5 | <SRL | U | 1.0 | 16.0 | 2.7 |





Atmospheric Analysis & Consulting, Inc.

Laboratory Analysis Report

CLIENT : Soil Water Air Protection Enterprise
PROJECT NO : 130597
MATRIX : AIR
UNITS : ug/m3

DATE RECEIVED : 05/17/2013
DATE REPORTED : 05/21/2013

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

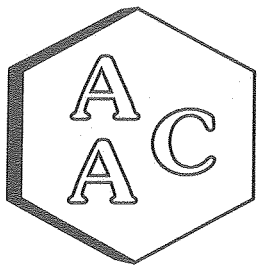
| Client ID | D-1 W4-Canister | | | Sample Reporting Limit (SRL) (MRLxDF's) | D-2 K-Canister | | | Sample Reporting Limit (SRL) (MRLxDF's) | Method Reporting Limit (MRL) |
|-----------------------------------|-----------------|--------|-----------|---|----------------|--------|-----------|---|------------------------------|
| | AAC ID | Result | Qualifier | | Analysis DF | Result | Qualifier | | |
| | 130597-63050 | | | | 130597-63051 | | | | |
| Date Sampled | 05/14/2013 | | | | 05/14/2013 | | | | |
| Date Analyzed | 05/20/2013 | | | | 05/20/2013 | | | | |
| Can Dilution Factor | 1.66 | | | | 5.88 | | | | |
| Benzene | 7.5 | | 1.0 | 2.6 | 10.0 | | 1.0 | 9.4 | 1.6 |
| Carbon Tetrachloride | 0.6 | J | 1.0 | 5.2 | <SRL | U | 1.0 | 18.5 | 3.1 |
| Cyclohexane | 0.5 | J | 1.0 | 2.9 | 2.6 | J | 1.0 | 10.1 | 1.7 |
| 1,2-Dichloropropane | <SRL | U | 1.0 | 3.8 | <SRL | U | 1.0 | 13.6 | 2.3 |
| Bromodichloromethane | <SRL | U | 1.0 | 5.5 | <SRL | U | 1.0 | 19.7 | 3.4 |
| 1,4-Dioxane | <SRL | U | 1.0 | 3.0 | <SRL | U | 1.0 | 10.6 | 1.8 |
| Trichloroethene (TCE) | <SRL | U | 1.0 | 4.5 | <SRL | U | 1.0 | 15.8 | 2.7 |
| 2,2,4-Trimethylpentane | 3.0 | J | 1.0 | 3.9 | 18.1 | | 1.0 | 13.7 | 2.3 |
| Heptane | 1.9 | J | 1.0 | 3.4 | 10.1 | J | 1.0 | 12.1 | 2.0 |
| cis-1,3-Dichloropropene | <SRL | U | 1.0 | 3.8 | <SRL | U | 1.0 | 13.3 | 2.3 |
| 4-Methyl-2-pentanone (MIBK) | 0.4 | J | 1.0 | 3.4 | 1.0 | J | 1.0 | 12.0 | 2.0 |
| trans-1,3-Dichloropropene | <SRL | U | 1.0 | 3.8 | <SRL | U | 1.0 | 13.3 | 2.3 |
| 1,1,2-Trichloroethane | <SRL | U | 1.0 | 4.5 | <SRL | U | 1.0 | 16.0 | 2.7 |
| Toluene | 5.2 | | 1.0 | 3.1 | 32.1 | | 1.0 | 11.1 | 1.9 |
| 2-Hexanone (MBK) | <SRL | U | 1.0 | 3.4 | <SRL | U | 1.0 | 12.0 | 2.0 |
| Dibromochloromethane | <SRL | U | 1.0 | 7.1 | <SRL | U | 1.0 | 25.1 | 4.3 |
| 1,2-Dibromoethane | <SRL | U | 1.0 | 6.4 | <SRL | U | 1.0 | 22.6 | 3.8 |
| Tetrachloroethene (PCE) | <SRL | U | 1.0 | 5.6 | <SRL | U | 1.0 | 19.9 | 3.4 |
| Chlorobenzene | <SRL | U | 1.0 | 3.8 | <SRL | U | 1.0 | 13.5 | 2.3 |
| Ethylbenzene | 1.7 | J | 1.0 | 3.6 | 6.6 | J | 1.0 | 12.8 | 2.2 |
| m & p-Xylenes | 4.8 | J | 1.0 | 7.2 | 20.4 | J | 1.0 | 25.5 | 4.3 |
| Bromoform | <SRL | U | 1.0 | 8.6 | <SRL | U | 1.0 | 30.4 | 5.2 |
| Styrene | <SRL | U | 1.0 | 3.5 | 3.8 | J | 1.0 | 12.5 | 2.1 |
| 1,1,2,2-Tetrachloroethane | <SRL | U | 1.0 | 5.7 | <SRL | U | 1.0 | 20.2 | 3.4 |
| o-Xylene | 1.9 | J | 1.0 | 3.6 | 9.7 | J | 1.0 | 12.8 | 2.2 |
| 4-Ethyltoluene | 0.9 | J | 1.0 | 4.1 | 4.1 | J | 1.0 | 14.5 | 2.5 |
| 1,3,5-Trimethylbenzene | 0.9 | J | 1.0 | 4.1 | 4.9 | J | 1.0 | 14.5 | 2.5 |
| 1,2,4-Trimethylbenzene | 3.2 | J | 1.0 | 4.1 | 15.6 | | 1.0 | 14.5 | 2.5 |
| Benzyl Chloride (a-Chlorotoluene) | <SRL | U | 1.0 | 4.3 | <SRL | U | 1.0 | 15.2 | 2.6 |
| 1,3-Dichlorobenzene | <SRL | U | 1.0 | 5.0 | <SRL | U | 1.0 | 17.7 | 3.0 |
| 1,4-Dichlorobenzene | <SRL | U | 1.0 | 5.0 | <SRL | U | 1.0 | 17.7 | 3.0 |
| 1,2-Dichlorobenzene | <SRL | U | 1.0 | 5.0 | <SRL | U | 1.0 | 17.7 | 3.0 |
| 1,2,4-Trichlorobenzene | <SRL | U | 1.0 | 6.1 | <SRL | U | 1.0 | 21.8 | 3.7 |
| Hexachlorobutadiene | <SRL | U | 1.0 | 8.8 | <SRL | U | 1.0 | 31.4 | 5.3 |
| BFB-Surrogate Std. % Recovery | 104% | | | | 103% | | | | 70-130% |

U - Compound was analyzed for, but was not detected at or above the SRL.
 J - Analyte was detected. However the analyte concentration is an estimated value, which is between the Sample Reporting Limit (SRL) and the Sample Quantitation Limit (SQL).
 SQL - Sample Quantitation Limit is the Limit Of Quantitation (LOQ) x Dilution Factors.
 NR - Not Reported on these analysis.


 Marcus Hueppe
 Laboratory Director



TO-15 QC REPORT



Atmospheric Analysis & Consulting, Inc.

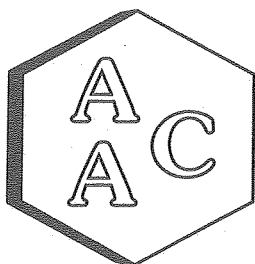
ANALYSIS DATE : 05/20/2013
ANALYST : JJG

INSTRUMENT ID : GC/MS-03
CALIBRATION STD ID : PS040413-01

VOLATILE ORGANIC COMPOUNDS BY EPA METHOD TO-15
Continuing Calibration Verification of the 05/15/2013 Calibration

| <i>Compounds</i> | <i>Conc</i> | <i>Daily Conc</i> | <i>%REC*</i> |
|--------------------------------|-------------|-------------------|--------------|
| 4-BFB (surrogate standard) | 10.00 | 10.18 | 102 |
| Chlorodifluoromethane | 10.10 | 10.26 | 102 |
| Propene | 11.00 | 10.92 | 99 |
| Dichlorodifluoromethane | 9.80 | 10.08 | 103 |
| Chloromethane | 10.10 | 10.06 | 100 |
| Dichlorotetrafluoroethane | 10.10 | 10.13 | 100 |
| Vinyl Chloride | 10.20 | 10.13 | 99 |
| Methanol | 4.90 | 5.27 | 108 |
| 1,3-Butadiene | 10.50 | 10.48 | 100 |
| Bromomethane | 10.20 | 9.15 | 90 |
| Chloroethane | 10.00 | 9.61 | 96 |
| Dichlorofluoromethane | 10.00 | 10.28 | 103 |
| Ethanol | 9.80 | 10.34 | 106 |
| Vinyl Bromide | 10.20 | 10.56 | 104 |
| Acetone | 10.80 | 9.73 | 90 |
| Trichlorofluoromethane | 10.10 | 11.18 | 111 |
| 2-Propanol (IPA) | 11.00 | 10.46 | 95 |
| Acrylonitrile | 10.50 | 11.08 | 106 |
| 1,1-Dichloroethene | 10.50 | 10.27 | 98 |
| Methylene Chloride (DCM) | 10.40 | 9.95 | 96 |
| Allyl Chloride | 11.00 | 11.62 | 106 |
| Carbon Disulfide | 10.50 | 9.94 | 95 |
| Trichlorotrifluoroethane | 10.40 | 10.41 | 100 |
| trans-1,2-Dichloroethene | 10.40 | 10.75 | 103 |
| 1,1-Dichloroethane | 10.40 | 10.47 | 101 |
| Methyl Tert Butyl Ether (MTBE) | 10.60 | 11.19 | 106 |
| Vinyl Acetate | 9.70 | 10.24 | 106 |
| 2-Butanone (MEK) | 10.60 | 11.00 | 104 |
| cis-1,2-Dichloroethene | 10.60 | 10.48 | 99 |
| Hexane | 10.70 | 10.81 | 101 |
| Chloroform | 10.60 | 11.01 | 104 |
| Ethyl Acetate | 11.00 | 11.51 | 105 |
| Tetrahydrofuran | 10.80 | 10.77 | 100 |
| 1,2-Dichloroethane | 10.40 | 11.21 | 108 |
| 1,1,1-Trichloroethane | 10.50 | 11.32 | 108 |





Atmospheric Analysis & Consulting, Inc.

ANALYSIS DATE : 05/20/2013

INSTRUMENT ID : GC/MS-03

ANALYST : JJG


CALIBRATION STD ID : PS040413-01

VOLATILE ORGANIC COMPOUNDS BY EPA METHOD TO-15

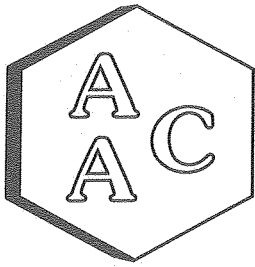
Continuing Calibration Verification of the 05/15/2013 Calibration

| Compounds | Conc | Daily Conc | %REC* |
|-----------------------------------|-------|------------|-------|
| Benzene | 10.50 | 9.62 | 92 |
| Carbon Tetrachloride | 10.10 | 10.51 | 104 |
| Cyclohexane | 10.50 | 9.94 | 95 |
| 1,2-Dichloropropane | 10.50 | 9.92 | 94 |
| Bromodichloromethane | 10.30 | 10.87 | 106 |
| 1,4-Dioxane | 10.30 | 9.91 | 96 |
| Trichloroethene (TCE) | 10.30 | 10.22 | 99 |
| 2,2,4-Trimethylpentane | 10.90 | 10.56 | 97 |
| Heptane | 10.70 | 10.28 | 96 |
| cis-1,3-Dichloropropene | 11.00 | 10.84 | 99 |
| 4-Methyl-2-pentanone (MiBK) | 10.30 | 10.39 | 101 |
| trans-1,3-Dichloropropene | 9.80 | 9.88 | 101 |
| 1,1,2-Trichloroethane | 10.60 | 10.26 | 97 |
| Toluene | 10.60 | 10.25 | 97 |
| 2-Hexanone (MBK) | 10.80 | 10.57 | 98 |
| Dibromochloromethane | 11.00 | 11.51 | 105 |
| 1,2-Dibromoethane | 10.40 | 10.30 | 99 |
| Tetrachloroethene (PCE) | 10.40 | 10.21 | 98 |
| Chlorobenzene | 10.60 | 10.41 | 98 |
| Ethylbenzene | 10.50 | 10.21 | 97 |
| m & p-Xylenes | 20.60 | 19.59 | 95 |
| Bromoform | 10.30 | 10.43 | 101 |
| Styrene | 10.40 | 9.89 | 95 |
| 1,1,2,2-Tetrachloroethane | 10.60 | 10.21 | 96 |
| o-Xylene | 10.60 | 9.94 | 94 |
| 4-Ethyltoluene | 10.40 | 10.41 | 100 |
| 1,3,5-Trimethylbenzene | 10.20 | 9.59 | 94 |
| 1,2,4-Trimethylbenzene | 10.20 | 10.58 | 104 |
| Benzyl Chloride (a-Chlorotoluene) | 10.00 | 11.13 | 111 |
| 1,3-Dichlorobenzene | 10.00 | 10.59 | 106 |
| 1,4-Dichlorobenzene | 10.00 | 9.95 | 100 |
| 1,2-Dichlorobenzene | 10.00 | 10.24 | 102 |
| 1,2,4-Trichlorobenzene | 9.30 | 9.60 | 103 |
| Hexachlorobutadiene | 9.80 | 10.13 | 103 |

* - %REC should be 70-130%


Marcus Hueppe
Laboratory Director





Atmospheric Analysis & Consulting, Inc.

Quality Control/Quality Assurance Report

CLIENT ID : Laboratory Control Spike DATE ANALYZED : 05/20/2013
AAC ID : LCS/LCSD DATE REPORTED : 05/20/2013
MEDIA : Air UNITS : ppbv

TO-15 Laboratory Control Spike Recovery

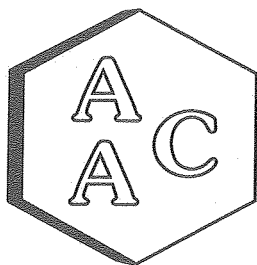
| Compound | Sample Conc. | Spike Added | Spike Res | Dup Spike Res | Spike % Rec * | Spike Dup % Rec * | RPD** % |
|--------------------------|--------------|-------------|-----------|---------------|---------------|-------------------|---------|
| 1,1-Dichloroethene | 0.0 | 10.50 | 10.27 | 9.91 | 98 | 94 | 3.6 |
| Methylene Chloride (DCM) | 0.0 | 10.40 | 9.95 | 9.60 | 96 | 92 | 3.6 |
| Benzene | 0.0 | 10.50 | 9.62 | 9.95 | 92 | 95 | 3.4 |
| Trichloroethene (TCE) | 0.0 | 10.30 | 10.22 | 10.18 | 99 | 99 | 0.4 |
| Toluene | 0.0 | 10.60 | 10.25 | 10.19 | 97 | 96 | 0.6 |
| Tetrachloroethene (PCE) | 0.0 | 10.40 | 10.21 | 10.18 | 98 | 98 | 0.3 |
| Chlorobenzene | 0.0 | 10.60 | 10.41 | 10.35 | 98 | 98 | 0.6 |
| Ethylbenzene | 0.0 | 10.50 | 10.21 | 10.04 | 97 | 96 | 1.7 |
| m & p-Xylenes | 0.0 | 20.60 | 19.59 | 19.20 | 95 | 93 | 2.0 |
| o-Xylene | 0.0 | 10.60 | 9.94 | 10.06 | 94 | 95 | 1.2 |

* Must be 70-130%

** Must be < 25%

Marcus Hueppe
Laboratory Director





Atmospheric Analysis & Consulting, Inc.

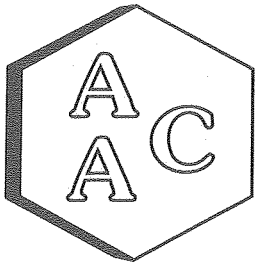
Method Blank Analysis Report

MATRIX : AIR ANALYSIS DATE : 05/20/2013
 UNITS : ppbv REPORT DATE : 05/20/2013

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

| <i>Client ID</i> | Method Blank | RL |
|--------------------------------|--------------|-----|
| <i>AAC ID</i> | MB 052013 | |
| Chlorodifluoromethane | <RL | 0.5 |
| Propene | <RL | 1.0 |
| Dichlorodifluoromethane | <RL | 0.5 |
| Chloromethane | <RL | 0.5 |
| Dichlorotetrafluoroethane | <RL | 0.5 |
| Vinyl Chloride | <RL | 0.5 |
| Methanol | <RL | 5.0 |
| 1,3-Butadiene | <RL | 0.5 |
| Bromomethane | <RL | 0.5 |
| Chloroethane | <RL | 0.5 |
| Dichlorofluoromethane | <RL | 0.5 |
| Ethanol | <RL | 2.0 |
| Vinyl Bromide | <RL | 0.5 |
| Acetone | <RL | 2.0 |
| Trichlorofluoromethane | <RL | 0.5 |
| 2-Propanol (IPA) | <RL | 2.0 |
| Acrylonitrile | <RL | 1.0 |
| 1,1-Dichloroethene | <RL | 0.5 |
| Methylene Chloride (DCM) | <RL | 1.0 |
| Allyl Chloride | <RL | 0.5 |
| Carbon Disulfide | <RL | 0.5 |
| Trichlorotrifluoroethane | <RL | 0.5 |
| trans-1,2-Dichloroethene | <RL | 0.5 |
| 1,1-Dichloroethane | <RL | 0.5 |
| Methyl Tert Butyl Ether (MTBE) | <RL | 0.5 |
| Vinyl Acetate | <RL | 1.0 |
| 2-Butanone (MEK) | <RL | 1.0 |
| cis-1,2-Dichloroethene | <RL | 0.5 |
| Hexane | <RL | 0.5 |
| Chloroform | <RL | 0.5 |
| Ethyl Acetate | <RL | 0.5 |
| Tetrahydrofuran | <RL | 0.5 |
| 1,2-Dichloroethane | <RL | 0.5 |
| 1,1,1-Trichloroethane | <RL | 0.5 |
| Benzene | <RL | 0.5 |
| Carbon Tetrachloride | <RL | 0.5 |
| Cyclohexane | <RL | 0.5 |
| 1,2-Dichloropropane | <RL | 0.5 |
| Bromodichloromethane | <RL | 0.5 |
| 1,4-Dioxane | <RL | 0.5 |
| Trichloroethene (TCE) | <RL | 0.5 |
| 2,2,4-Trimethylpentane | <RL | 0.5 |
| Heptane | <RL | 0.5 |





Atmospheric Analysis & Consulting, Inc.

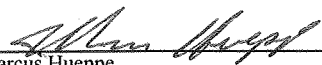
Method Blank Analysis Report

MATRIX : AIR ANALYSIS DATE : 05/20/2013
UNITS : ppbv REPORT DATE : 05/20/2013

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

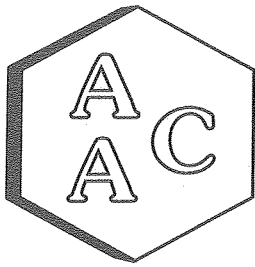
| <i>Client ID</i> | <i>Method Blank</i> | <i>RL</i> |
|------------------------------------|---------------------|-----------|
| <i>AAC ID</i> | <i>MB 052013</i> | |
| cis-1,3-Dichloropropene | <RL | 0.5 |
| 4-Methyl-2-pentanone (MIBK) | <RL | 0.5 |
| trans-1,3-Dichloropropene | <RL | 0.5 |
| 1,1,2-Trichloroethane | <RL | 0.5 |
| Toluene | <RL | 0.5 |
| 2-Hexanone (MBK) | <RL | 0.5 |
| Dibromochloromethane | <RL | 0.5 |
| 1,2-Dibromoethane | <RL | 0.5 |
| Tetrachloroethene (PCE) | <RL | 0.5 |
| Chlorobenzene | <RL | 0.5 |
| Ethylbenzene | <RL | 0.5 |
| m & p-Xylenes | <RL | 1.0 |
| Bromoform | <RL | 0.5 |
| Styrene | <RL | 0.5 |
| 1,1,2,2-Tetrachloroethane | <RL | 0.5 |
| o-Xylene | <RL | 0.5 |
| 4-Ethyltoluene | <RL | 0.5 |
| 1,3,5-Trimethylbenzene | <RL | 0.5 |
| 1,2,4-Trimethylbenzene | <RL | 0.5 |
| Benzyl Chloride (a-Chlorotoluene) | <RL | 0.5 |
| 1,3-Dichlorobenzene | <RL | 0.5 |
| 1,4-Dichlorobenzene | <RL | 0.5 |
| 1,2-Dichlorobenzene | <RL | 0.5 |
| 1,2,4-Trichlorobenzene | <RL | 0.5 |
| Hexachlorobutadiene | <RL | 0.5 |
| System Monitoring Compounds | | |
| BFB-Surrogate Std. % Recovery | 102% | -- |

RL - Reporting Limit



Marcus Hueppe
Laboratory Director





Atmospheric Analysis & Consulting, Inc.

