

Atmospheric Analysis & Consulting, Inc.

CLIENT : Soil Water Air Protection Enterprise
PROJECT NAME : Bridgeton Sanitary Landfill Air Quality Assessment
AAC PROJECT NO. : 130650
REPORT DATE : 06/03/2013

On May 31, 2013, Atmospheric Analysis & Consulting, Inc. received six (6) 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 (mmHg)
BZ-1-Canister	130650-63200	593.4
BZ-2-Canister	130650-63209	485.6
U-1-Canister	130650-63218	618.8
U-2-Canister	130650-63227	602.7
D-1-Canister	130650-63236	526.9
D-2-Canister	130650-63245	487.2


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.

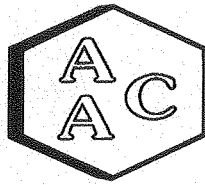
I certify that this data is technically accurate, complete and in compliance with the terms and conditions of the contract. The Laboratory Director or his designee, as verified by the following signature, has authorized the release of the data contained in this hardcopy data package.

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


Marcus Hueppe
Laboratory Director

This report consists of 100 pages.





CANISTER PRESSURE LOG

Client: Soil Water Air Protection Ent Project No.: 130650
Date: 5/31/2013

Canister #	Sample #	Initial Pressure	Final Pressure
575	63200	593.4	1032.5
576	63209	485.6	1017.2
777	63218	618.8	1019.4
716	63227	602.7	1030.6
695	63236	526.9	1023.3
653	63245	487.2	1024.5

130650
AR# 130427

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
 Telephone No. / Fax No.: (310) 434-0110 / (310) 434-0011

REQUESTED TESTS / ANALYSES

LAB ID	SAMPLE ID NUMBER	Type	Date	Time	VOCS - EPA TO-15	Reduced Sulfur Compounds - ASTM D5504	Aldehydes - 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	
63200	BZ-1-Caulster	Can	5/29/13	12:09	X	X								X				
63201	-DNPH	Tube		11:57			X											
63202	-Aews			12:08			X											
63203	-HCL			11:59			X											
63204	-AMMONIA			12:02				X										
63205	-SO2			12:05					X									
63206	-HEN			12:01					X									
63207	-AMINES			12:04						X								
63208	-MERCURY			12:07							X							

Special Instructions / Conditions of Receipt

Date: 5/29/13 Page 1 of 1

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

Relinquished By: Paul Rosefeld	Date: 5/29/13	Time: 10:30	Received By: [Signature]	Date: 5/29/13	Time: 14:15
Relinquished By: [Signature]	Date:	Time:	Received By: [Signature]	Date:	Time:
Relinquished By:	Date:	Time:	Received By:	Date:	Time:

APC # 130650

CHAIN OF CUSTODY RECORD / ANALYTICAL REQUEST FORM

Bridgeton Sanitary Landfill Air Quality Assessment

Client Name: SOIL / WATER AIR PROTECTION ENTERPRISE Telephone No. / Fax No.: (310) 434-0110 / (310) 434-0011 Date: 5/29/13 Page 1 of 1

Project Manager: PAUL ROSENFELD, PH.D. REQUESTED TESTS / ANALYSES

Address: 1640 FIFTH STREET, SUITE 204, SANTA MONICA, CA 90401

Project Name and Location: BRIDGETON SANITARY LANDFILL AIR QUALITY ASSESSMENT

Sampled By: Paul Rosenfeld Sample Signature: Paul Rosenfeld Special Instructions / Conditions of Receipt

LAB ID	SAMPLE ID NUMBER	Type	Date	Time	VOCS - EPA TO-15	Reduced Sulfur Compounds - ASTM D5504	Aldehydes - 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
63209	BZ-2-CANISTER	Can	5/29/13	16:25	X	X								X			
63210	-DNRH	TUBE		16:16			X										
63211	-ACIDS			16:19			X										
63212	-HCL			16:14				X									
63213	-AMMONIA			16:21					X								
63214	-SO2			16:22						X							
63215	-HEN			16:17							X						
63216	-AMINES			16:18								X					
63217	-MERCURY			16:21											X		

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

Relinquished By: [Signature] Date: 5/29/13 Time: 10:30 Received By: [Signature] Date: 5/30/13 Time: 14:15

Relinquished By: [Signature] Date: [] Time: [] Received By: [Signature] Date: [] Time: []

Relinquished By: [Signature] Date: [] Time: [] Received By: [Signature] Date: [] Time: []