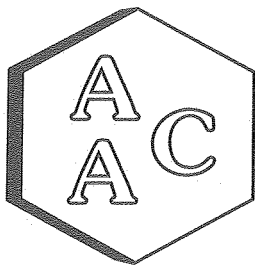
Quality Control/Quality Assurance Report

AAC ID : 130597-63048 **DATE ANALYZED** : 05/20/2013
MATRIX : Air **DATE REPORTED** : 05/20/2013
 UNITS : ppbv

TO-15 Duplicate Analysis

| Compound | Sample Conc | Duplicate Conc | % RPD |
|--------------------------------|-------------|----------------|-------|
| Chlorodifluoromethane | <SRL | <SRL | 0.0 |
| Propene | <SRL | <SRL | 0.0 |
| Dichlorodifluoromethane | <SRL | <SRL | 0.0 |
| Chloromethane | <SRL | <SRL | 0.0 |
| Dichlorotetrafluoroethane | <SRL | <SRL | 0.0 |
| Vinyl Chloride | <SRL | <SRL | 0.0 |
| Methanol | 17.2 | 17.3 | 0.6 |
| 1,3-Butadiene | <SRL | <SRL | 0.0 |
| Bromomethane | <SRL | <SRL | 0.0 |
| Chloroethane | <SRL | <SRL | 0.0 |
| Dichlorofluoromethane | <SRL | <SRL | 0.0 |
| Ethanol | 5.54 | 5.56 | 0.4 |
| Vinyl Bromide | <SRL | <SRL | 0.0 |
| Acetone | 8.24 | 8.04 | 2.5 |
| Trichlorofluoromethane | <SRL | <SRL | 0.0 |
| 2-Propanol (IPA) | <SRL | <SRL | 0.0 |
| Acrylonitrile | <SRL | <SRL | 0.0 |
| 1,1-Dichloroethene | <SRL | <SRL | 0.0 |
| Methylene Chloride (DCM) | <SRL | <SRL | 0.0 |
| Allyl Chloride | <SRL | <SRL | 0.0 |
| Carbon Disulfide | <SRL | <SRL | 0.0 |
| Trichlorotrifluoroethane | <SRL | <SRL | 0.0 |
| trans-1,2-Dichloroethene | <SRL | <SRL | 0.0 |
| 1,1-Dichloroethane | <SRL | <SRL | 0.0 |
| Methyl Tert Butyl Ether (MTBE) | <SRL | <SRL | 0.0 |
| Vinyl Acetate | <SRL | <SRL | 0.0 |
| 2-Butanone (MEK) | <SRL | <SRL | 0.0 |
| cis-1,2-Dichloroethene | <SRL | <SRL | 0.0 |
| Hexane | 1.02 | 0.98 | 4.0 |
| Chloroform | <SRL | <SRL | 0.0 |
| Ethyl Acetate | <SRL | <SRL | 0.0 |
| Tetrahydrofuran | <SRL | <SRL | 0.0 |
| 1,2-Dichloroethane | <SRL | <SRL | 0.0 |
| 1,1,1-Trichloroethane | <SRL | <SRL | 0.0 |
| Benzene | <SRL | <SRL | 0.0 |
| Carbon Tetrachloride | <SRL | <SRL | 0.0 |





Atmospheric Analysis & Consulting, Inc.

Quality Control/Quality Assurance Report

AAC ID : 130597-63048 DATE ANALYZED : 05/20/2013
MATRIX : Air DATE REPORTED : 05/20/2013
UNITS : ppbv

TO-15 Duplicate Analysis

| Compound | Sample Conc | Duplicate Conc | % RPD |
|------------------------------------|-------------|----------------|-------|
| Cyclohexane | <SRL | <SRL | 0.0 |
| 1,2-Dichloropropane | <SRL | <SRL | 0.0 |
| Bromodichloromethane | <SRL | <SRL | 0.0 |
| 1,4-Dioxane | <SRL | <SRL | 0.0 |
| Trichloroethene (TCE) | <SRL | <SRL | 0.0 |
| 2,2,4-Trimethylpentane | <SRL | <SRL | 0.0 |
| Heptane | <SRL | <SRL | 0.0 |
| cis-1,3-Dichloropropene | <SRL | <SRL | 0.0 |
| 4-Methyl-2-pentanone (MIBK) | <SRL | <SRL | 0.0 |
| trans-1,3-Dichloropropene | <SRL | <SRL | 0.0 |
| 1,1,2-Trichloroethane | <SRL | <SRL | 0.0 |
| Toluene | 1.17 | 1.15 | 1.7 |
| 2-Hexanone (MBK) | <SRL | <SRL | 0.0 |
| Dibromochloromethane | <SRL | <SRL | 0.0 |
| 1,2-Dibromoethane | <SRL | <SRL | 0.0 |
| Tetrachloroethene (PCE) | <SRL | <SRL | 0.0 |
| Chlorobenzene | <SRL | <SRL | 0.0 |
| Ethylbenzene | <SRL | <SRL | 0.0 |
| m & p-Xylenes | <SRL | <SRL | 0.0 |
| Bromoform | <SRL | <SRL | 0.0 |
| Styrene | <SRL | <SRL | 0.0 |
| 1,1,2,2-Tetrachloroethane | <SRL | <SRL | 0.0 |
| o-Xylene | <SRL | <SRL | 0.0 |
| 4-Ethyltoluene | <SRL | <SRL | 0.0 |
| 1,3,5-Trimethylbenzene | <SRL | <SRL | 0.0 |
| 1,2,4-Trimethylbenzene | <SRL | <SRL | 0.0 |
| Benzyl Chloride (a-Chlorotoluene) | <SRL | <SRL | 0.0 |
| 1,3-Dichlorobenzene | <SRL | <SRL | 0.0 |
| 1,4-Dichlorobenzene | <SRL | <SRL | 0.0 |
| 1,2-Dichlorobenzene | <SRL | <SRL | 0.0 |
| 1,2,4-Trichlorobenzene | <SRL | <SRL | 0.0 |
| Hexachlorobutadiene | <SRL | <SRL | 0.0 |
| System Monitoring Compounds | | | |
| BFB-Surrogate Std. % Recovery | 102% | 104% | 2.0 |

SRL - Sample Reporting Limit


Marcus Hueppe
Laboratory Director



**TO-15
RAW
DATA**

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201305.D
 Acq On : 20 May 2013 11:44
 Operator : JJG
 Sample : 130597-63048 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 20 15:18:58 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------|--------|------|----------|-------|-------|----------|
| Internal Standards | | | | | | |
| 1) Bromochloromethane | 12.350 | 128 | 145590 | 10.00 | ppbv | 0.00 |
| 36) 1,4-Difluorobenzene | 14.579 | 114 | 790345 | 10.00 | ppbv | 0.00 |
| 55) Chlorobenzene-d5 | 20.285 | 117 | 750102 | 10.00 | ppbv | 0.00 |

| | | | | | | |
|--------------------------------|--------|-----|--------|-------|------|------|
| System Monitoring Compounds | | | | | | |
| 63) 4-Bromofluorobenzene (BFB) | 22.710 | 174 | 479494 | 10.21 | ppbv | 0.00 |

Spiked Amount 10.000 Recovery = 102.10%

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue | Dev(Min) |
|---------------------------------|--------|------|----------|------|-------|--------|----------|
| 2) Chlorodifluoromethane | 4.818 | 51 | 4516 | 0.16 | ppbv | # 96 | |
| 3) Propene | 0.000 | | 0 | N.D. | d | | |
| 4) Dichlorodifluoromethane | 4.890 | 85 | 14589 | 0.29 | ppbv | 98 | |
| 5) Chloromethane | 5.288 | 52 | 975 | 0.21 | ppbv | # 10 | |
| 6) Dichlorotetrafluoroethane | 5.324 | 135 | 204 | N.D. | | | |
| 7) VinylChloride | 0.000 | | 0 | N.D. | | | |
| 8) Methanol | 5.867 | 31 | 39778 | 8.56 | ppbv | | |
| 9) 1,3-Butadiene | 0.000 | | 0 | N.D. | | | |
| 10) Bromomethane | 0.000 | | 0 | N.D. | ppbv | | 0.00 |
| 11) Chloroethane | 0.000 | | 0 | N.D. | ppbv | | 0.00 |
| 12) Dichlorofluoromethane | 0.000 | | 0 | N.D. | ppbv | | 0.00 |
| 13) Ethanol | 7.134 | 45 | 17153 | 2.76 | ppbv | | |
| 14) VinylBromide | 0.000 | | 0 | N.D. | | | |
| 15) Acetone | 7.984 | 58 | 32335 | 4.10 | ppbv | | 0.00 |
| 16) Trichlorofluoromethane | 7.659 | 103 | 4061 | 0.14 | ppbv | # 95 | |
| 17) 2-Propanol (IPA) | 8.238 | 45 | 6570 | 0.24 | ppbv | 10% | |
| 18) Acrylonitrile | 0.000 | | 0 | N.D. | | | |
| 19) 1,1-Dichloroethene | 0.000 | | 0 | N.D. | | Qvalue | |
| 20) MethyleneChloride (DCM) | 0.000 | | 0 | N.D. | ppbv | # 96 | |
| 21) AllylChloride | 0.000 | | 0 | N.D. | d | | |
| 22) CarbonDisulfide | 0.000 | | 0 | N.D. | ppbv | 98 | |
| 23) Trichlorotrifluoroethane | 0.000 | | 0 | N.D. | d | # 10 | |
| 24) trans-1,2-Dichloroethene | 0.000 | | 0 | N.D. | | | |
| 25) 1,1-Dichloroethane | 0.000 | | 0 | N.D. | | | |
| 26) MethylTertButylether (M...) | 0.000 | | 0 | N.D. | ppbv | | |
| 27) VinylAcetate | 0.000 | | 0 | N.D. | d | | |
| 28) 2-Butanone (MEK) | 11.494 | 72 | 3961 | 0.49 | ppbv | # 35 | |
| 29) cis-1,2-Dichloroethene | 0.000 | | 0 | N.D. | | | |
| 30) Hexane | 11.458 | 86 | 1972 | 0.51 | ppbv | # 49 | |
| 31) Chloroform | 12.493 | 83 | 995 | N.D. | ppbv | | |
| 32) EthylAcetate | 12.118 | 43 | 1448 | N.D. | | | |

Handwritten signature

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201305.D
 Acq On : 20 May 2013 11:44
 Operator : JJG
 Sample : 130597-63048 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 20 15:18:58 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

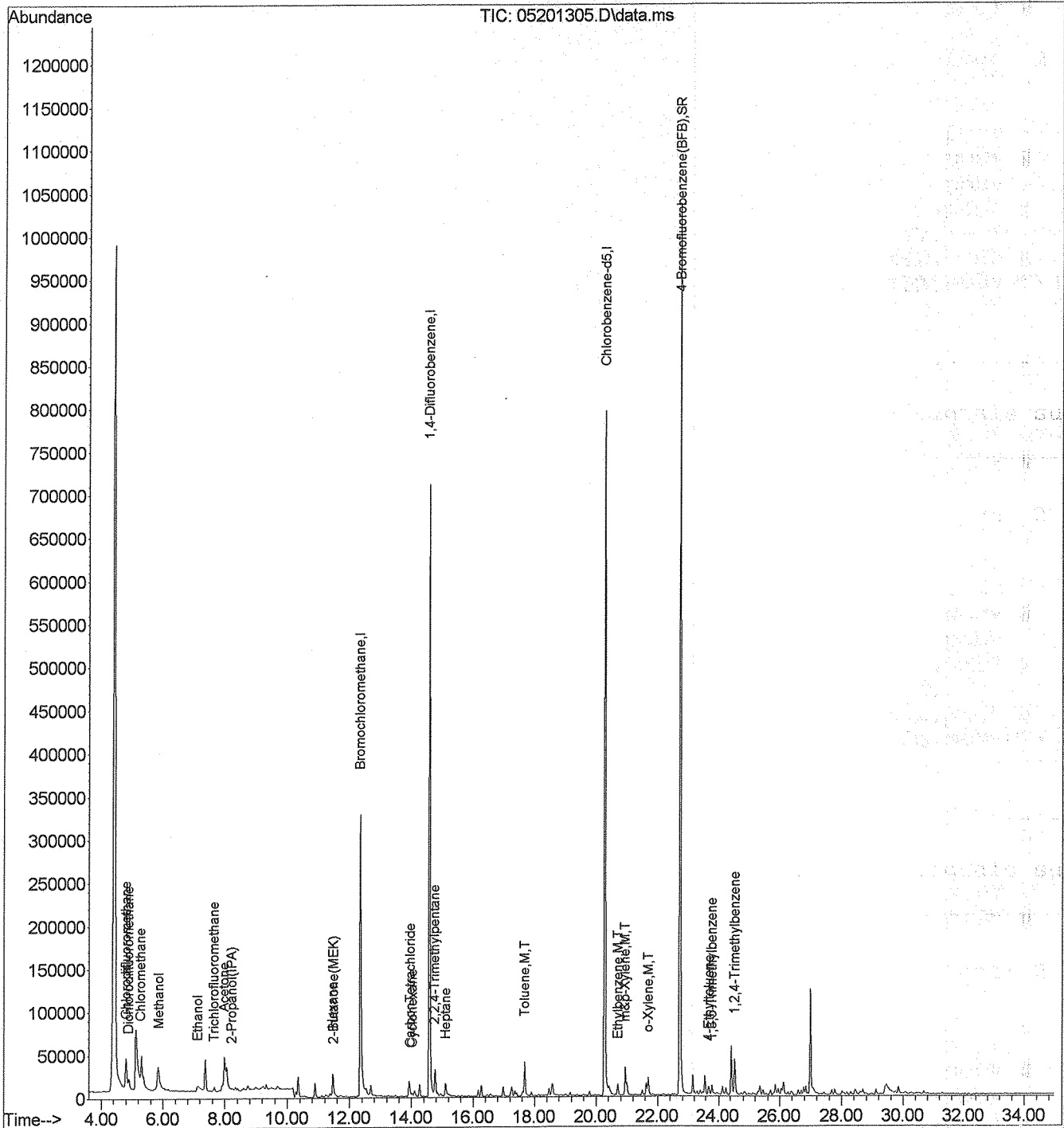
| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------------|--------|------|----------|------|--------------|----------|
| 33) Tetrahydrofuran | 0.000 | | 0 | N.D. | | |
| 34) 1,2-Dichloroethane | 13.634 | 62 | 123 | N.D. | | |
| 35) 1,1,1-Trichloroethane | 0.000 | | 0 | N.D. | | |
| 37) Benzene | 0.000 | | 0 | N.D. | d | |
| 38) CarbonTetrachloride | 13.973 | 117 | 2480 | 0.05 | ppbv | 93 |
| 39) Cyclohexane | 14.026 | 69 | 612 | 0.07 | ppbv # | 92 |
| 40) 1,2-Dichloropropane | 15.257 | 63 | 116 | N.D. | | |
| 41) Bromodichloromethane | 0.000 | | 0 | N.D. | | |
| 42) 1,4-Dioxane | 0.000 | | 0 | N.D. | | |
| 43) Trichloroethene (TCE) | 15.292 | 130 | 259 | N.D. | | |
| 44) 2,2,4-Trimethylpentane | 14.757 | 57 | 39593 | 0.38 | ppbv | 96 |
| 45) Heptane | 15.114 | 71 | 4039 | 0.22 | ppbv # | 86 |
| 46) cis-1,3-Dichloropropene | 0.000 | | 0 | N.D. | | |
| 47) 4-Methyl-2-pentanone (M...) | 16.576 | 58 | 710 | N.D. | | |
| 48) trans-1,3-Dichloropropene | 17.682 | 75 | 154 | N.D. | | |
| 49) 1,1,2-Trichloroethane | 0.000 | | 0 | N.D. | | |
| 50) Toluene | 17.682 | 91 | 42772 | 0.58 | ppbv Dev (M) | 99 |
| 51) 2-Hexanone (MBK) | 0.000 | | 0 | N.D. | | |
| 52) Dibromochloromethane | 0.000 | | 0 | N.D. | | |
| 53) 1,2-Dibromoethane | 0.000 | | 0 | N.D. | | |
| 54) Tetrachloroethene (PCE) | 19.019 | 166 | 116 | N.D. | | |
| 56) Chlorobenzene | 0.000 | | 0 | N.D. | | |
| 57) Ethylbenzene | 20.713 | 91 | 14549 | 0.15 | ppbv | 98 |
| 58) m&p-Xylene | 20.945 | 106 | 19939 | 0.51 | ppbv # | 89 |
| 59) Bromoform | 0.000 | | 0 | N.D. | | |
| 60) Styrene | 21.694 | 104 | 1167 | N.D. | | |
| 61) 1,1,2,2-Tetrachloroethane | 22.354 | 83 | 122 | N.D. | | |
| 62) o-Xylene | 21.694 | 91 | 17217 | 0.22 | ppbv | 97 |
| 64) 4-Ethyltoluene | 23.691 | 120 | 3288 | 0.10 | ppbv | 95 |
| 65) 1,3,5-Trimethylbenzene | 23.780 | 120 | 4534 | 0.10 | ppbv # | 91 |
| 66) 1,2,4-Trimethylbenzene | 24.529 | 120 | 16063 | 0.35 | ppbv | 94 |
| 67) BenzylChloride (a-Chlor...) | 25.189 | 91 | 424 | N.D. | | |
| 68) 1,3-Dichlorobenzene | 25.064 | 146 | 678 | N.D. | | |
| 69) 1,4-Dichlorobenzene | 25.296 | 146 | 1203 | N.D. | | |
| 70) 1,2-Dichlorobenzene | 25.867 | 146 | 924 | N.D. | ppbv Dev (M) | 99 |
| 71) 1,2,4-Trichlorobenzene | 29.451 | 180 | 2361 | N.D. | | |
| 72) Hexachlorobutadiene | 30.075 | 225 | 876 | N.D. | | |

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

Postols

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201305.D
 Acq On : 20 May 2013 11:44
 Operator : JJG
 Sample : 130597-63048 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 20 15:18:58 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration



Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201307.D
 Acq On : 20 May 2013 13:20
 Operator : JJG
 Sample : 130597-63049 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: May 20 17:50:50 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev (Min) |
|-------------------------|--------|------|----------|-------|-------|-----------|
| Internal Standards | | | | | | |
| 1) Bromochloromethane | 12.350 | 128 | 142515 | 10.00 | ppbv | 0.00 |
| 36) 1,4-Difluorobenzene | 14.579 | 114 | 791178 | 10.00 | ppbv | 0.00 |
| 55) Chlorobenzene-d5 | 20.285 | 117 | 740079 | 10.00 | ppbv | 0.00 |

| | | | | | | |
|--------------------------------|--------|-----|--------|-------|------|------|
| System Monitoring Compounds | | | | | | |
| 63) 4-Bromofluorobenzene (BFB) | 22.710 | 174 | 467395 | 10.09 | ppbv | 0.00 |

Spiked Amount 10.000 Recovery = 100.90%

| Target Compounds | R.T. | QIon | Response | Conc | Units | Dev (Min) | Qvalue |
|---------------------------------|--------|------|----------|------|-------|-----------|-----------|
| 2) Chlorodifluoromethane | 4.835 | 51 | 3076 | 0.11 | ppbv | | # 91 |
| 3) Propene | 0.000 | | 0 | N.D. | d | | |
| 4) Dichlorodifluoromethane | 4.908 | 85 | 10044 | 0.21 | ppbv | | 97 |
| 5) Chloromethane | 5.288 | 52 | 818 | 0.18 | ppbv | | # 1 |
| 6) Dichlorotetrafluoroethane | 0.000 | | 0 | N.D. | | | |
| 7) VinylChloride | 0.000 | | 0 | N.D. | | | |
| 8) Methanol | 5.867 | 31 | 27227 | 5.93 | ppbv | | |
| 9) 1,3-Butadiene | 0.000 | | 0 | N.D. | | | |
| 10) Bromomethane | 0.000 | | 0 | N.D. | d | | 0.00 |
| 11) Chloroethane | 0.000 | | 0 | N.D. | d | | 0.00 |
| 12) Dichlorofluoromethane | 0.000 | | 0 | N.D. | ppbv | | 0.00 |
| 13) Ethanol | 7.134 | 45 | 15223 | 2.50 | ppbv | | |
| 14) VinylBromide | 0.000 | | 0 | N.D. | | | |
| 15) Acetone | 8.002 | 58 | 19738 | 2.56 | ppbv | | 0.00 |
| 16) Trichlorofluoromethane | 7.659 | 103 | 2361 | 0.08 | ppbv | | # 93 |
| 17) 2-Propanol (IPA) | 8.238 | 45 | 14188 | 0.54 | ppbv | | 100% |
| 18) Acrylonitrile | 0.000 | | 0 | N.D. | | | |
| 19) 1,1-Dichloroethene | 0.000 | | 0 | N.D. | | | Qvalue |
| 20) MethyleneChloride (DCM) | 0.000 | | 0 | N.D. | ppbv | | # 91 |
| 21) AllylChloride | 0.000 | | 0 | N.D. | d | | |
| 22) CarbonDisulfide | 0.000 | | 0 | N.D. | d | | 97 |
| 23) Trichlorotrifluoroethane | 8.998 | 103 | 719 | N.D. | ppbv | | # 1 |
| 24) trans-1,2-Dichloroethene | 0.000 | | 0 | N.D. | | | |
| 25) 1,1-Dichloroethane | 0.000 | | 0 | N.D. | | | Dev (Min) |
| 26) MethylTertButylEther (M...) | 0.000 | | 0 | N.D. | | | |
| 27) VinylAcetate | 10.888 | 43 | 1396 | N.D. | | | |
| 28) 2-Butanone (MEK) | 11.512 | 72 | 2185 | 0.28 | ppbv | | # 67 |
| 29) cis-1,2-Dichloroethene | 0.000 | | 0 | N.D. | | | 0.00 |
| 30) Hexane | 0.000 | | 0 | N.D. | d | | 0.00 |
| 31) Chloroform | 12.493 | 83 | 334 | N.D. | | | |
| 32) EthylAcetate | 12.136 | 43 | 825 | N.D. | | | |

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201307.D
 Acq On : 20 May 2013 13:20
 Operator : JJG
 Sample : 130597-63049 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 3 Sample Multiplier: 1

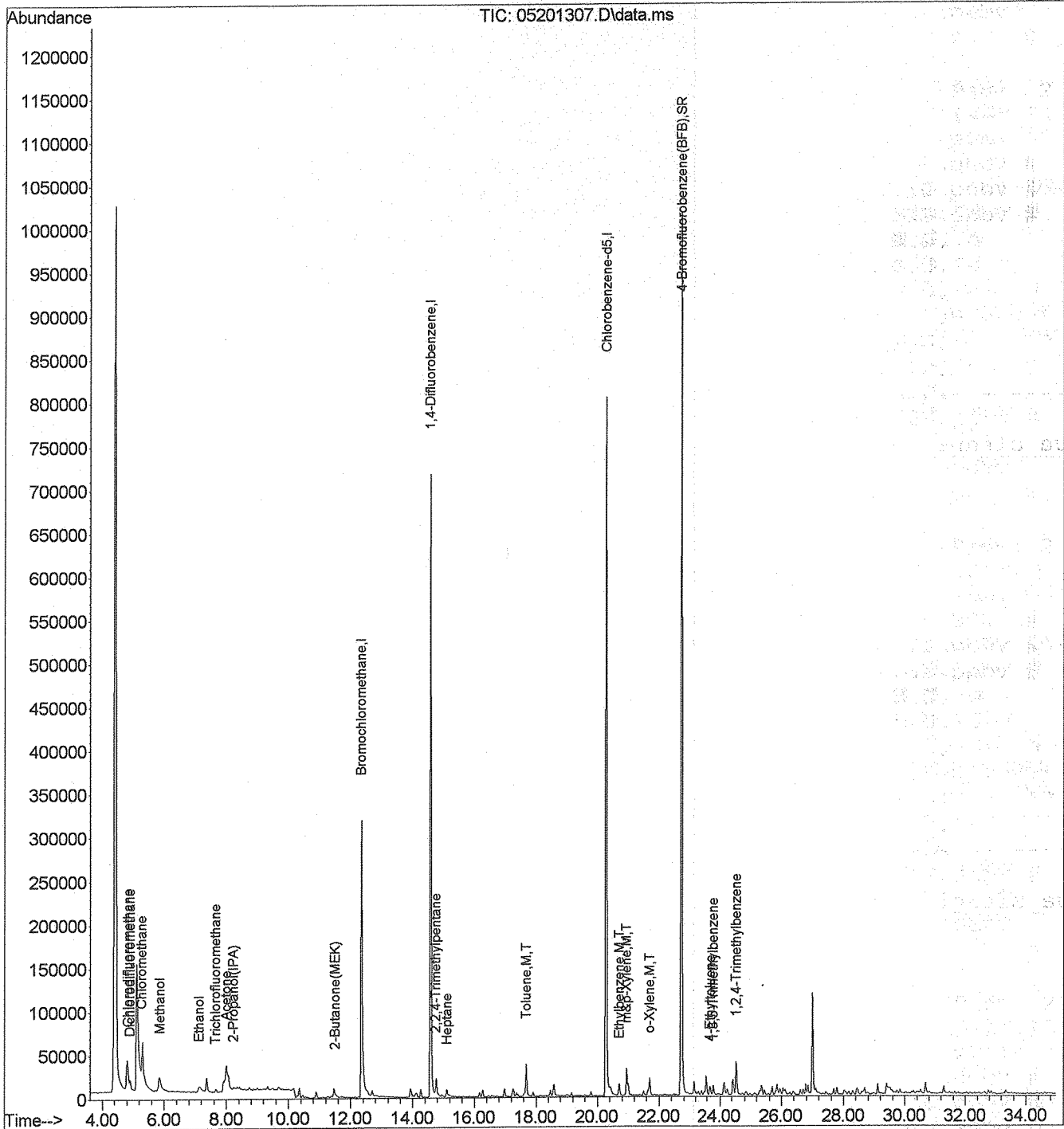
Quant Time: May 20 17:50:50 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev (Min) |
|---------------------------------|--------|------|----------|------|--------|-----------|
| 33) Tetrahydrofuran | 0.000 | | 0 | N.D. | | |
| 34) 1,2-Dichloroethane | 0.000 | | 0 | N.D. | | |
| 35) 1,1,1-Trichloroethane | 0.000 | | 0 | N.D. | | |
| 37) Benzene | 0.000 | | 0 | N.D. | d | |
| 38) CarbonTetrachloride | 13.973 | 117 | 1270 | N.D. | | |
| 39) Cyclohexane | 14.026 | 69 | 124 | N.D. | | |
| 40) 1,2-Dichloropropane | 0.000 | | 0 | N.D. | | |
| 41) Bromodichloromethane | 0.000 | | 0 | N.D. | | |
| 42) 1,4-Dioxane | 0.000 | | 0 | N.D. | | |
| 43) Trichloroethene (TCE) | 15.292 | 130 | 294 | N.D. | | |
| 44) 2,2,4-Trimethylpentane | 14.757 | 57 | 26450 | 0.25 | ppbv | 98 |
| 45) Heptane | 15.114 | 71 | 2449 | 0.13 | ppbv | 92 |
| 46) cis-1,3-Dichloropropene | 0.000 | | 0 | N.D. | | |
| 47) 4-Methyl-2-pentanone (M...) | 16.576 | 58 | 256 | N.D. | | |
| 48) trans-1,3-Dichloropropene | 17.682 | 75 | 111 | N.D. | | |
| 49) 1,1,2-Trichloroethane | 0.000 | | 0 | N.D. | | |
| 50) Toluene | 17.682 | 91 | 40216 | 0.54 | ppbv | Dev (Min) |
| 51) 2-Hexanone (MBK) | 0.000 | | 0 | N.D. | | |
| 52) Dibromochloromethane | 0.000 | | 0 | N.D. | | |
| 53) 1,2-Dibromoethane | 0.000 | | 0 | N.D. | | |
| 54) Tetrachloroethene (PCE) | 0.000 | | 0 | N.D. | | |
| 56) Chlorobenzene | 20.285 | 114 | 139 | N.D. | | |
| 57) Ethylbenzene | 20.713 | 91 | 15193 | 0.16 | ppbv | 97 |
| 58) m&p-Xylene | 20.945 | 106 | 20041 | 0.52 | ppbv # | 94 |
| 59) Bromoform | 0.000 | | 0 | N.D. | | |
| 60) Styrene | 21.676 | 104 | 862 | N.D. | | |
| 61) 1,1,2,2-Tetrachloroethane | 0.000 | | 0 | N.D. | | |
| 62) o-Xylene | 21.694 | 91 | 17786 | 0.23 | ppbv | 92 |
| 64) 4-Ethyltoluene | 23.691 | 120 | 2876 | 0.09 | ppbv # | 89 |
| 65) 1,3,5-Trimethylbenzene | 23.780 | 120 | 4500 | 0.10 | ppbv # | 92 |
| 66) 1,2,4-Trimethylbenzene | 24.529 | 120 | 15887 | 0.35 | ppbv # | 93 |
| 67) BenzylChloride (a-Chlor...) | 25.189 | 91 | 107 | N.D. | | |
| 68) 1,3-Dichlorobenzene | 25.082 | 146 | 228 | N.D. | | |
| 69) 1,4-Dichlorobenzene | 25.296 | 146 | 764 | N.D. | | |
| 70) 1,2-Dichlorobenzene | 25.849 | 146 | 128 | N.D. | | |
| 71) 1,2,4-Trichlorobenzene | 29.451 | 180 | 798 | N.D. | | |
| 72) Hexachlorobutadiene | 30.075 | 225 | 181 | N.D. | | |

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

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201307.D
 Acq On : 20 May 2013 13:20
 Operator : JJG
 Sample : 130597-63049 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: May 20 17:50:50 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration



Handwritten signature

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201308.D
 Acq On : 20 May 2013 14:08
 Operator : JJG
 Sample : 130597-63050 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: May 20 18:41:17 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev (Min) |
|-------------------------|--------|------|----------|-------|-------|-----------|
| Internal Standards | | | | | | |
| 1) Bromochloromethane | 12.350 | 128 | 141218 | 10.00 | ppbv | 0.00 |
| 36) 1,4-Difluorobenzene | 14.579 | 114 | 778335 | 10.00 | ppbv | 0.00 |
| 55) Chlorobenzene-d5 | 20.285 | 117 | 732351 | 10.00 | ppbv | 0.00 |

| | | | | | | |
|--------------------------------|--------|-----|--------|-------|------|------|
| System Monitoring Compounds | | | | | | |
| 63) 4-Bromofluorobenzene (BFB) | 22.710 | 174 | 475368 | 10.37 | ppbv | 0.00 |

Spiked Amount 10.000 Recovery = 103.70%

| Target Compounds | R.T. | QIon | Response | Conc | Units | Dev (Min) | Qvalue |
|---------------------------------|--------|------|----------|-------|-------|-----------|-----------|
| 2) Chlorodifluoromethane | 4.817 | 51 | 5268 | 0.19 | ppbv | | # 94 |
| 3) Propene | 0.000 | | 0 | N.D. | d | | |
| 4) Dichlorodifluoromethane | 4.908 | 85 | 17441 | 0.36 | ppbv | | 98 |
| 5) Chloromethane | 5.288 | 52 | 1289 | 0.28 | ppbv | | # 1 |
| 6) Dichlorotetrafluoroethane | 5.324 | 135 | 257 | N.D. | | | |
| 7) VinylChloride | 0.000 | | 0 | N.D. | | | |
| 8) Methanol | 5.849 | 31 | 96765 | 22.53 | ppbv | | |
| 9) 1,3-Butadiene | 0.000 | | 0 | N.D. | | | |
| 10) Bromomethane | 0.000 | | 0 | N.D. | d | | 0.00 |
| 11) Chloroethane | 0.000 | | 0 | N.D. | | | 0.00 |
| 12) Dichlorofluoromethane | 0.000 | | 0 | N.D. | | | 0.00 |
| 13) Ethanol | 7.079 | 45 | 22744 | 3.78 | ppbv | | |
| 14) VinylBromide | 0.000 | | 0 | N.D. | | | |
| 15) Acetone | 7.966 | 58 | 39431 | 5.16 | ppbv | | 0.00 |
| 16) Trichlorofluoromethane | 7.658 | 103 | 4517 | 0.16 | ppbv | | # 98 |
| 17) 2-Propanol (IPA) | 8.219 | 45 | 12522 | 0.48 | ppbv | | |
| 18) Acrylonitrile | 0.000 | | 0 | N.D. | | | |
| 19) 1,1-Dichloroethene | 0.000 | | 0 | N.D. | | | Qvalue |
| 20) MethyleneChloride (DCM) | 0.000 | | 0 | N.D. | dbv | | # 94 |
| 21) AllylChloride | 0.000 | | 0 | N.D. | d | | |
| 22) CarbonDisulfide | 0.000 | | 0 | N.D. | dbv | | 98 |
| 23) Trichlorotrifluoroethane | 0.000 | | 0 | N.D. | d | | # 1 |
| 24) trans-1,2-Dichloroethene | 0.000 | | 0 | N.D. | | | |
| 25) 1,1-Dichloroethane | 0.000 | | 0 | N.D. | | | Dev (Min) |
| 26) MethylTertButylEther (M...) | 0.000 | | 0 | N.D. | ppbv | | |
| 27) VinylAcetate | 0.000 | | 0 | N.D. | d | | |
| 28) 2-Butanone (MEK) | 11.459 | 72 | 8787 | 1.12 | ppbv | | # 01 |
| 29) cis-1,2-Dichloroethene | 0.000 | | 0 | N.D. | | | 0.00 |
| 30) Hexane | 11.459 | 86 | 2051 | 0.55 | ppbv | | # 42 |
| 31) Chloroform | 12.511 | 83 | 891 | N.D. | ppbv | | |
| 32) EthylAcetate | 0.000 | | 0 | N.D. | d | | |

05/20/13

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201308.D
 Acq On : 20 May 2013 14:08
 Operator : JJG
 Sample : 130597-63050 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: May 20 18:41:17 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

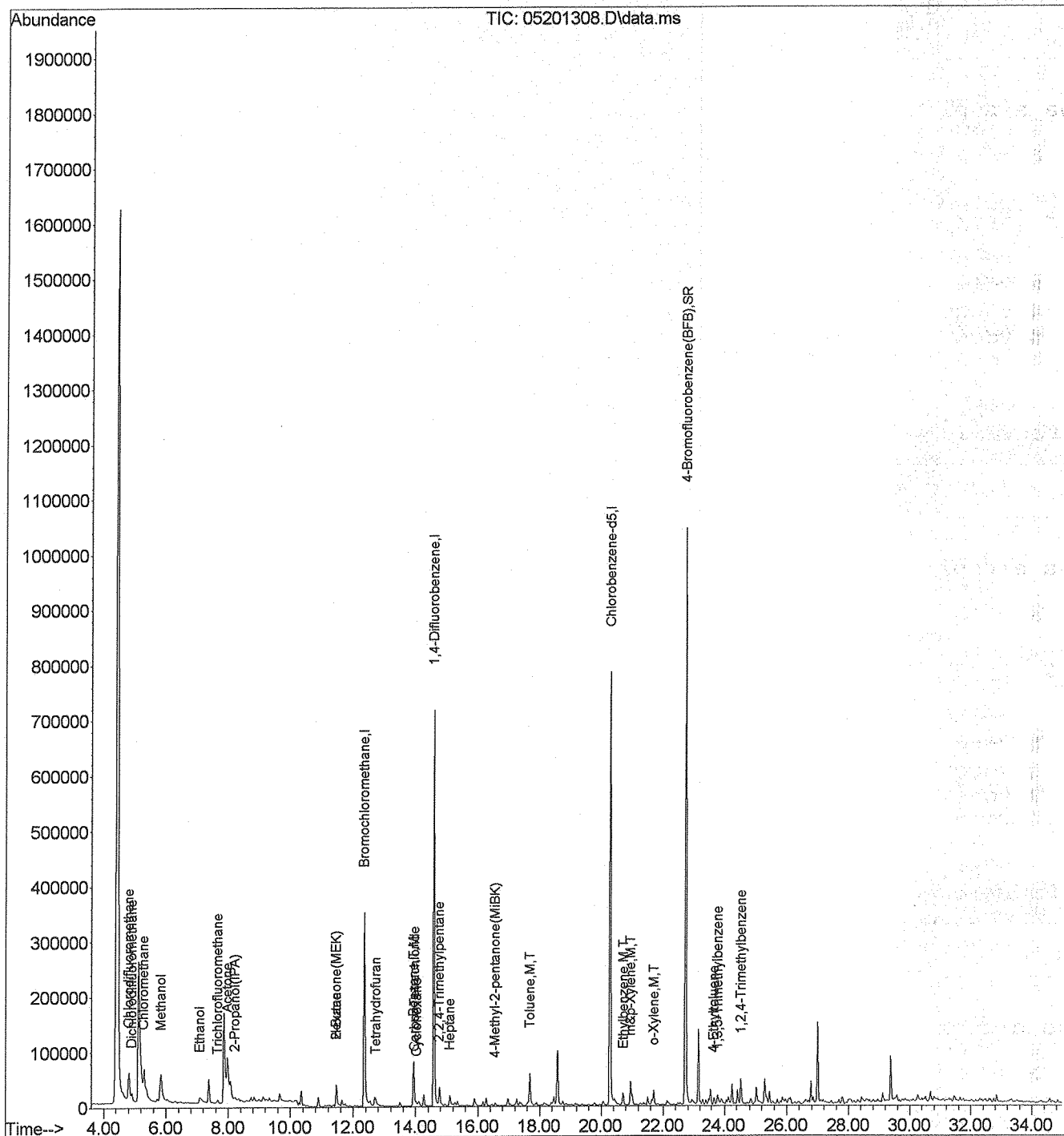
| Compound | R.T. | QIon | Response | Conc | Units | Dev (Min) |
|---------------------------------|--------|------|----------|------|--------|-----------|
| 33) Tetrahydrofuran | 12.725 | 72 | 4525 | 0.57 | ppbv # | 78 |
| 34) 1,2-Dichloroethane | 13.616 | 62 | 112 | N.D. | | |
| 35) 1,1,1-Trichloroethane | 0.000 | | 0 | N.D. | | |
| 37) Benzene | 13.937 | 78 | 80800 | 1.41 | ppbv | 97 |
| 38) CarbonTetrachloride | 13.973 | 117 | 2708 | 0.06 | ppbv # | 97 |
| 39) Cyclohexane | 14.026 | 69 | 713 | 0.08 | ppbv # | 1 |
| 40) 1,2-Dichloropropane | 15.417 | 63 | 120 | N.D. | | |
| 41) Bromodichloromethane | 0.000 | | 0 | N.D. | | |
| 42) 1,4-Dioxane | 0.000 | | 0 | N.D. | | |
| 43) Trichloroethene (TCE) | 0.000 | | 0 | N.D. | | |
| 44) 2,2,4-Trimethylpentane | 14.757 | 57 | 39602 | 0.39 | ppbv # | 94 |
| 45) Heptane | 15.096 | 71 | 5093 | 0.28 | ppbv # | 73 |
| 46) cis-1,3-Dichloropropene | 0.000 | | 0 | N.D. | | |
| 47) 4-Methyl-2-pentanone (M...) | 16.558 | 58 | 1332 | 0.06 | ppbv # | 68 |
| 48) trans-1,3-Dichloropropene | 17.682 | 75 | 401 | N.D. | | |
| 49) 1,1,2-Trichloroethane | 0.000 | | 0 | N.D. | d | |
| 50) Toluene | 17.682 | 91 | 61267(m) | 0.84 | ppbv | DEV (MIN) |
| 51) 2-Hexanone (MBK) | 18.199 | 58 | 672 | N.D. | | |
| 52) Dibromochloromethane | 19.019 | 129 | 119 | N.D. | ppbv # | 78 |
| 53) 1,2-Dibromoethane | 0.000 | | 0 | N.D. | | |
| 54) Tetrachloroethene (PCE) | 19.019 | 166 | 114 | N.D. | | |
| 56) Chlorobenzene | 0.000 | | 0 | N.D. | | 97 |
| 57) Ethylbenzene | 20.713 | 91 | 22552 | 0.23 | ppbv # | 98 |
| 58) m&p-Xylene | 20.945 | 106 | 25377 | 0.67 | ppbv # | 93 |
| 59) Bromoform | 0.000 | | 0 | N.D. | | |
| 60) Styrene | 21.658 | 104 | 1440 | N.D. | | |
| 61) 1,1,2,2-Tetrachloroethane | 0.000 | | 0 | N.D. | | |
| 62) o-Xylene | 21.694 | 91 | 20962 | 0.27 | ppbv | 99 |
| 64) 4-Ethyltoluene | 23.673 | 120 | 3435 | 0.11 | ppbv # | 87 |
| 65) 1,3,5-Trimethylbenzene | 23.780 | 120 | 5194 | 0.11 | ppbv # | 70 |
| 66) 1,2,4-Trimethylbenzene | 24.529 | 120 | 17570 | 0.39 | ppbv # | 87 |
| 67) BenzylChloride (a-Chlor...) | 25.189 | 91 | 117 | N.D. | ppbv # | 63 |
| 68) 1,3-Dichlorobenzene | 25.064 | 146 | 141 | N.D. | | |
| 69) 1,4-Dichlorobenzene | 0.000 | | 0 | N.D. | d | |
| 70) 1,2-Dichlorobenzene | 25.849 | 146 | 281 | N.D. | ppbv # | 107 |
| 71) 1,2,4-Trichlorobenzene | 29.451 | 180 | 611 | N.D. | | |
| 72) Hexachlorobutadiene | 30.075 | 225 | 118 | N.D. | | 78 |