AAE # 130650

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.
 Telephone No. / Fax No.: (310) 434-0110 / (310) 434-0011
 Date: 5/29/13 Page 1 of 1

Address: 1640 FIFTH STREET, SUITE 204, SANTA MONICA, CA 90401
 Project Name and Location: BRIDGETON SANITARY LANDFILL AIR QUALITY ASSESSMENT
 Requested Tests / Analyses

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

LAB ID	SAMPLE ID NUMBER	Type	Date	Time	VOCS - EPA TO-15	Reduced Sulfur Compounds - ASTM D5504	Aldehydes - 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	Special Instructions / Conditions of Receipt
603218	U-1 - Vanister	Can	5/29/13	12:46	X	X								X				
603219	-DNPH	Tube		12:42		X												
603220	-Acids			12:47			X											
603221	-HEL			12:49				X										
603222	-AMMONIA			12:50					X									
603223	-SO2			12:46						X								
603224	-HEN			12:51							X							
603225	-AMINES			12:45								X						
603226	-MERCURY			12:48												X		

Relinquished By: *Paul Rosenfeld* Date: 5/29/13 Time: 18:30
 Relinquished By: *Paul Rosenfeld* Date: Date: Time: Time:
 Relinquished By: Date: Time: Received By: *WJ77* Date: 5/30/13 Time: 14:15

SOIL / WATER / AIR PROTECTION ENTERPRISE
 SAMPLES RECEIVED @ 7.0 '0

APP # 130650

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
 Telephone No. / Fax No.: (310) 434-0110 / (310) 434-0011
 Date: 5/29/13 Page 1 of 1

Project Name and Location: BRIDGETON SANITARY LANDFILL AIR QUALITY ASSESSMENT
 Sampled By: Paul Rosenfeld
 Sample Signature: Paul Rosenfeld
 Requested Tests / Analyses:
 VOCs - EPA TO-15
 Reduced Sulfur Compounds - ASTM D5504
 Aldehydes - 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
 Special Instructions / Conditions of Receipt

LAB ID	SAMPLE ID NUMBER	Type	Date	Time	VOCs - EPA TO-15	Reduced Sulfur Compounds - ASTM D5504	Aldehydes - 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
03227	U-2-CHARGER	CAN	5/21/13	16:46	X	X								X			
03228	- DIRT	TUBE		16:43			X										
03229	- ACIDS			16:41			X										
03230	- HCL			16:40				X									
03231	- AMMONIA			16:44					X								
03232	- SO2			16:45						X							
03233	- HCN			16:42							X						
03234	- AMINES			16:42								X					
03235	- MERCURY			16:39											X		

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

Relinquished By: ROB HESS	Date: 5/29/13	Time: 18:30	Received By: [Signature]	Date: 5/30/13	Time: 14:5
Relinquished By: [Signature]	Date:	Time:	Received By: [Signature]	Date:	Time:
Relinquished By:	Date:	Time:	Received By:	Date:	Time:

AAE # 130650

CHAIN OF CUSTODY RECORD / ANALYTICAL REQUEST FORM

Bridgeton Sanitary Landfill Air Quality Assessment

Client Name: SOIL / WATER AIR PROTECTION ENTERPRISE Telephone No. / Fax No.: (310) 434-0110 / (310) 434-0011
 Project Manager: PAUL ROSENFELD, PH.D. Date: 5/29/13 Page 1 of 1

Address: 1640 FIFTH STREET, SUITE 204, SANTA MONICA, CA 90401
 Project Name and Location: BRIDGETON SANITARY LANDFILL AIR QUALITY ASSESSMENT
 Requested Tests / Analyses: VOCs - EPA TO-15, Reduced Sulfur Compounds - ASTM D5504, Aldehydes - 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

Sampled By: Paul Rosenfeld Signature: Paul Rosenfeld
 Special Instructions / Conditions of Receipt

LAB ID	SAMPLE ID NUMBER	Type	Date	Time	VOCs - EPA TO-15	Reduced Sulfur Compounds - ASTM D5504	Aldehydes - 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
03230	D-1 - CANISTER	CAN	5/29/13	13:28	X									X			
03237	- DNPH	TUBE		13:29			X										
03238	- ACIDS			13:26			X										
03239	- HEL			13:23				X									
03240	- AMMONIA			13:18					X								
03241	- SO2			13:20						X							
03242	- HCN			13:22							X						
03243	- AMINES			13:25								X					
03244	- MERCURY		5/29	13:21												X	