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

Handwritten signature and date:
 05/20/13

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201308.D
 Acq On : 20 May 2013 14:08
 Operator : JJG
 Sample : 130597-63050 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: May 20 18:41:17 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration



Handwritten signature and date: JJG 05/20/13

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201309.D
 Acq On : 20 May 2013 14:56
 Operator : JJG
 Sample : 130597-63051 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: May 20 18:44:03 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------|--------|------|----------|-------|-------|----------|
| Internal Standards | | | | | | |
| 1) Bromochloromethane | 12.350 | 128 | 142936 | 10.00 | ppbv | 0.00 |
| 36) 1,4-Difluorobenzene | 14.579 | 114 | 779054 | 10.00 | ppbv | 0.00 |
| 55) Chlorobenzene-d5 | 20.285 | 117 | 733160 | 10.00 | ppbv | 0.00 |

| | | | | | | |
|--------------------------------|--------|-----|--------|-------|------|------|
| System Monitoring Compounds | | | | | | |
| 63) 4-Bromofluorobenzene (BFB) | 22.711 | 174 | 473626 | 10.32 | ppbv | 0.00 |

Spiked Amount 10.000 Recovery = 103.20%

| Target Compounds | R.T. | QIon | Response | Conc | Units | Dev(Min) | Qvalue |
|---------------------------------|--------|------|----------|-------|-------|----------|-----------|
| 2) Chlorodifluoromethane | 4.836 | 51 | 1751 | 0.06 | ppbv | # | 69 |
| 3) Propene | 4.781 | 42 | 6370 | 0.86 | ppbv | # | 81 |
| 4) Dichlorodifluoromethane | 4.908 | 85 | 5626 | 0.12 | ppbv | | 97 |
| 5) Chloromethane | 5.288 | 52 | 543 | 0.12 | ppbv | # | 71 |
| 6) Dichlorotetrafluoroethane | 0.000 | | 0 | N.D. | | | |
| 7) VinylChloride | 0.000 | | 0 | N.D. | | | |
| 8) Methanol | 5.849 | 31 | 47984 | 10.59 | ppbv | | |
| 9) 1,3-Butadiene | 0.000 | | 0 | N.D. | | | |
| 10) Bromomethane | 0.000 | | 0 | N.D. | | | 0.00 |
| 11) Chloroethane | 0.000 | | 0 | N.D. | | | 0.00 |
| 12) Dichlorofluoromethane | 0.000 | | 0 | N.D. | | | 0.00 |
| 13) Ethanol | 7.080 | 45 | 33682 | 5.53 | ppbv | | |
| 14) VinylBromide | 0.000 | | 0 | N.D. | | | |
| 15) Acetone | 7.984 | 58 | 32740 | 4.23 | ppbv | | 0.00 |
| 16) Trichlorofluoromethane | 0.000 | | 0 | N.D. | | | |
| 17) 2-Propanol (IPA) | 8.183 | 45 | 140470 | 5.30 | ppbv | 20% | 75 |
| 18) Acrylonitrile | 0.000 | | 0 | N.D. | | | |
| 19) 1,1-Dichloroethene | 0.000 | | 0 | N.D. | | | Qvalue |
| 20) MethyleneChloride (DCM) | 0.000 | | 0 | N.D. | | | # 69 |
| 21) AllylChloride | 0.000 | | 0 | N.D. | | | # 81 |
| 22) CarbonDisulfide | 0.000 | | 0 | N.D. | | | 97 |
| 23) Trichlorotrifluoroethane | 8.998 | 103 | 375 | N.D. | | | # 71 |
| 24) trans-1,2-Dichloroethene | 10.424 | 96 | 2082 | 0.12 | ppbv | # | 81 |
| 25) 1,1-Dichloroethane | 0.000 | | 0 | N.D. | | | Dev (Min) |
| 26) MethylTertButylether (M...) | 10.496 | 73 | 114 | N.D. | | | |
| 27) VinylAcetate | 0.000 | | 0 | N.D. | | | d |
| 28) 2-Butanone (MEK) | 11.494 | 72 | 3564 | 0.45 | ppbv | # | 71 |
| 29) cis-1,2-Dichloroethene | 0.000 | | 0 | N.D. | | | 0.00 |
| 30) Hexane | 11.459 | 86 | 2803 | 0.74 | ppbv | # | 76 |
| 31) Chloroform | 12.511 | 83 | 806 | N.D. | | | |
| 32) EthylAcetate | 0.000 | | 0 | N.D. | | | d |

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201309.D
 Acq On : 20 May 2013 14:56
 Operator : JJG
 Sample : 130597-63051 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: May 20 18:44:03 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

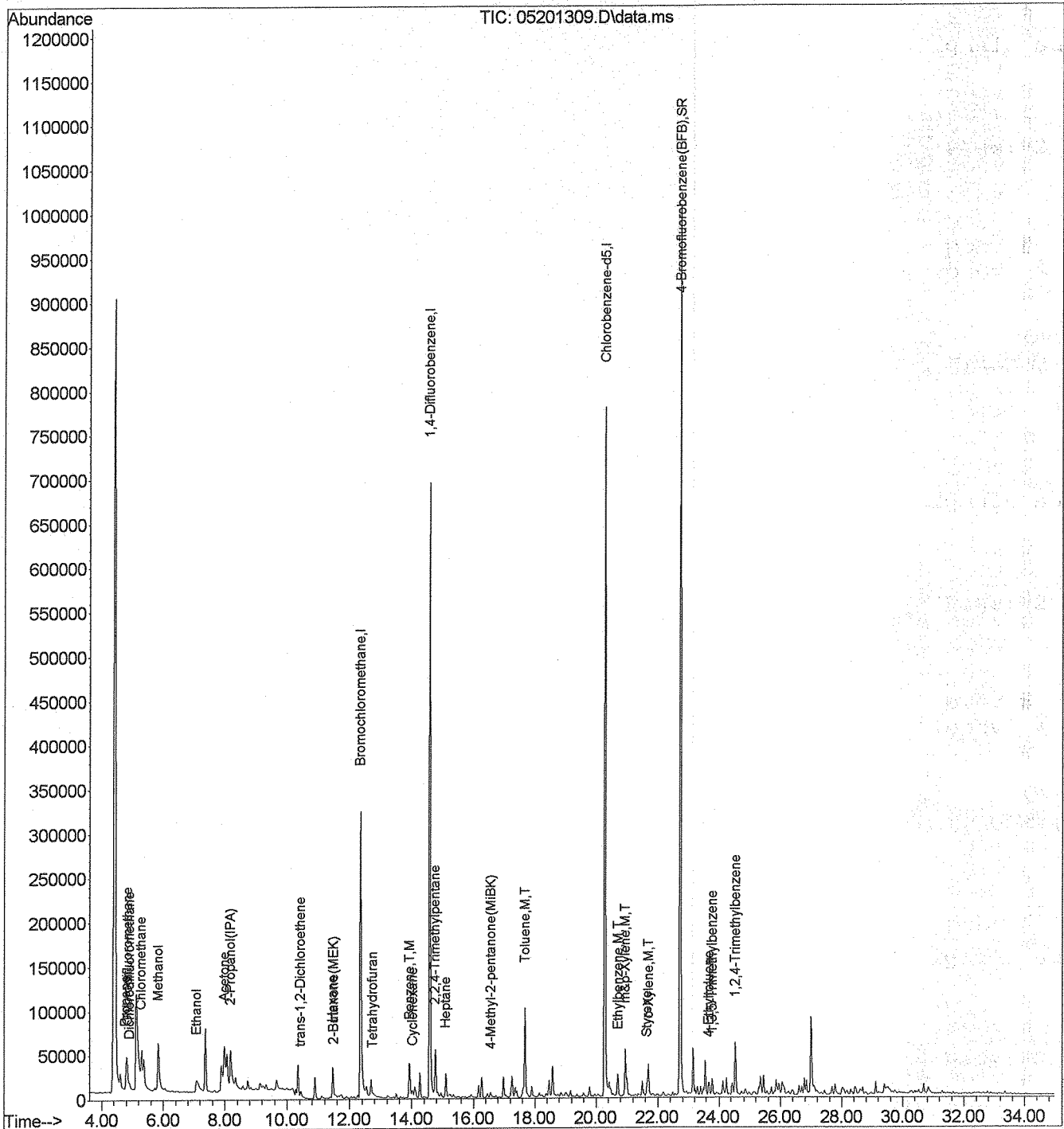
| Compound | R.T. | QIon | Response | Conc | Units | Dev (Min) |
|---------------------------------|--------|------|----------|------|--------|-----------|
| 33) Tetrahydrofuran | 12.743 | 72 | 1461 | 0.18 | ppbv # | 51 |
| 34) 1,2-Dichloroethane | 0.000 | | 0 | N.D. | | |
| 35) 1,1,1-Trichloroethane | 0.000 | | 0 | N.D. | | |
| 37) Benzene | 13.937 | 78 | 30241 | 0.53 | ppbv | 95 |
| 38) CarbonTetrachloride | 13.973 | 117 | 662 | N.D. | | |
| 39) Cyclohexane | 14.027 | 69 | 1082 | 0.13 | ppbv # | 63 |
| 40) 1,2-Dichloropropane | 0.000 | | 0 | N.D. | | |
| 41) Bromodichloromethane | 0.000 | | 0 | N.D. | | |
| 42) 1,4-Dioxane | 0.000 | | 0 | N.D. | | |
| 43) Trichloroethene (TCE) | 15.310 | 130 | 236 | N.D. | | |
| 44) 2,2,4-Trimethylpentane | 14.758 | 57 | 67357 | 0.66 | ppbv # | 93 |
| 45) Heptane | 15.096 | 71 | 7709 | 0.42 | ppbv # | 86 |
| 46) cis-1,3-Dichloropropene | 0.000 | | 0 | N.D. | | |
| 47) 4-Methyl-2-pentanone (M...) | 16.576 | 58 | 844 | 0.04 | ppbv # | 61 |
| 48) trans-1,3-Dichloropropene | 17.682 | 75 | 910 | N.D. | | |
| 49) 1,1,2-Trichloroethane | 17.860 | 97 | 354 | N.D. | | |
| 50) Toluene | 17.682 | 91 | 105241 | 1.45 | ppbv | |
| 51) 2-Hexanone (MBK) | 0.000 | | 0 | N.D. | | |
| 52) Dibromochloromethane | 19.019 | 129 | 703 | N.D. | | |
| 53) 1,2-Dibromoethane | 0.000 | | 0 | N.D. | | |
| 54) Tetrachloroethene (PCE) | 19.019 | 166 | 930 | N.D. | | |
| 56) Chlorobenzene | 20.285 | 114 | 135 | N.D. | | |
| 57) Ethylbenzene | 20.713 | 91 | 25085 | 0.26 | ppbv | 97 |
| 58) m&p-Xylene | 20.945 | 106 | 30471 | 0.80 | ppbv # | 92 |
| 59) Bromoform | 0.000 | | 0 | N.D. | | |
| 60) Styrene | 21.659 | 104 | 9505 | 0.15 | ppbv # | 94 |
| 61) 1,1,2,2-Tetrachloroethane | 22.318 | 83 | 108 | N.D. | | |
| 62) o-Xylene | 21.694 | 91 | 29268 | 0.38 | ppbv | 96 |
| 64) 4-Ethyltoluene | 23.691 | 120 | 4528 | 0.14 | ppbv # | 96 |
| 65) 1,3,5-Trimethylbenzene | 23.781 | 120 | 7763 | 0.17 | ppbv # | 89 |
| 66) 1,2,4-Trimethylbenzene | 24.529 | 120 | 24410 | 0.54 | ppbv | 91 |
| 67) BenzylChloride (a-Chlor...) | 25.189 | 91 | 140 | N.D. | | |
| 68) 1,3-Dichlorobenzene | 0.000 | | 0 | N.D. | | |
| 69) 1,4-Dichlorobenzene | 25.296 | 146 | 184 | N.D. | | |
| 70) 1,2-Dichlorobenzene | 0.000 | | 0 | N.D. | | |
| 71) 1,2,4-Trichlorobenzene | 29.469 | 180 | 338 | N.D. | | |
| 72) Hexachlorobutadiene | 0.000 | | 0 | N.D. | | |

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

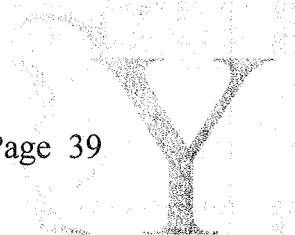
Handwritten signature
 96

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201309.D
 Acq On : 20 May 2013 14:56
 Operator : JJG
 Sample : 130597-63051 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: May 20 18:44:03 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration



**TO-15
RAW QC
& ICAL
SUMMARY**



MS #3 Instrument Logbook

8/20/13

Sequence Name: C:\msdchem\1\sequence\2013\052013.S

Comment: GCMS-03

Operator: JJG

Data Path: C:\MSDCHEM\1\MS03\2013\052013\

Instrument Control Pre-Seq Cmd:

Data Analysis Pre-Seq Cmd:

Instrument Control Post-Seq Cmd:

Data Analysis Post-Seq Cmd:

| Line | Sample Name/Misc Info |
|------------|--|
| 1) Sample | 1 05201301 TO15-5MS TO15 BFB 052013 |
| 2) Sample | 1 05201302 TO15-5MS TO15 CCV 052013 |
| 3) Sample | 1 05201303 TO15-5MS TO15 LCSD 052013 |
| 4) Sample | 1 05201304 TO15-5MS TO15 MB 052013 |
| 5) Sample | 2 05201305 TO15-5MS 130597-63048 x1 |
| 6) Sample | 2 05201306 TO15-5MS 130597-63048 x1 dp |
| 7) Sample | 3 05201307 TO15-5MS 130597-63049 x1 |
| 8) Sample | 4 05201308 TO15-5MS 130597-63050 x1 |
| 9) Sample | 5 05201309 TO15-5MS 130597-63051 x1 |
| 10) Sample | 6 05201310 TO15-5MS Flow Check#051713-01 |
| 11) Sample | 16 05201311 TO15-5MS TO15 0.25ppbv MDL#1 |
| 12) Sample | 16 05201312 TO15-5MS TO15 0.25ppbv MDL#2 |
| 13) Sample | 16 05201313 TO15-5MS TO15 0.25ppbv MDL#3 |
| 14) Sample | 16 05201314 TO15-5MS TO15 0.10ppbv MDL#1 |
| 15) Sample | 16 05201315 TO15-5MS TO15 0.10ppbv MDL#2 |
| 16) Sample | 16 05201316 TO15-5MS TO15 0.10ppbv MDL#3 |
| 17) Sample | 16 05201317 TO15-5MS TO15 0.05ppbv MDL#1 |
| 18) Sample | 16 05201318 TO15-5MS TO15 0.05ppbv MDL#2 |
| 19) Sample | 16 05201319 TO15-5MS TO15 0.05ppbv MDL#3 |
| 20) Sample | 16 05201320 TO15-5MS TO15 0.05ppbv MDL#4 |
| 21) Sample | 16 05201321 TO15-5MS TO15 0.05ppbv MDL#5 |
| 22) Sample | 16 05201322 TO15-5MS TO15 0.05ppbv MDL#6 |
| 23) Sample | 16 05201323 TO15-5MS TO15 0.05ppbv MDL#7 |
| 24) Sample | 1 05201324 TO15-5MS TO15 0.50ppbv MDL#1 |
| 25) Sample | 1 05201325 TO15-5MS TO15 0.50ppbv MDL#2 |
| 26) Sample | 1 05201326 TO15-5MS TO15 0.50ppbv MDL#3 |
| 27) Sample | 1 05201327 TO15-5MS TO15 0.50ppbv MDL#4 |
| 28) Sample | 1 05201328 TO15-5MS TO15 2.00ppbv MDL#1 |
| 29) Sample | 1 05201329 TO15-5MS TO15 2.00ppbv MDL#2 |
| 30) Sample | 1 05201330 TO15-5MS TO15 2.00ppbv MDL#3 |
| 31) Sample | 1 05201331 TO15-5MS Can Check#000663 |

8/20/13

Comments: _____

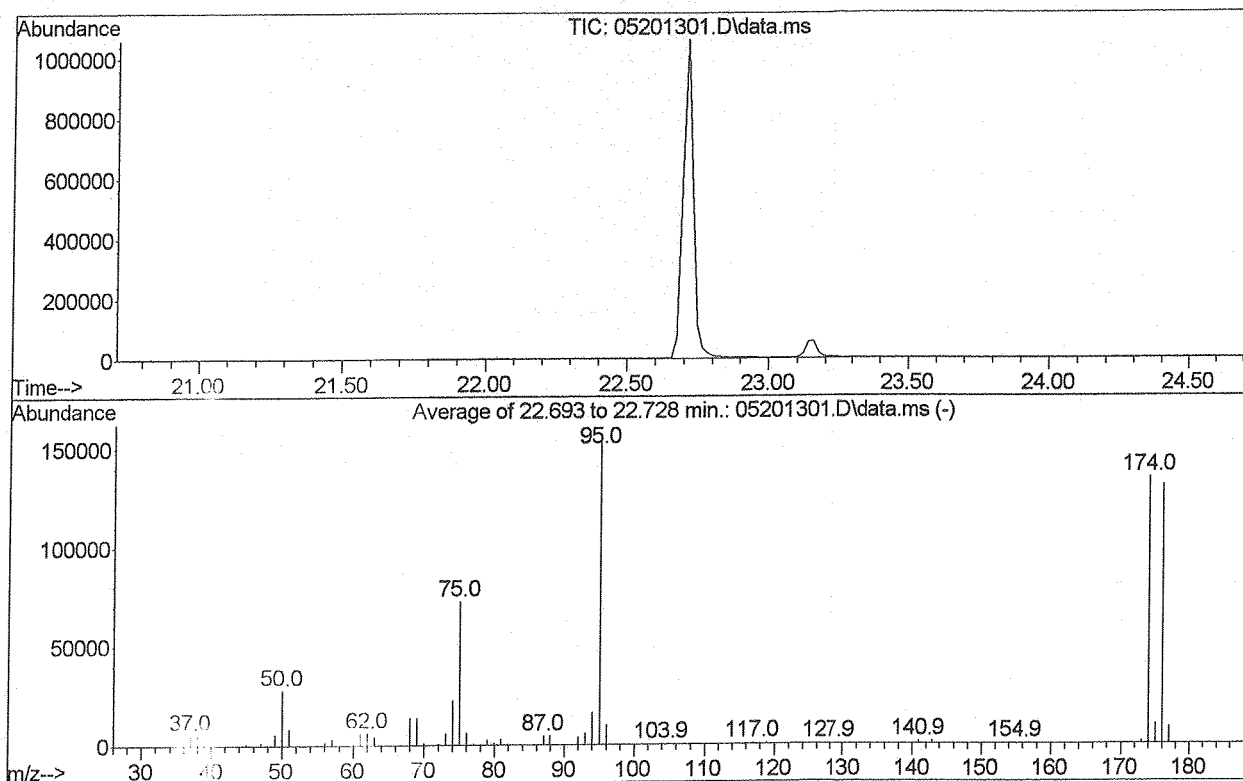
Analyst: _____

Date: _____

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201301.D
 Acq On : 20 May 2013 8:30 am
 Operator : JJG
 Sample : TO15 BFB 052013
 Misc : IS/Surr: PS082712-02 + 50mL
 ALS Vial : 1 Sample Multiplier: 1

Integration File: PAMS.P

Method : C:\msdchem\1\METHODS\2013\051513.M
 Title : TO-15/TO-14
 Last Update : Thu May 16 10:13:39 2013



AutoFind: Scans 1064, 1065, 1066; Background Corrected with Scan 1059

| Target Mass | Rel. to Mass | Lower Limit% | Upper Limit% | Rel. Abn% | Raw Abn | Result Pass/Fail |
|-------------|--------------|--------------|--------------|-----------|---------|------------------|
| 50 | 95 | 15 | 40 | 18.2 | 28152 | PASS |
| 75 | 95 | 30 | 60 | 46.9 | 72677 | PASS |
| 95 | 95 | 100 | 100 | 100.0 | 155029 | PASS |
| 96 | 95 | 5 | 9 | 6.7 | 10315 | PASS |
| 173 | 174 | 0.00 | 2 | 0.9 | 1250 | PASS |
| 174 | 95 | 50 | 100 | 86.9 | 134797 | PASS |
| 175 | 174 | 5 | 9 | 7.5 | 10114 | PASS |
| 176 | 174 | 95 | 101 | 97.3 | 131200 | PASS |
| 177 | 176 | 5 | 9 | 6.5 | 8588 | PASS |

[Handwritten signature]
 052013

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201302.D
 Acq On : 20 May 2013 9:16
 Operator : JJG
 Sample : TO15 CCV 052013
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 20 15:10:48 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev (Min) |
|--------------------------------|--------|------|----------|-------|-------|-----------|
| Internal Standards | | | | | | |
| 1) Bromochloromethane | 12.350 | 128 | 143576 | 10.00 | ppbv | 0.00 |
| 36) 1,4-Difluorobenzene | 14.579 | 114 | 771909 | 10.00 | ppbv | 0.00 |
| 55) Chlorobenzene-d5 | 20.285 | 117 | 728376 | 10.00 | ppbv | 0.00 |
| System Monitoring Compounds | | | | | | |
| 63) 4-Bromofluorobenzene (BFB) | 22.710 | 174 | 464060 | 10.18 | ppbv | 0.00 |

Spiked Amount 10.000 Recovery = 101.80%

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|---------------------------------|--------|------|----------|-------|-------|--------|
| 2) Chlorodifluoromethane | 4.818 | 51 | 291794 | 10.26 | ppbv | 98 |
| 3) Propene | 4.781 | 42 | 81603m | 10.92 | ppbv | |
| 4) Dichlorodifluoromethane | 4.908 | 85 | 494692 | 10.08 | ppbv | 99 |
| 5) Chloromethane | 5.288 | 52 | 46353m | 10.06 | ppbv | |
| 6) Dichlorotetrafluoroethane | 5.324 | 135 | 338768 | 10.13 | ppbv | 87 |
| 7) VinylChloride | 5.668 | 62 | 165430m | 10.13 | ppbv | |
| 8) Methanol | 5.867 | 31 | 24437m | 5.27 | ppbv | |
| 9) 1,3-Butadiene | 5.867 | 54 | 108659m | 10.48 | ppbv | |
| 10) Bromomethane | 6.446 | 96 | 107576m | 9.15 | ppbv | 98 |
| 11) Chloroethane | 6.736 | 66 | 25524 | 9.61 | ppbv | 98 |
| 12) Dichlorofluoromethane | 7.007 | 67 | 361081 | 10.28 | ppbv | 98 |
| 13) Ethanol | 7.043 | 45 | 63310m | 10.34 | ppbv | |
| 14) VinylBromide | 7.261 | 108 | 145824m | 10.56 | ppbv | |
| 15) Acetone | 7.966 | 58 | 75676m | 9.73 | ppbv | 99 |
| 16) Trichlorofluoromethane | 7.677 | 103 | 320744 | 11.18 | ppbv | 99 |
| 17) 2-Propanol (IPA) | 8.165 | 45 | 278535m | 10.46 | ppbv | |
| 18) Acrylonitrile | 8.962 | 52 | 128929m | 11.08 | ppbv | |
| 19) 1,1-Dichloroethene | 8.726 | 96 | 167885 | 10.27 | ppbv | 95 |
| 20) MethyleneChloride (DCM) | 9.324 | 84 | 149347m | 9.95 | ppbv | 98 |
| 21) AllylChloride | 9.305 | 39 | 152903m | 11.62 | ppbv | |
| 22) CarbonDisulfide | 9.486 | 76 | 482461m | 9.94 | ppbv | 98 |
| 23) Trichlorotrifluoroethane | 8.998 | 103 | 244998 | 10.41 | ppbv | 98 |
| 24) trans-1,2-Dichloroethene | 10.424 | 96 | 188989m | 10.75 | ppbv | 87 |
| 25) 1,1-Dichloroethane | 10.906 | 63 | 375435 | 10.47 | ppbv | 100 |
| 26) MethylTertButylether (M...) | 10.442 | 73 | 532942 | 11.19 | ppbv | 98 |
| 27) VinylAcetate | 10.888 | 43 | 453781m | 10.24 | ppbv | |
| 28) 2-Butanone (MEK) | 11.423 | 72 | 87526 | 11.00 | ppbv | 93 |
| 29) cis-1,2-Dichloroethene | 11.904 | 96 | 198379 | 10.48 | ppbv | 98 |
| 30) Hexane | 11.476 | 86 | 41018 | 10.81 | ppbv | 89 |
| 31) Chloroform | 12.511 | 83 | 438727 | 11.01 | ppbv | 98 |
| 32) EthylAcetate | 12.011 | 43 | 461179 | 11.51 | ppbv | 96 |

[Handwritten signature]

Data Path : C:\msdchem\1\MS03\2013\052013\
Data File : 05201302.D
Acq On : 20 May 2013 9:16
Operator : JJG
Sample : TO15 CCV 052013
Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
ALS Vial : 1 Sample Multiplier: 1

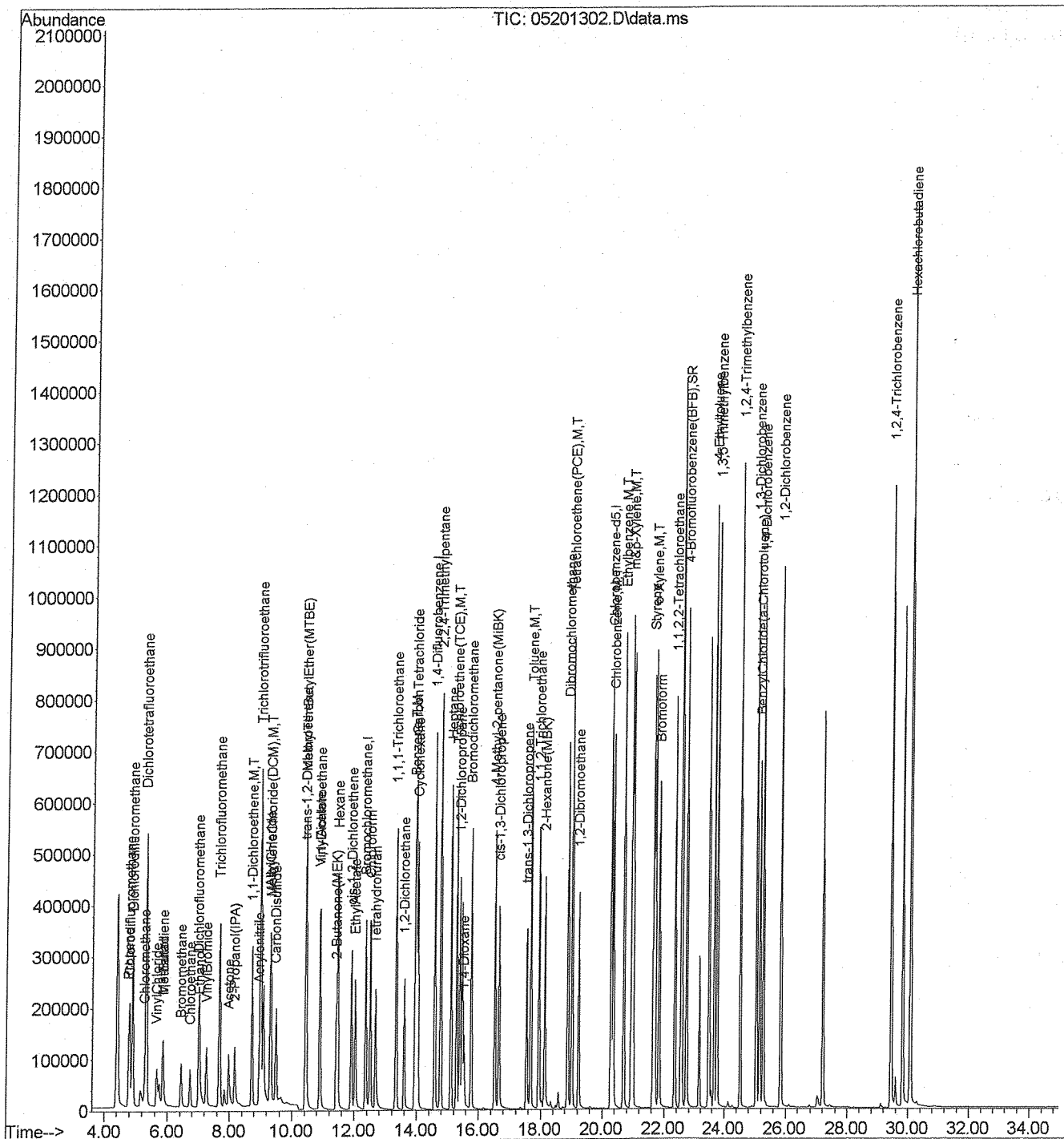
Quant Time: May 20 15:10:48 2013
Quant Method : C:\msdchem\1\METHODS\2013\051513.M
Quant Title : TO-15/TO-14
QLast Update : Thu May 16 10:13:39 2013
Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------------|--------|------|----------|-------|--------|----------|
| 33) Tetrahydrofuran | 12.671 | 72 | 86339 | 10.77 | ppbv | 91 |
| 34) 1,2-Dichloroethane | 13.598 | 62 | 324164 | 11.21 | ppbv | 98 |
| 35) 1,1,1-Trichloroethane | 13.331 | 97 | 499351 | 11.32 | ppbv | 99 |
| 37) Benzene | 13.937 | 78 | 547051 | 9.62 | ppbv | 98 |
| 38) CarbonTetrachloride | 13.973 | 117 | 494267 | 10.51 | ppbv | 99 |
| 39) Cyclohexane | 14.026 | 69 | 83769 | 9.94 | ppbv | 93 |
| 40) 1,2-Dichloropropane | 15.399 | 63 | 226882 | 9.92 | ppbv | 96 |
| 41) Bromodichloromethane | 15.756 | 85 | 318819 | 10.87 | ppbv | 98 |
| 42) 1,4-Dioxane | 15.524 | 88 | 1329020 | 9.91 | ppbv | |
| 43) Trichloroethene (TCE) | 15.292 | 130 | 281501 | 10.22 | ppbv | 99 |
| 44) 2,2,4-Trimethylpentane | 14.775 | 57 | 1069045 | 10.56 | ppbv | 98 |
| 45) Heptane | 15.114 | 71 | 185709 | 10.28 | ppbv | 98 |
| 46) cis-1,3-Dichloropropene | 16.648 | 75 | 353115 | 10.84 | ppbv | 97 |
| 47) 4-Methyl-2-pentanone (M...) | 16.523 | 58 | 215485 | 10.39 | ppbv | 96 |
| 48) trans-1,3-Dichloropropene | 17.539 | 75 | 327805 | 9.88 | ppbv | 99 |
| 49) 1,1,2-Trichloroethane | 17.931 | 97 | 260422 | 10.26 | ppbv | 97 |
| 50) Toluene | 17.682 | 91 | 739649 | 10.25 | ppbv | 100 |
| 51) 2-Hexanone (MBK) | 18.128 | 58 | 271737 | 10.57 | ppbv | 94 |
| 52) Dibromochloromethane | 18.877 | 129 | 561094 | 11.51 | ppbv | 99 |
| 53) 1,2-Dibromoethane | 19.233 | 107 | 428469 | 10.30 | ppbv | 100 |
| 54) Tetrachloroethene (PCE) | 19.019 | 166 | 412614 | 10.21 | ppbv | 98 |
| 56) Chlorobenzene | 20.357 | 114 | 197081 | 10.41 | ppbv | 98 |
| 57) Ethylbenzene | 20.695 | 91 | 981614 | 10.21 | ppbv | 99 |
| 58) m&p-Xylene | 20.945 | 106 | 740845 | 19.59 | ppbv # | 93 |
| 59) Bromoform | 21.837 | 173 | 529166 | 10.43 | ppbv | 99 |
| 60) Styrene | 21.640 | 104 | 604919 | 9.89 | ppbv | 98 |
| 61) 1,1,2,2-Tetrachloroethane | 22.336 | 83 | 559748 | 10.21 | ppbv | 98 |
| 62) o-Xylene | 21.694 | 91 | 764291 | 9.94 | ppbv | 99 |
| 64) 4-Ethyltoluene | 23.673 | 120 | 328644 | 10.41 | ppbv | 98 |
| 65) 1,3,5-Trimethylbenzene | 23.780 | 120 | 442338 | 9.59 | ppbv | 97 |
| 66) 1,2,4-Trimethylbenzene | 24.529 | 120 | 475453 | 10.58 | ppbv | 98 |
| 67) BenzylChloride (a-Chlor...) | 25.153 | 91 | 753014 | 11.13 | ppbv | 99 |
| 68) 1,3-Dichlorobenzene | 25.046 | 146 | 748687 | 10.59 | ppbv | 99 |
| 69) 1,4-Dichlorobenzene | 25.260 | 146 | 708025m | 9.95 | ppbv | 97 |
| 70) 1,2-Dichlorobenzene | 25.831 | 146 | 756953m | 10.24 | ppbv | 90 |
| 71) 1,2,4-Trichlorobenzene | 29.433 | 180 | 702250m | 9.60 | ppbv | 94 |
| 72) Hexachlorobutadiene | 30.075 | 225 | 599913m | 10.13 | ppbv | 94 |

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

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201302.D
 Acq On : 20 May 2013 9:16
 Operator : JJG
 Sample : TO15 CCV 052013
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 20 15:10:48 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration



Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201303.D
 Acq On : 20 May 2013 10:08
 Operator : JJG
 Sample : TO15 LCSD 052013
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 20 15:13:30 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------|--------|------|----------|-------|-------|----------|
| Internal Standards | | | | | | |
| 1) Bromochloromethane | 12.350 | 128 | 152556 | 10.00 | ppbv | 0.00 |
| 36) 1,4-Difluorobenzene | 14.579 | 114 | 794135 | 10.00 | ppbv | 0.00 |
| 55) Chlorobenzene-d5 | 20.285 | 117 | 761722 | 10.00 | ppbv | 0.00 |

| | | | | | | |
|--------------------------------|--------|-----|--------|-------|------|------|
| System Monitoring Compounds | | | | | | |
| 63) 4-Bromofluorobenzene (BFB) | 22.711 | 174 | 476998 | 10.01 | ppbv | 0.00 |

Spiked Amount 10.000 Recovery = 100.10%

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|---------------------------------|--------|------|----------|-------|--------|------------|
| 2) Chlorodifluoromethane | 4.817 | 51 | 303974 | 10.06 | ppbv | 100 |
| 3) Propene | 4.781 | 42 | 87596 | 11.04 | ppbv | 99 |
| 4) Dichlorodifluoromethane | 4.908 | 85 | 520885m | 9.99 | ppbv | |
| 5) Chloromethane | 5.288 | 52 | 51259m | 10.47 | ppbv | |
| 6) Dichlorotetrafluoroethane | 5.342 | 135 | 368027 | 10.35 | ppbv | 99 |
| 7) VinylChloride | 5.668 | 62 | 182843m | 10.54 | ppbv | Qvalue 100 |
| 8) Methanol | 5.867 | 31 | 24827m | 5.04 | ppbv | |
| 9) 1,3-Butadiene | 5.867 | 54 | 116950m | 10.62 | ppbv | |
| 10) Bromomethane | 6.446 | 96 | 124047m | 9.93 | ppbv | 0.00 |
| 11) Chloroethane | 6.736 | 66 | 28238 | 10.00 | ppbv | 0.99 |
| 12) Dichlorofluoromethane | 7.025 | 67 | 379194 | 10.16 | ppbv | 100 |
| 13) Ethanol | 7.061 | 45 | 66127m | 10.16 | ppbv | |
| 14) VinylBromide | 7.261 | 108 | 157402m | 10.73 | ppbv | |
| 15) Acetone | 7.966 | 58 | 77640m | 9.40 | ppbv | 0.00 |
| 16) Trichlorofluoromethane | 7.677 | 103 | 321049 | 10.54 | ppbv | 98 |
| 17) 2-Propanol (IPA) | 8.165 | 45 | 284498m | 10.06 | ppbv | 100 |
| 18) Acrylonitrile | 8.962 | 52 | 133005m | 10.75 | ppbv | |
| 19) 1,1-Dichloroethene | 8.726 | 96 | 172274 | 9.91 | ppbv | Qvalue 94 |
| 20) MethyleneChloride (DCM) | 9.323 | 84 | 153136m | 9.60 | ppbv | 100 |
| 21) AllylChloride | 9.305 | 39 | 150889m | 10.79 | ppbv | 99 |
| 22) CarbonDisulfide | 9.486 | 76 | 497311m | 9.64 | ppbv | |
| 23) Trichlorotrifluoroethane | 8.998 | 103 | 254299 | 10.17 | ppbv | 99 |
| 24) trans-1,2-Dichloroethene | 10.424 | 96 | 202904m | 10.86 | ppbv | 99 |
| 25) 1,1-Dichloroethane | 10.906 | 63 | 386178 | 10.13 | ppbv | 99 |
| 26) MethylTertButylEther (M...) | 10.460 | 73 | 547978 | 10.83 | ppbv | 99 |
| 27) VinylAcetate | 10.888 | 43 | 462337m | 9.82 | ppbv | |
| 28) 2-Butanone (MEK) | 11.423 | 72 | 89975 | 10.64 | ppbv | 89 |
| 29) cis-1,2-Dichloroethene | 11.905 | 96 | 210020 | 10.44 | ppbv | 98 |
| 30) Hexane | 11.477 | 86 | 41847 | 10.38 | ppbv | 98 |
| 31) Chloroform | 12.493 | 83 | 454880 | 10.74 | ppbv | 96 |
| 32) EthylAcetate | 12.011 | 43 | 479187 | 11.25 | ppbv # | 97 |

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201303.D
 Acq On : 20 May 2013 10:08
 Operator : JJG
 Sample : TO15 LCSD 052013
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

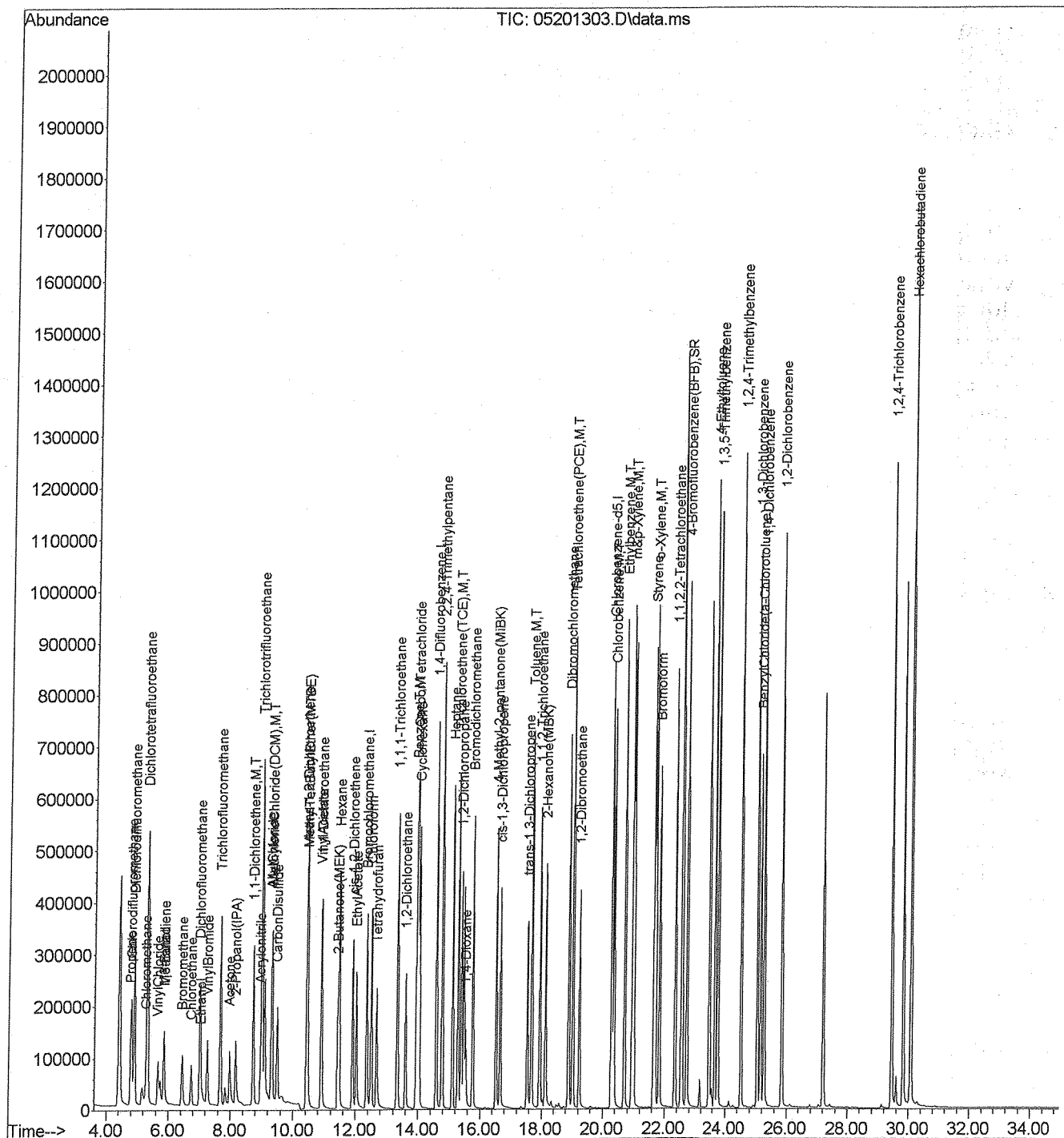
Quant Time: May 20 15:13:30 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------------|--------|------|----------|-------|--------|----------|
| 33) Tetrahydrofuran | 12.671 | 72 | 904420 | 10.62 | ppbv | |
| 34) 1,2-Dichloroethane | 13.599 | 62 | 335654 | 10.93 | ppbv | 98 |
| 35) 1,1,1-Trichloroethane | 13.331 | 97 | 511877 | 10.92 | ppbv | 100 |
| 37) Benzene | 13.937 | 78 | 581905 | 9.95 | ppbv | 99 |
| 38) CarbonTetrachloride | 13.973 | 117 | 512676 | 10.60 | ppbv | 98 |
| 39) Cyclohexane | 14.026 | 69 | 88291 | 10.18 | ppbv | 93 |
| 40) 1,2-Dichloropropane | 15.400 | 63 | 236548 | 10.05 | ppbv | 97 |
| 41) Bromodichloromethane | 15.756 | 85 | 320799 | 10.63 | ppbv | 100 |
| 42) 1,4-Dioxane | 15.524 | 88 | 1416830 | 10.27 | ppbv | |
| 43) Trichloroethene (TCE) | 15.293 | 130 | 288607 | 10.18 | ppbv | 98 |
| 44) 2,2,4-Trimethylpentane | 14.775 | 57 | 1124746 | 10.80 | ppbv | 99 |
| 45) Heptane | 15.114 | 71 | 194332 | 10.46 | ppbv | 95 |
| 46) cis-1,3-Dichloropropene | 16.648 | 75 | 372944 | 11.13 | ppbv | 98 |
| 47) 4-Methyl-2-pentanone (M...) | 16.523 | 58 | 220138 | 10.32 | ppbv | 98 |
| 48) trans-1,3-Dichloropropene | 17.539 | 75 | 346666 | 10.16 | ppbv | 99 |
| 49) 1,1,2-Trichloroethane | 17.932 | 97 | 269003 | 10.30 | ppbv | 98 |
| 50) Toluene | 17.682 | 91 | 756607 | 10.19 | ppbv | 100 |
| 51) 2-Hexanone (MBK) | 18.128 | 58 | 287314 | 10.87 | ppbv | 98 |
| 52) Dibromochloromethane | 18.877 | 129 | 572137 | 11.41 | ppbv | 99 |
| 53) 1,2-Dibromoethane | 19.233 | 107 | 439448 | 10.27 | ppbv | 98 |
| 54) Tetrachloroethene (PCE) | 19.019 | 166 | 423610 | 10.18 | ppbv | 98 |
| 56) Chlorobenzene | 20.357 | 114 | 205021 | 10.35 | ppbv | 99 |
| 57) Ethylbenzene | 20.696 | 91 | 1008992 | 10.04 | ppbv | 100 |
| 58) m&p-Xylene | 20.945 | 106 | 759479 | 19.20 | ppbv | 95 |
| 59) Bromoform | 21.819 | 173 | 546426 | 10.29 | ppbv # | 96 |
| 60) Styrene | 21.641 | 104 | 651232 | 10.18 | ppbv | 99 |
| 61) 1,1,2,2-Tetrachloroethane | 22.336 | 83 | 601822 | 10.50 | ppbv | 99 |
| 62) o-Xylene | 21.694 | 91 | 808676 | 10.06 | ppbv | 100 |
| 64) 4-Ethyltoluene | 23.674 | 120 | 346452 | 10.49 | ppbv | 99 |
| 65) 1,3,5-Trimethylbenzene | 23.781 | 120 | 460933 | 9.55 | ppbv | 98 |
| 66) 1,2,4-Trimethylbenzene | 24.529 | 120 | 489304 | 10.41 | ppbv | 99 |
| 67) BenzylChloride (a-Chlor...) | 25.154 | 91 | 779146 | 11.02 | ppbv | 99 |
| 68) 1,3-Dichlorobenzene | 25.047 | 146 | 764676 | 10.34 | ppbv | 99 |
| 69) 1,4-Dichlorobenzene | 25.261 | 146 | 728879m | 9.80 | ppbv | 98 |
| 70) 1,2-Dichlorobenzene | 25.831 | 146 | 767808m | 9.93 | ppbv | 99 |
| 71) 1,2,4-Trichlorobenzene | 29.433 | 180 | 729797m | 9.54 | ppbv | 99 |
| 72) Hexachlorobutadiene | 30.075 | 225 | 603707m | 9.75 | ppbv | 99 |

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

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201303.D
 Acq On : 20 May 2013 10:08
 Operator : JJG
 Sample : TO15 LCSD 052013
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 20 15:13:30 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration



Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201304.D
 Acq On : 20 May 2013 10:56
 Operator : JJG
 Sample : TO15 MB 052013
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 20 15:16:45 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------------------|--------|------|----------|-------|-------|----------|
| Internal Standards | | | | | | |
| 1) Bromochloromethane | 12.350 | 128 | 147959 | 10.00 | ppbv | 0.00 |
| 36) 1,4-Difluorobenzene | 14.579 | 114 | 803870 | 10.00 | ppbv | 0.00 |
| 55) Chlorobenzene-d5 | 20.285 | 117 | 747823 | 10.00 | ppbv | 0.00 |
| System Monitoring Compounds | | | | | | |
| 63) 4-Bromofluorobenzene (BFB) | 22.710 | 174 | 477532 | 10.20 | ppbv | 0.00 |

Spiked Amount 10.000 Recovery = 102.00%

| Target Compounds | R.T. | QIon | Response | Conc | Units | Dev(Min) | Qvalue |
|---------------------------------|-------|------|----------|------|-------|----------|--------|
| 2) Chlorodifluoromethane | 0.000 | | 0 | | N.D. | | |
| 3) Propene | 4.835 | 42 | 245 | | N.D. | | |
| 4) Dichlorodifluoromethane | 0.000 | | 0 | | N.D. | | |
| 5) Chloromethane | 0.000 | | 0 | | N.D. | | |
| 6) Dichlorotetrafluoroethane | 0.000 | | 0 | | N.D. | | |
| 7) VinylChloride | 0.000 | | 0 | | N.D. | | |
| 8) Methanol | 0.000 | | 0 | | N.D. | | |
| 9) 1,3-Butadiene | 0.000 | | 0 | | N.D. | | |
| 10) Bromomethane | 0.000 | | 0 | | N.D. | 0.00 | |
| 11) Chloroethane | 0.000 | | 0 | | N.D. | 0.00 | |
| 12) Dichlorofluoromethane | 0.000 | | 0 | | N.D. | 0.00 | |
| 13) Ethanol | 0.000 | | 0 | | N.D. | | |
| 14) VinylBromide | 0.000 | | 0 | | N.D. | | |
| 15) Acetone | 0.000 | | 0 | | N.D. | 0.00 | |
| 16) Trichlorofluoromethane | 0.000 | | 0 | | N.D. | | |
| 17) 2-Propanol (IPA) | 8.310 | 45 | 357 | | N.D. | 100.00% | |
| 18) Acrylonitrile | 0.000 | | 0 | | N.D. | | |
| 19) 1,1-Dichloroethene | 0.000 | | 0 | | N.D. | | Qvalue |
| 20) MethyleneChloride (DCM) | 0.000 | | 0 | | N.D. | d | |
| 21) AllylChloride | 0.000 | | 0 | | N.D. | | |
| 22) CarbonDisulfide | 9.522 | 76 | 921 | | N.D. | | |
| 23) Trichlorotrifluoroethane | 0.000 | | 0 | | N.D. | | |
| 24) trans-1,2-Dichloroethene | 0.000 | | 0 | | N.D. | | |
| 25) 1,1-Dichloroethane | 0.000 | | 0 | | N.D. | Dev(Min) | |
| 26) MethylTertButylether (M...) | 0.000 | | 0 | | N.D. | d | |
| 27) VinylAcetate | 0.000 | | 0 | | N.D. | | |
| 28) 2-Butanone (MEK) | 0.000 | | 0 | | N.D. | 0.00 | |
| 29) cis-1,2-Dichloroethene | 0.000 | | 0 | | N.D. | 0.00 | |
| 30) Hexane | 0.000 | | 0 | | N.D. | 0.00 | |
| 31) Chloroform | 0.000 | | 0 | | N.D. | | |
| 32) EthylAcetate | 0.000 | | 0 | | N.D. | | |

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201304.D
 Acq On : 20 May 2013 10:56
 Operator : JJG
 Sample : TO15 MB 052013
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 20 15:16:45 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

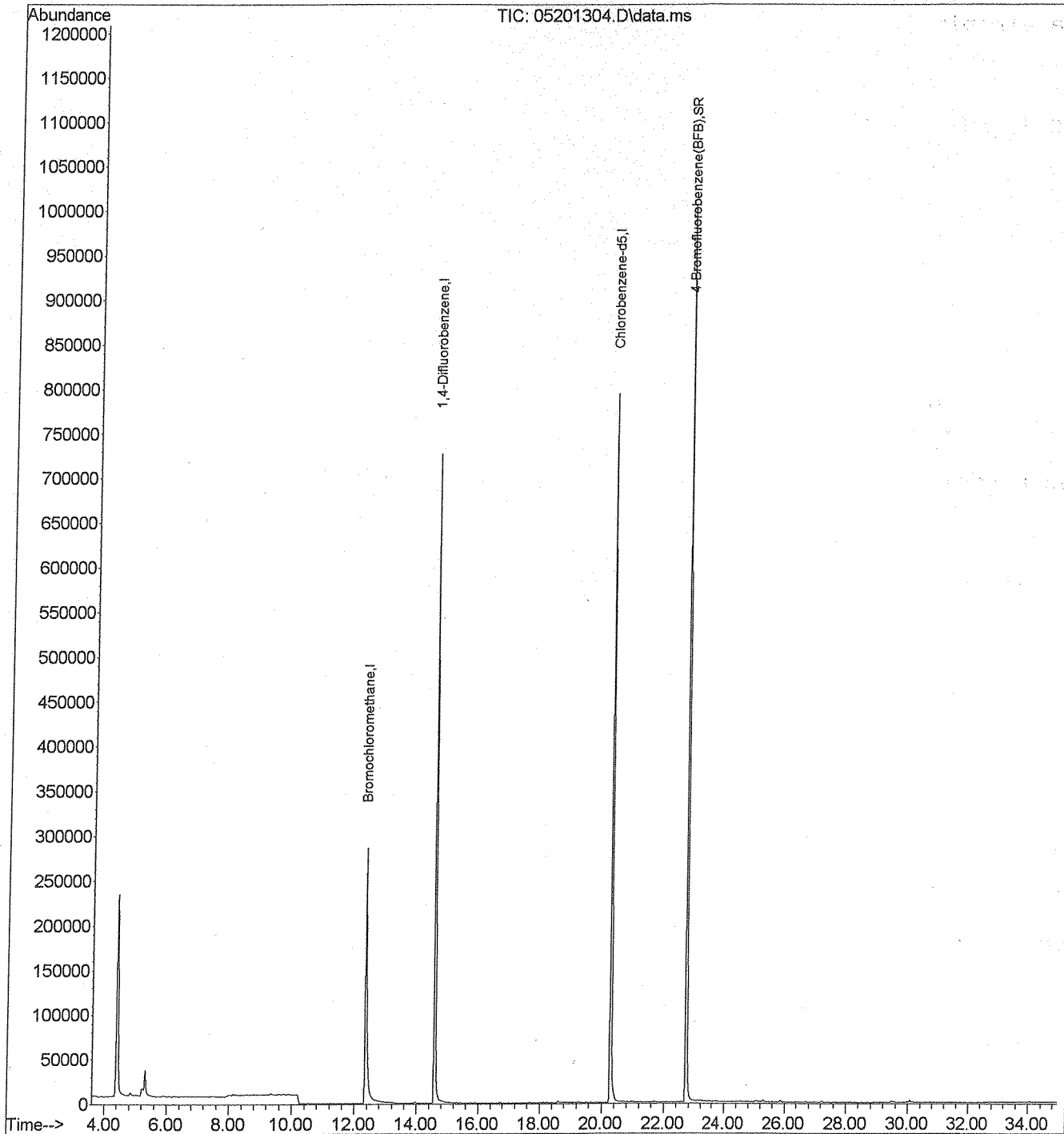
| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------------------|--------|------|----------|------|-------|----------|
| 33) Tetrahydrofuran | 0.000 | | 0 | | N.D. | |
| 34) 1,2-Dichloroethane | 0.000 | | 0 | | N.D. | |
| 35) 1,1,1-Trichloroethane | 0.000 | | 0 | | N.D. | |
| 37) Benzene | 13.955 | 78 | 1768 | | N.D. | |
| 38) CarbonTetrachloride | 0.000 | | 0 | | N.D. | |
| 39) Cyclohexane | 0.000 | | 0 | | N.D. | |
| 40) 1,2-Dichloropropane | 0.000 | | 0 | | N.D. | |
| 41) Bromodichloromethane | 0.000 | | 0 | | N.D. | |
| 42) 1,4-Dioxane | 0.000 | | 0 | | N.D. | |
| 43) Trichloroethene (TCE) | 0.000 | | 0 | | N.D. | |
| 44) 2,2,4-Trimethylpentane | 0.000 | | 0 | | N.D. | |
| 45) Heptane | 0.000 | | 0 | | N.D. | |
| 46) cis-1,3-Dichloropropene | 0.000 | | 0 | | N.D. | |
| 47) 4-Methyl-2-pentanone (M... | 0.000 | | 0 | | N.D. | |
| 48) trans-1,3-Dichloropropene | 0.000 | | 0 | | N.D. | |
| 49) 1,1,2-Trichloroethane | 0.000 | | 0 | | N.D. | |
| 50) Toluene | 17.700 | 91 | 480 | | N.D. | |
| 51) 2-Hexanone (MBK) | 0.000 | | 0 | | N.D. | |
| 52) Dibromochloromethane | 0.000 | | 0 | | N.D. | |
| 53) 1,2-Dibromoethane | 0.000 | | 0 | | N.D. | |
| 54) Tetrachloroethene (PCE) | 0.000 | | 0 | | N.D. | |
| 56) Chlorobenzene | 20.357 | 114 | 288 | | N.D. | |
| 57) Ethylbenzene | 20.713 | 91 | 649 | | N.D. | |
| 58) m&p-Xylene | 21.034 | 106 | 245 | | N.D. | |
| 59) Bromoform | 0.000 | | 0 | | N.D. | |
| 60) Styrene | 21.694 | 104 | 478 | | N.D. | |
| 61) 1,1,2,2-Tetrachloroethane | 22.354 | 83 | 131 | | N.D. | |
| 62) o-Xylene | 21.712 | 91 | 483 | | N.D. | |
| 64) 4-Ethyltoluene | 23.709 | 120 | 123 | | N.D. | |
| 65) 1,3,5-Trimethylbenzene | 23.709 | 120 | 123 | | N.D. | |
| 66) 1,2,4-Trimethylbenzene | 24.565 | 120 | 281 | | N.D. | |
| 67) BenzylChloride (a-Chlor... | 25.225 | 91 | 888 | | N.D. | |
| 68) 1,3-Dichlorobenzene | 25.064 | 146 | 1801 | | N.D. | |
| 69) 1,4-Dichlorobenzene | 25.296 | 146 | 2498 | | N.D. | |
| 70) 1,2-Dichlorobenzene | 25.867 | 146 | 1709 | | N.D. | |
| 71) 1,2,4-Trichlorobenzene | 29.469 | 180 | 2820 | | N.D. | |
| 72) Hexachlorobutadiene | 30.075 | 225 | 1034 | | N.D. | |

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

Handwritten signature and date: 05/20/13

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201304.D
 Acq On : 20 May 2013 10:56
 Operator : JJG
 Sample : TO15 MB 052013
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 20 15:16:45 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration



Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201305.D
 Acq On : 20 May 2013 11:44
 Operator : JJG
 Sample : 130597-63048 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 20 15:18:58 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev (Min) |
|-------------------------|--------|------|----------|-------|-------|-----------|
| Internal Standards | | | | | | |
| 1) Bromochloromethane | 12.350 | 128 | 145590 | 10.00 | ppbv | 0.00 |
| 36) 1,4-Difluorobenzene | 14.579 | 114 | 790345 | 10.00 | ppbv | 0.00 |
| 55) Chlorobenzene-d5 | 20.285 | 117 | 750102 | 10.00 | ppbv | 0.00 |

| | | | | | | |
|--------------------------------|--------|-----|--------|-------|------|------|
| System Monitoring Compounds | | | | | | |
| 63) 4-Bromofluorobenzene (BFB) | 22.710 | 174 | 479494 | 10.21 | ppbv | 0.00 |

Spiked Amount 10.000 Recovery = 102.10%

| Target Compounds | R.T. | QIon | Response | Conc | Units | Dev (Min) | Qvalue |
|---------------------------------|--------|------|----------|------|--------|-----------|--------|
| 2) Chlorodifluoromethane | 4.818 | 51 | 4516 | 0.16 | ppbv # | | 96 |
| 3) Propene | 0.000 | | 0 | N.D. | d | | |
| 4) Dichlorodifluoromethane | 4.890 | 85 | 14589 | 0.29 | ppbv | | 98 |
| 5) Chloromethane | 5.288 | 52 | 975 | 0.21 | ppbv # | | 10 |
| 6) Dichlorotetrafluoroethane | 5.324 | 135 | 204 | N.D. | | | |
| 7) VinylChloride | 0.000 | | 0 | N.D. | Units | | |
| 8) Methanol | 5.867 | 31 | 39778 | 8.56 | ppbv | | |
| 9) 1,3-Butadiene | 0.000 | | 0 | N.D. | | | |
| 10) Bromomethane | 0.000 | | 0 | N.D. | ppbv | | 0.00 |
| 11) Chloroethane | 0.000 | | 0 | N.D. | ppbv | | 0.00 |
| 12) Dichlorofluoromethane | 0.000 | | 0 | N.D. | | | 0.00 |
| 13) Ethanol | 7.134 | 45 | 17153 | 2.76 | ppbv | | |
| 14) VinylBromide | 0.000 | | 0 | N.D. | | | |
| 15) Acetone | 7.984 | 58 | 32335 | 4.10 | ppbv | | 0.00 |
| 16) Trichlorofluoromethane | 7.659 | 103 | 4061 | 0.14 | ppbv # | | 95 |
| 17) 2-Propanol (IPA) | 8.238 | 45 | 6570 | 0.24 | ppbv | | 0.00 |
| 18) Acrylonitrile | 0.000 | | 0 | N.D. | | | |
| 19) 1,1-Dichloroethene | 0.000 | | 0 | N.D. | | | Qvalue |
| 20) MethyleneChloride (DCM) | 0.000 | | 0 | N.D. | d | | 96 |
| 21) AllylChloride | 0.000 | | 0 | N.D. | d | | |
| 22) CarbonDisulfide | 0.000 | | 0 | N.D. | d | | 98 |
| 23) Trichlorotrifluoroethane | 0.000 | | 0 | N.D. | d | | 10 |
| 24) trans-1,2-Dichloroethene | 0.000 | | 0 | N.D. | | | |
| 25) 1,1-Dichloroethane | 0.000 | | 0 | N.D. | | | |
| 26) MethylTertButylEther (M...) | 0.000 | | 0 | N.D. | ppbv | | |
| 27) VinylAcetate | 0.000 | | 0 | N.D. | d | | |
| 28) 2-Butanone (MEK) | 11.494 | 72 | 3961 | 0.49 | ppbv # | | 35 |
| 29) cis-1,2-Dichloroethene | 0.000 | | 0 | N.D. | | | 0.00 |
| 30) Hexane | 11.458 | 86 | 1972 | 0.51 | ppbv # | | 49 |
| 31) Chloroform | 12.493 | 83 | 995 | N.D. | | | |
| 32) EthylAcetate | 12.118 | 43 | 1448 | N.D. | | | |

Handwritten signature and date: JJG 5/20/13

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201305.D
 Acq On : 20 May 2013 11:44
 Operator : JJG
 Sample : 130597-63048 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 20 15:18:58 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

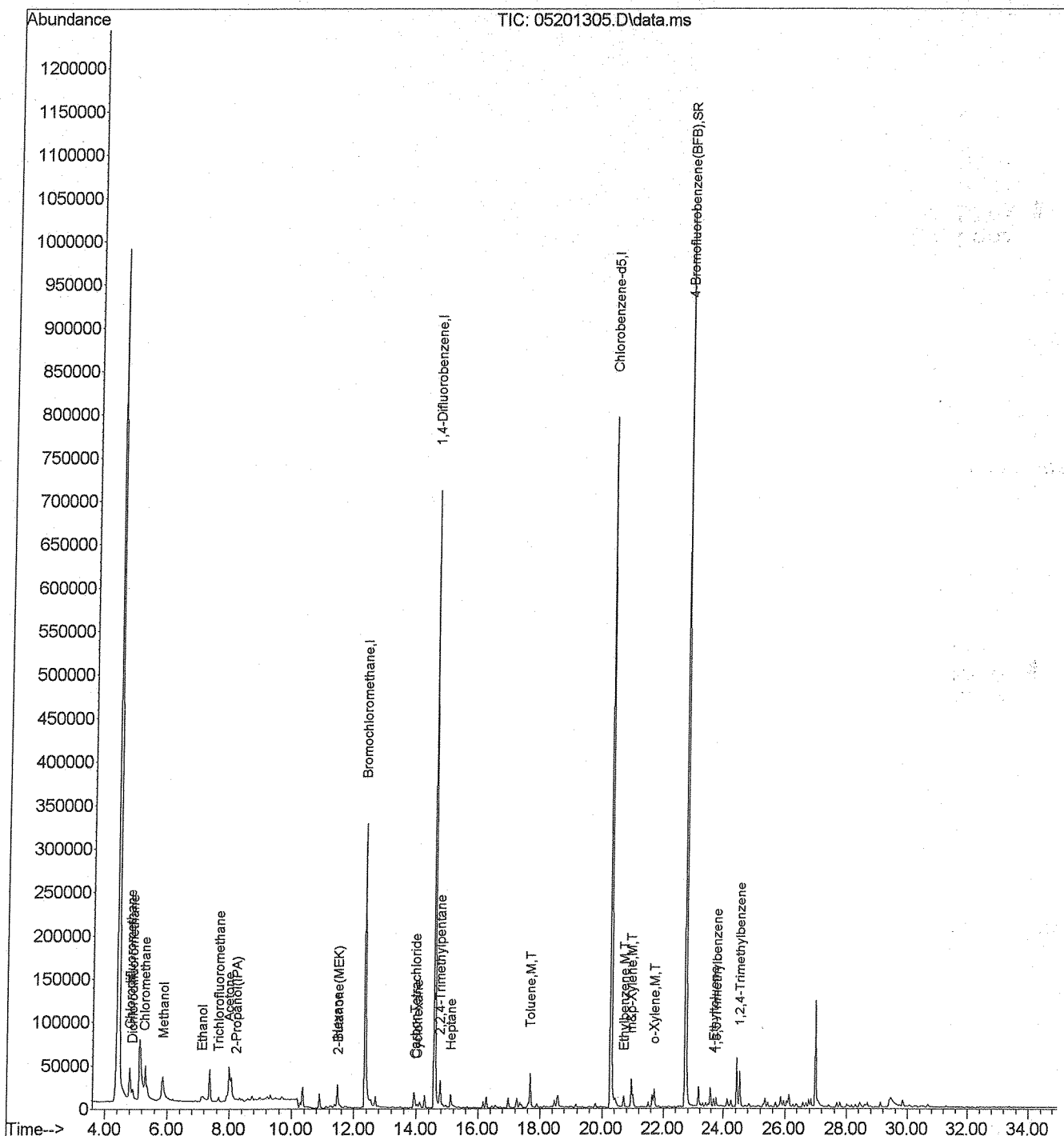
| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------------|--------|------|----------|------|--------|----------|
| 33) Tetrahydrofuran | 0.000 | | 0 | N.D. | | |
| 34) 1,2-Dichloroethane | 13.634 | 62 | 123 | N.D. | | |
| 35) 1,1,1-Trichloroethane | 0.000 | | 0 | N.D. | | |
| 37) Benzene | 0.000 | | 0 | N.D. | d | |
| 38) CarbonTetrachloride | 13.973 | 117 | 2480 | 0.05 | ppbv | 93 |
| 39) Cyclohexane | 14.026 | 69 | 612 | 0.07 | ppbv # | 92 |
| 40) 1,2-Dichloropropane | 15.257 | 63 | 116 | N.D. | | |
| 41) Bromodichloromethane | 0.000 | | 0 | N.D. | | |
| 42) 1,4-Dioxane | 0.000 | | 0 | N.D. | | |
| 43) Trichloroethene (TCE) | 15.292 | 130 | 259 | N.D. | | |
| 44) 2,2,4-Trimethylpentane | 14.757 | 57 | 39593 | 0.38 | ppbv | 96 |
| 45) Heptane | 15.114 | 71 | 4039 | 0.22 | ppbv # | 86 |
| 46) cis-1,3-Dichloropropene | 0.000 | | 0 | N.D. | | |
| 47) 4-Methyl-2-pentanone (M...) | 16.576 | 58 | 710 | N.D. | | |
| 48) trans-1,3-Dichloropropene | 17.682 | 75 | 154 | N.D. | | |
| 49) 1,1,2-Trichloroethane | 0.000 | | 0 | N.D. | | |
| 50) Toluene | 17.682 | 91 | 42772 | 0.58 | ppbv | 99 |
| 51) 2-Hexanone (MBK) | 0.000 | | 0 | N.D. | | |
| 52) Dibromochloromethane | 0.000 | | 0 | N.D. | | |
| 53) 1,2-Dibromoethane | 0.000 | | 0 | N.D. | | |
| 54) Tetrachloroethene (PCE) | 19.019 | 166 | 116 | N.D. | | |
| 56) Chlorobenzene | 0.000 | | 0 | N.D. | | |
| 57) Ethylbenzene | 20.713 | 91 | 14549 | 0.15 | ppbv | 98 |
| 58) m&p-Xylene | 20.945 | 106 | 19939 | 0.51 | ppbv # | 89 |
| 59) Bromoform | 0.000 | | 0 | N.D. | | |
| 60) Styrene | 21.694 | 104 | 1167 | N.D. | | |
| 61) 1,1,2,2-Tetrachloroethane | 22.354 | 83 | 122 | N.D. | | |
| 62) o-Xylene | 21.694 | 91 | 17217 | 0.22 | ppbv | 97 |
| 64) 4-Ethyltoluene | 23.691 | 120 | 3288 | 0.10 | ppbv | 95 |
| 65) 1,3,5-Trimethylbenzene | 23.780 | 120 | 4534 | 0.10 | ppbv # | 91 |
| 66) 1,2,4-Trimethylbenzene | 24.529 | 120 | 16063 | 0.35 | ppbv | 94 |
| 67) BenzylChloride (a-Chlor...) | 25.189 | 91 | 424 | N.D. | | |
| 68) 1,3-Dichlorobenzene | 25.064 | 146 | 678 | N.D. | | |
| 69) 1,4-Dichlorobenzene | 25.296 | 146 | 1203 | N.D. | | |
| 70) 1,2-Dichlorobenzene | 25.867 | 146 | 924 | N.D. | | |
| 71) 1,2,4-Trichlorobenzene | 29.451 | 180 | 2361 | N.D. | | |
| 72) Hexachlorobutadiene | 30.075 | 225 | 876 | N.D. | | |

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

Handwritten signature and date:
 05/20/13

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201305.D
 Acq On : 20 May 2013 11:44
 Operator : JJG
 Sample : 130597-63048 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 20 15:18:58 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration



Handwritten signature/initials
 05/20/13

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201306.D
 Acq On : 20 May 2013 12:32
 Operator : JJG
 Sample : 130597-63048 x1 dp
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 20 15:22:35 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev (Min) |
|-------------------------|--------|------|----------|-------|-------|-----------|
| Internal Standards | | | | | | |
| 1) Bromochloromethane | 12.350 | 128 | 142078 | 10.00 | ppbv | 0.00 |
| 36) 1,4-Difluorobenzene | 14.579 | 114 | 794580 | 10.00 | ppbv | 0.00 |
| 55) Chlorobenzene-d5 | 20.285 | 117 | 739144 | 10.00 | ppbv | 0.00 |

| | | | | | | |
|--------------------------------|--------|-----|--------|-------|------|------|
| System Monitoring Compounds | | | | | | |
| 63) 4-Bromofluorobenzene (BFB) | 22.710 | 174 | 482007 | 10.42 | ppbv | 0.00 |

Spiked Amount 10.000 Recovery = 104.20%

| Target Compounds | R.T. | QIon | Response | Conc | Units | Dev (Min) | Qvalue |
|---------------------------------|--------|------|----------|------|-------|-----------|--------|
| 2) Chlorodifluoromethane | 4.817 | 51 | 4288 | 0.15 | ppbv | | # 96 |
| 3) Propene | 0.000 | | 0 | N.D. | d | | |
| 4) Dichlorodifluoromethane | 4.890 | 85 | 14587 | 0.30 | ppbv | | # 97 |
| 5) Chloromethane | 5.288 | 52 | 1200 | 0.26 | ppbv | | # 9 |
| 6) Dichlorotetrafluoroethane | 5.306 | 135 | 115 | N.D. | | | |
| 7) VinylChloride | 0.000 | | 0 | N.D. | | | |
| 8) Methanol | 5.867 | 31 | 390750 | 8.62 | ppbv | | |
| 9) 1,3-Butadiene | 0.000 | | 0 | N.D. | | | |
| 10) Bromomethane | 0.000 | | 0 | N.D. | | | 0.00 |
| 11) Chloroethane | 0.000 | | 0 | N.D. | | | 0.00 |
| 12) Dichlorofluoromethane | 0.000 | | 0 | N.D. | | | 0.00 |
| 13) Ethanol | 7.116 | 45 | 167660 | 2.77 | ppbv | | |
| 14) VinylBromide | 0.000 | | 0 | N.D. | | | |
| 15) Acetone | 7.984 | 58 | 307980 | 4.00 | ppbv | | 0.00 |
| 16) Trichlorofluoromethane | 7.658 | 103 | 3728 | 0.13 | ppbv | | 96 |
| 17) 2-Propanol (IPA) | 8.256 | 45 | 6065 | 0.23 | ppbv | | # 1 |
| 18) Acrylonitrile | 0.000 | | 0 | N.D. | | | |
| 19) 1,1-Dichloroethene | 0.000 | | 0 | N.D. | | | Qvalue |
| 20) MethyleneChloride (DCM) | 0.000 | | 0 | N.D. | | | # 96 |
| 21) AllylChloride | 0.000 | | 0 | N.D. | d | | |
| 22) CarbonDisulfide | 0.000 | | 0 | N.D. | | | 97 |
| 23) Trichlorotrifluoroethane | 0.000 | | 0 | N.D. | | | # 9 |
| 24) trans-1,2-Dichloroethene | 0.000 | | 0 | N.D. | | | |
| 25) 1,1-Dichloroethane | 0.000 | | 0 | N.D. | | | |
| 26) MethylTertButylEther (M...) | 0.000 | | 0 | N.D. | | | |
| 27) VinylAcetate | 0.000 | | 0 | N.D. | d | | |
| 28) 2-Butanone (MEK) | 11.494 | 72 | 32720 | 0.42 | ppbv | | 93 |
| 29) cis-1,2-Dichloroethene | 0.000 | | 0 | N.D. | | | 95 |
| 30) Hexane | 11.459 | 86 | 1825 | 0.49 | ppbv | | # 60 |
| 31) Chloroform | 12.493 | 83 | 688 | N.D. | | | |
| 32) EthylAcetate | 12.136 | 43 | 1347 | N.D. | | | |

Handwritten signature

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201306.D
 Acq On : 20 May 2013 12:32
 Operator : JJG
 Sample : 130597-63048 x1 dp
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 20 15:22:35 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

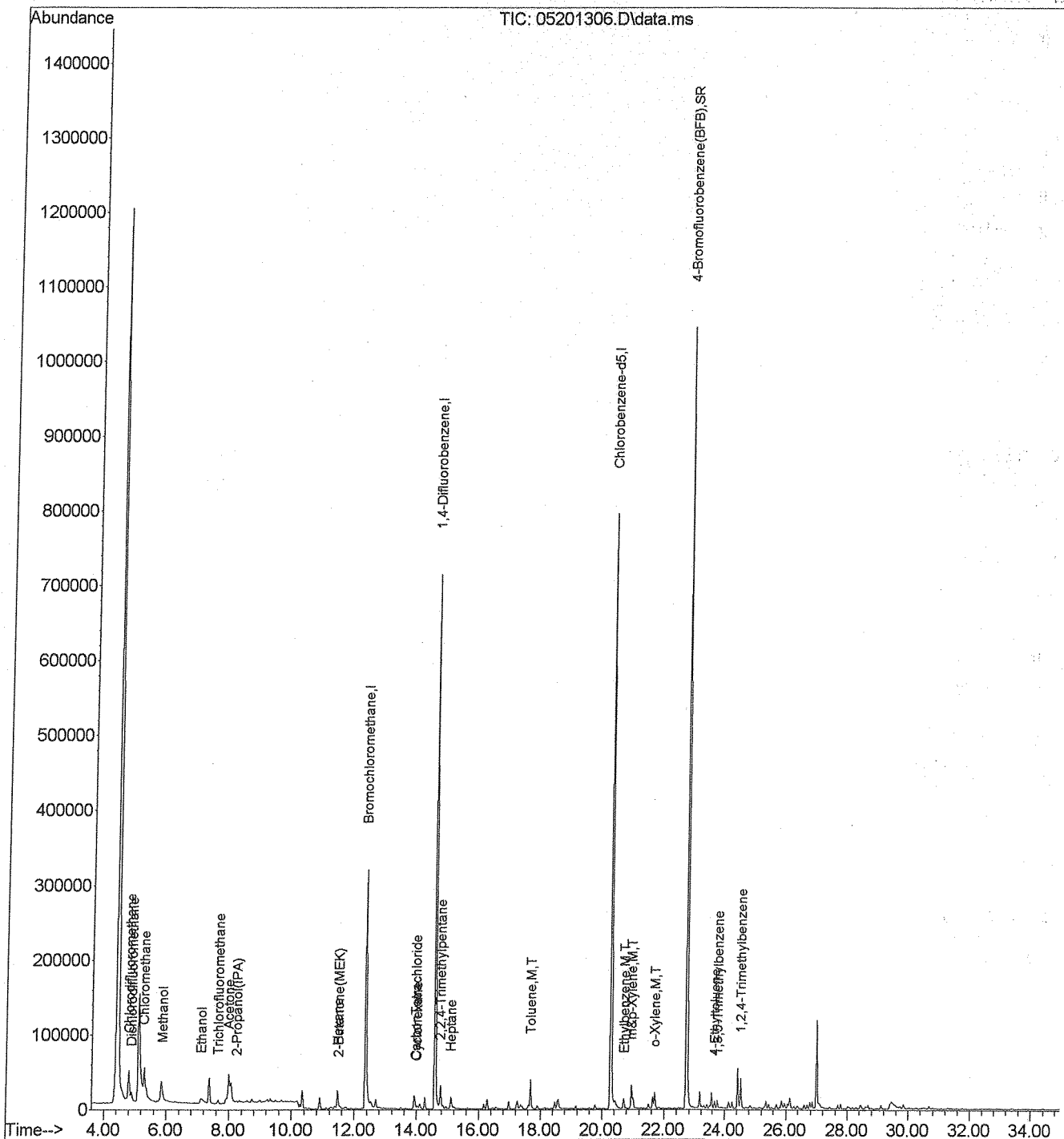
| Compound | R.T. | QIon | Response | Conc | Units | Dev (Min) |
|--------------------------------|--------|------|----------|------|----------|-----------|
| 33) Tetrahydrofuran | 0.000 | | 0 | N.D. | | |
| 34) 1,2-Dichloroethane | 0.000 | | 0 | N.D. | | |
| 35) 1,1,1-Trichloroethane | 0.000 | | 0 | N.D. | | |
| 37) Benzene | 0.000 | | 0 | N.D. | d | |
| 38) CarbonTetrachloride | 13.973 | 117 | 2283 | 0.05 | ppbv | 95 |
| 39) Cyclohexane | 14.008 | 69 | 600 | 0.07 | ppbv # | 85 |
| 40) 1,2-Dichloropropane | 0.000 | | 0 | N.D. | | |
| 41) Bromodichloromethane | 0.000 | | 0 | N.D. | | |
| 42) 1,4-Dioxane | 0.000 | | 0 | N.D. | | |
| 43) Trichloroethene (TCE) | 0.000 | | 0 | N.D. | | |
| 44) 2,2,4-Trimethylpentane | 14.757 | 57 | 38580 | 0.37 | ppbv | 98 |
| 45) Heptane | 15.096 | 71 | 4205 | 0.23 | ppbv # | 65 |
| 46) cis-1,3-Dichloropropene | 0.000 | | 0 | N.D. | | |
| 47) 4-Methyl-2-pentanone (M... | 16.576 | 58 | 528 | N.D. | | |
| 48) trans-1,3-Dichloropropene | 17.682 | 75 | 125 | N.D. | | |
| 49) 1,1,2-Trichloroethane | 0.000 | | 0 | N.D. | | |
| 50) Toluene | 17.682 | 91 | 42407 | 0.57 | ppbv (m) | 99 |
| 51) 2-Hexanone (MBK) | 0.000 | | 0 | N.D. | | |
| 52) Dibromochloromethane | 0.000 | | 0 | N.D. | | |
| 53) 1,2-Dibromoethane | 0.000 | | 0 | N.D. | | |
| 54) Tetrachloroethene (PCE) | 0.000 | | 0 | N.D. | | |
| 56) Chlorobenzene | 20.285 | 114 | 234 | N.D. | | |
| 57) Ethylbenzene | 20.713 | 91 | 15014 | 0.15 | ppbv | 99 |
| 58) m&p-Xylene | 20.945 | 106 | 19638 | 0.51 | ppbv # | 91 |
| 59) Bromoform | 0.000 | | 0 | N.D. | | |
| 60) Styrene | 21.694 | 104 | 921 | N.D. | | |
| 61) 1,1,2,2-Tetrachloroethane | 0.000 | | 0 | N.D. | | |
| 62) o-Xylene | 21.694 | 91 | 17748 | 0.23 | ppbv # | 95 |
| 64) 4-Ethyltoluene | 23.691 | 120 | 2993 | 0.09 | ppbv | 99 |
| 65) 1,3,5-Trimethylbenzene | 23.780 | 120 | 4222 | 0.09 | ppbv # | 91 |
| 66) 1,2,4-Trimethylbenzene | 24.529 | 120 | 16049 | 0.35 | ppbv | 91 |
| 67) BenzylChloride (a-Chlor... | 25.189 | 91 | 125 | N.D. | | |
| 68) 1,3-Dichlorobenzene | 25.064 | 146 | 278 | N.D. | | |
| 69) 1,4-Dichlorobenzene | 25.278 | 146 | 711 | N.D. | | |
| 70) 1,2-Dichlorobenzene | 25.849 | 146 | 383 | N.D. | | 99 |
| 71) 1,2,4-Trichlorobenzene | 29.451 | 180 | 1267 | N.D. | | |
| 72) Hexachlorobutadiene | 30.075 | 225 | 299 | N.D. | | |

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

[Handwritten signature]
 05/20/13

Data Path : C:\msdchem\1\MS03\2013\052013\
 Data File : 05201306.D
 Acq On : 20 May 2013 12:32
 Operator : JJG
 Sample : 130597-63048 x1 dp
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 20 15:22:35 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration



[Handwritten signature]
 05/20/13

Method Path : C:\msdchem\1\METHODS\2013\
 Method File : 051513.M
 Title : TO-15/TO-14
 Last Update : Thu May 16 10:13:39 2013
 Response Via : Initial Calibration

| # | ID | Conc | ISTD Conc | Path\File |
|---|-----|------|--------------|--|
| 1 | 0.5 | 1 | 10 | C:\msdchem\1\MS03\2013\051513\05151311.D |
| 2 | 1.0 | 1 | 10 | C:\msdchem\1\MS03\2013\051513\05151310.D |
| 3 | 2.0 | 2 | 10 | C:\msdchem\1\MS03\2013\051513\05151309.D |
| 4 | 5.0 | 5 | 10 | C:\msdchem\1\MS03\2013\051513\05151308.D |
| 5 | 10 | 10 | 10 | C:\msdchem\1\MS03\2013\051513\05151307.D |
| 6 | 20 | 20 | 10 | C:\msdchem\1\MS03\2013\051513\05151306.D |
| 7 | 50 | 51 | 10 | C:\msdchem\1\MS03\2013\051513\05151305.D |

| # | ID | Update Time | Quant Time | Acquisition Time |
|---|-----|-------------------|-------------------|-------------------|
| 1 | 0.5 | May 16 10:04 2013 | May 15 17:01 2013 | 15 May 2013 16:25 |
| 2 | 1.0 | May 16 10:04 2013 | May 15 16:55 2013 | 15 May 2013 15:40 |
| 3 | 2.0 | May 16 10:04 2013 | May 15 16:52 2013 | 15 May 2013 14:53 |
| 4 | 5.0 | May 16 10:03 2013 | May 15 15:19 2013 | 15 May 2013 14:07 |
| 5 | 10 | May 16 10:03 2013 | May 15 13:55 2013 | 15 May 2013 13:21 |
| 6 | 20 | May 16 10:03 2013 | May 15 13:53 2013 | 15 May 2013 12:35 |
| 7 | 50 | May 16 10:03 2013 | May 15 13:50 2013 | 15 May 2013 11:48 |

051513.M Thu May 16 10:15:15 2013

Response Factor Report MS03 Enhanced

Method Path : C:\msdchem\1\METHODS\2013\
 Method File : 051513.M
 Title : TO-15/TO-14
 Last Update : Thu May 16 10:13:39 2013
 Response Via : Initial Calibration

Calibration Files
 0.5 =05151311.D 1.0 =05151310.D 2.0 =05151309.D 5.0 =05151308.D 10 =05151307.D 20 =05151306.D
 50 =05151305.D

| Compound | 0.5 | 1.0 | 2.0 | 5.0 | 10 | 20 | 50 | Avg | %RSD |
|---------------------------|-------|-------|-------|-------|-------|-------|-------|-------|-------|
| 1) I Bromochloromethane | 2.089 | 2.153 | 2.101 | 2.025 | 1.904 | 1.810 | 1.786 | 1.981 | 7.44 |
| 2) Chlorodifluoro... | 0.516 | 0.538 | 0.543 | 0.530 | 0.511 | 0.503 | 0.502 | 0.520 | 3.16 |
| 3) Propene | 3.667 | 3.778 | 3.625 | 3.449 | 3.255 | 3.109 | 3.048 | 3.419 | 8.40 |
| 4) Dichlorodifluo... | 0.306 | 0.351 | 0.359 | 0.346 | 0.335 | 0.307 | 0.243 | 0.321 | 12.54 |
| 5) Chloromethane | 2.437 | 2.578 | 2.525 | 2.366 | 2.249 | 2.146 | 2.008 | 2.330 | 8.87 |
| 6) Dichlorotetra... | 1.119 | 1.165 | 1.194 | 1.159 | 1.138 | 1.103 | 1.085 | 1.138 | 3.35 |
| 7) Vinylchloride | 0.467 | 0.401 | 0.345 | 0.325 | 0.314 | 0.302 | 0.359 | 17.59 | 17.59 |
| 8) Methanol | 0.661 | 0.757 | 0.783 | 0.769 | 0.750 | 0.696 | 0.638 | 0.722 | 7.88 |
| 9) 1,3-Butadiene | 0.912 | 0.907 | 0.885 | 0.831 | 0.790 | 0.768 | 0.638 | 0.819 | 11.91 |
| 10) Bromomethane | 0.201 | 0.195 | 0.200 | 0.184 | 0.171 | 0.177 | 0.167 | 0.185 | 7.45 |
| 11) Chloroethane | 2.569 | 2.604 | 2.585 | 2.470 | 2.365 | 2.317 | 2.217 | 2.447 | 6.14 |
| 12) Dichlorofluoro... | 0.438 | 0.486 | 0.443 | 0.421 | 0.414 | 0.410 | 0.372 | 0.426 | 8.19 |
| 13) Ethanol | 0.913 | 0.985 | 1.013 | 1.000 | 0.954 | 0.955 | 0.914 | 0.962 | 4.10 |
| 14) VinylBromide | 0.717 | 0.643 | 0.532 | 0.487 | 0.468 | 0.473 | 0.472 | 0.542 | 18.30 |
| 15) Acetone | 2.166 | 2.190 | 2.089 | 1.956 | 1.896 | 1.864 | 1.820 | 1.997 | 7.53 |
| 16) Trichlorofluor... | 2.202 | 2.238 | 2.186 | 1.792 | 1.640 | 1.508 | 1.412 | 1.854 | 18.98 |
| 17) 2-Propanol(IPA) | 0.759 | 0.846 | 0.865 | 0.860 | 0.819 | 0.805 | 0.720 | 0.811 | 6.70 |
| 18) Acrylonitrile | 1.126 | 1.244 | 1.222 | 1.158 | 1.131 | 1.093 | 1.000 | 1.139 | 7.17 |
| 19) M,T 1,1-Dichloroet... | 1.192 | 1.150 | 1.088 | 1.078 | 1.000 | 0.935 | 0.876 | 1.046 | 10.90 |
| 20) M,T Methylenechlor... | 1.018 | 0.993 | 0.990 | 0.947 | 0.912 | 0.791 | 0.764 | 0.917 | 11.04 |
| 21) AllylChloride | 3.799 | 3.684 | 3.581 | 3.365 | 3.184 | 3.094 | 2.965 | 3.382 | 9.36 |
| 22) Carbondisulfide | 1.791 | 1.899 | 1.777 | 1.674 | 1.586 | 1.456 | 1.291 | 1.639 | 12.89 |
| 23) Trichlorotrifl... | 1.186 | 1.375 | 1.235 | 1.308 | 1.238 | 1.156 | 1.076 | 1.225 | 8.05 |
| 24) trans-1,2-Dich... | 2.829 | 2.839 | 2.719 | 2.571 | 2.383 | 2.197 | 1.949 | 2.498 | 13.53 |
| 25) 1,1-Dichloroet... | 3.471 | 3.740 | 3.529 | 3.441 | 3.284 | 3.047 | 2.705 | 3.317 | 10.38 |
| 26) MethylTertButy... | 3.239 | 3.305 | 3.371 | 3.266 | 3.089 | 2.850 | 2.479 | 3.086 | 10.33 |
| 27) VinylAcetate | 0.480 | 0.548 | 0.589 | 0.592 | 0.595 | 0.548 | 0.526 | 0.554 | 7.60 |
| 28) 2-Butanone(MEK) | | | | | | | | | |

Response Factor Report MS03 Enhanced

Method Path : C:\msdchem\1\METHODS\2013\
 Method File : 051513.M
 Title : TO-15/TO-14

| | ISTD | | | | | | | | | | | | | | | |
|-----|------|---------------------|----------------|-------|-------|-------|-------|-------|-------|-------|-------|--|--|--|--|--|
| 29) | I | cis-1,2-Dichloro... | 1.339 | 1.387 | 1.385 | 1.382 | 1.308 | 1.264 | 1.167 | 1.319 | 6.17 | | | | | |
| 30) | I | Hexane | 0.270 | 0.298 | 0.295 | 0.272 | 0.256 | 0.239 | 0.219 | 0.264 | 10.89 | | | | | |
| 31) | I | Chloroform | 2.980 | 3.005 | 2.926 | 2.844 | 2.659 | 2.617 | 2.396 | 2.775 | 8.12 | | | | | |
| 32) | I | Ethylacetate | 2.836 | 3.052 | 2.959 | 2.914 | 2.819 | 2.582 | 2.380 | 2.792 | 8.36 | | | | | |
| 33) | I | Tetrahydrofuran | 0.574 | 0.569 | 0.601 | 0.585 | 0.559 | 0.535 | 0.483 | 0.558 | 7.01 | | | | | |
| 34) | I | 1,2-Dichloroet... | 2.043 | 2.248 | 2.117 | 2.034 | 1.930 | 1.874 | 1.847 | 2.013 | 7.07 | | | | | |
| 35) | I | 1,1,1-Trichlor... | 3.350 | 3.382 | 3.301 | 3.075 | 2.948 | 2.834 | 2.618 | 3.073 | 9.43 | | | | | |
| 36) | I | 1,4-Difluorobenzene | -----ISTD----- | | | | | | | | | | | | | |
| 37) | T,M | Benzene | 0.840 | 0.830 | 0.802 | 0.758 | 0.721 | 0.648 | 0.558 | 0.737 | 14.02 | | | | | |
| 38) | I | CarbonTetrachl... | 0.664 | 0.694 | 0.674 | 0.620 | 0.600 | 0.541 | 0.469 | 0.609 | 13.26 | | | | | |
| 39) | I | Cyclohexane | 0.121 | 0.121 | 0.117 | 0.111 | 0.109 | 0.098 | 0.087 | 0.109 | 11.74 | | | | | |
| 40) | I | 1,2-Dichloropr... | 0.324 | 0.333 | 0.328 | 0.306 | 0.291 | 0.265 | 0.226 | 0.296 | 13.14 | | | | | |
| 41) | I | Bromodichlorom... | 0.411 | 0.420 | 0.409 | 0.393 | 0.371 | 0.346 | 0.311 | 0.380 | 10.47 | | | | | |
| 42) | I | 1,4-Dioxane | 0.175 | 0.182 | 0.184 | 0.183 | 0.179 | 0.165 | 0.147 | 0.174 | 7.67 | | | | | |
| 43) | M,T | Trichloroethen... | 0.378 | 0.389 | 0.390 | 0.369 | 0.352 | 0.328 | 0.293 | 0.357 | 10.02 | | | | | |
| 44) | I | 2,2,4-Trimethy... | 1.494 | 1.513 | 1.438 | 1.352 | 1.312 | 1.136 | 0.935 | 1.312 | 16.01 | | | | | |
| 45) | I | Heptane | 0.250 | 0.247 | 0.252 | 0.241 | 0.240 | 0.218 | 0.189 | 0.234 | 9.72 | | | | | |
| 46) | I | cis-1,3-Dichlo... | 0.443 | 0.456 | 0.446 | 0.443 | 0.425 | 0.389 | 0.351 | 0.422 | 9.07 | | | | | |
| 47) | I | 4-Methyl-2-pen... | 0.264 | 0.293 | 0.298 | 0.287 | 0.277 | 0.248 | 0.212 | 0.269 | 11.25 | | | | | |
| 48) | I | trans-1,3-Dich... | 0.406 | 0.438 | 0.451 | 0.457 | 0.441 | 0.423 | 0.391 | 0.430 | 5.59 | | | | | |
| 49) | I | 1,1,2-Trichlor... | 0.350 | 0.366 | 0.351 | 0.346 | 0.330 | 0.297 | 0.262 | 0.329 | 11.16 | | | | | |
| 50) | M,T | Toluene | 1.036 | 1.035 | 1.017 | 0.944 | 0.952 | 0.835 | 0.723 | 0.935 | 12.54 | | | | | |
| 51) | I | 2-Hexanone(MBK) | 0.339 | 0.353 | 0.364 | 0.351 | 0.351 | 0.307 | 0.266 | 0.333 | 10.43 | | | | | |
| 52) | I | Dibromochlorom... | 0.667 | 0.684 | 0.678 | 0.663 | 0.647 | 0.577 | 0.504 | 0.632 | 10.56 | | | | | |
| 53) | I | 1,2-Dibromoethane | 0.584 | 0.579 | 0.577 | 0.568 | 0.528 | 0.494 | 0.443 | 0.539 | 9.97 | | | | | |
| 54) | M,T | Tetrachloroeth... | 0.591 | 0.589 | 0.566 | 0.551 | 0.517 | 0.458 | 0.394 | 0.524 | 14.01 | | | | | |
| 55) | I | Chlorobenzene-d5 | -----ISTD----- | | | | | | | | | | | | | |
| 56) | M,T | Chlorobenzene | 0.280 | 0.286 | 0.275 | 0.270 | 0.258 | 0.243 | 0.208 | 0.260 | 10.37 | | | | | |
| 57) | M,T | Ethylbenzene | 1.493 | 1.508 | 1.428 | 1.373 | 1.283 | 1.189 | 0.964 | 1.320 | 14.66 | | | | | |
| 58) | M,T | m&p-Xylene | 0.595 | 0.611 | 0.556 | 0.534 | 0.502 | 0.455 | 0.383 | 0.519 | 15.49 | | | | | |
| 59) | I | Bromoform | 0.762 | 0.809 | 0.762 | 0.719 | 0.685 | 0.627 | 0.514 | 0.697 | 14.37 | | | | | |
| 60) | I | Styrene | 0.895 | 0.947 | 0.887 | 0.890 | 0.834 | 0.768 | 0.660 | 0.840 | 11.58 | | | | | |
| 61) | I | 1,1,2,2-Tetrac... | 0.834 | 0.866 | 0.828 | 0.795 | 0.730 | 0.671 | 0.544 | 0.753 | 15.13 | | | | | |
| 62) | M,T | o-Xylene | 1.255 | 1.276 | 1.150 | 1.089 | 0.982 | 0.893 | 0.746 | 1.056 | 18.39 | | | | | |
| 63) | SR | 4-Bromofluorob... | 0.633 | 0.653 | 0.621 | 0.628 | 0.618 | 0.610 | 0.617 | 0.626 | 2.29 | | | | | |
| 64) | I | 4-Ethyltoluene | 0.453 | 0.499 | 0.463 | 0.465 | 0.439 | 0.394 | 0.321 | 0.433 | 13.60 | | | | | |

051513.M Thu May 16 10:15:11 2013

Response Factor Report MS03 Enhanced

Method Path : C:\msdchem\1\METHODS\2013\
 Method File : 051513.M
 Title : TO-15/TO-14

| | | | | | | | | | | |
|-----|-------------------|-------|-------|-------|-------|-------|-------|-------|-------|-------|
| 65) | 1,3,5-Trimethy... | 0.738 | 0.738 | 0.689 | 0.655 | 0.614 | 0.551 | 0.450 | 0.633 | 16.58 |
| 66) | 1,2,4-Trimethy... | 0.683 | 0.701 | 0.662 | 0.647 | 0.606 | 0.555 | 0.465 | 0.617 | 13.44 |
| 67) | BenzylChloride... | 0.792 | 0.849 | 0.988 | 0.987 | 1.014 | 0.994 | 0.876 | 0.928 | 9.45 |
| 68) | 1,3-Dichlorobe... | 1.007 | 1.071 | 1.056 | 1.046 | 0.966 | 0.902 | 0.748 | 0.971 | 11.78 |
| 69) | 1,4-Dichlorobe... | 1.123 | 1.125 | 1.062 | 1.020 | 0.936 | 0.861 | 0.709 | 0.977 | 15.58 |
| 70) | 1,2-Dichlorobe... | 1.116 | 1.179 | 1.130 | 1.079 | 0.969 | 0.899 | 0.735 | 1.015 | 15.50 |
| 71) | 1,2,4-Trichlor... | 0.991 | 1.125 | 1.099 | 1.117 | 1.025 | 0.921 | 0.751 | 1.004 | 13.37 |
| 72) | Hexachlorobuta... | 0.954 | 0.987 | 0.907 | 0.861 | 0.768 | 0.693 | 0.522 | 0.813 | 20.26 |

(#) = Out of Range