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

Relinquished By: Rob Hesse Date: 5/29/13 Time: 18:30
 Received By: [Signature] Date: 5/29/13 Time: 14:15

Relinquished By: [Signature] Date: [] Time: []
 Received By: [Signature] Date: [] Time: []

Relinquished By: [Signature] Date: [] Time: []
 Received By: [Signature] Date: [] Time: []

SOIL / WATER / AIR PROTECTION ENTERPRISE

ARC# 132650

CHAIN OF CUSTODY RECORD / ANALYTICAL REQUEST FORM

Bridgeton Sanitary Landfill Air Quality Assessment

Client Name: SOIL / WATER AIR PROTECTION ENTERPRISE
 Telephone No. / Fax No.: (310) 434-0110 / (310) 434-0011
 Project Manager: PAUL ROSENFELD, PH.D.
 Address: 1640 FIFTH STREET, SUITE 204, SANTA MONICA, CA 90401
 Date: 5/29/13 Page 1 of 1

Project Name and Location: BRIDGETON SANITARY LANDFILL AIR QUALITY ASSESSMENT
 Requested Tests / Analyses: VOCs - EPA TO-15, Reduced Sulfur Compounds - ASTM D5504, Aldehydes - 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
 Special Instructions / Conditions of Receipt

LAB ID	SAMPLE ID NUMBER	Type	Date	Time	VOCs - EPA TO-15	Reduced Sulfur Compounds - ASTM D5504	Aldehydes - 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
63245	D-2 - CANISTER	Can	5/29/13	17:20	X		X							X			
63246	- DNRPH	TOBE		17:20													
63247	- AC105			17:21			X										
63248	- HCL			17:24				X									
63249	- AMMONIA			17:25					X								
63250	- SO2			17:22						X							
63251	- HCN			17:26					X								
63252	- AMINES			17:27							X						
63253	- MERCURY	Y		17:23											X		

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

Relinquished By: KOBHACK	Date: 5/29/13	Time: 18:30	Received By: [Signature]	Date: 5/30/13	Time: 14:15
Relinquished By:	Date:	Time:	Received By:	Date:	Time:
Relinquished By:	Date:	Time:	Received By:	Date:	Time:

AAE # 130650

CHAIN OF CUSTODY RECORD / ANALYTICAL REQUEST FORM

Bridgeton Sanitary Landfill Air Quality Assessment

Client Name: SOIL / WATER AIR PROTECTION ENTERPRISE
 Telephone No. / Fax No.: (310) 434-0110 / (310) 434-0011
 Date: 5/29/12 Page 1 of 1

Project Manager: PAUL ROSENFELD, PH.D.
REQUESTED TESTS / ANALYSES

Address: 1640 FIFTH STREET, SUITE 204, SANTA MONICA, CA 90401

Project Name and Location: BRIDGETON SANITARY LANDFILL AIR QUALITY ASSESSMENT

Sampled By: Paul Rosenfeld
 Sampler Signature: Paul Rosenfeld

LAB ID	SAMPLE ID NUMBER	Type	Date	Time	VOCS - EPA TO-15	Reduced Sulfur Compounds - ASTM D5504	Aldehydes - 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	Special Instructions / Conditions of Receipt
63254	TRIP BLANK - DNPH	TUBE	5/29/12	15:00			X											
63255	-ACIDS			15:30			X											
63256	-HCL						X											
63257	-AMMONIA							X										
63258	-SO2								X									
63259	-HCN								X									
63260	-AMINES									X								
63261	-MERCURY										X							

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

Relinquished By: ROBHESSE
 Date: 5/29/12 Time: 18:30
 Received By: [Signature]
 Date: 5/29/12 Time: 14:15

Relinquished By: [Signature]
 Date: [Blank] Time: [Blank]
 Received By: [Signature]
 Date: 5/29/12 Time: 14:15

Atmospheric Analysis and Consulting Inc.

Canister Sampling Field Data Sheet

GENERAL INFORMATION

Project Name and/or ID No.: SWAPE Project: 601 Bridgeton Landfill
 Site Address and/or ID No.: Bridgeton Sanitary Landfill, Bridgeton, Missouri
 Sample Name and/or ID No.: BZ-1 - Canister
 AAC Batch ID: 130650 AAC Sample ID: 63200

SAMPLING INFORMATION

Start Date/Time: 5/29/13 7:51 Stop Date/Time: 5/29/13 12:09
 Start Temp/Pressure*: 25°C / 29.96 Stop Temp/Pressure*: 30.6°C / 29.90
 Initial Can Pressure**: -29 Final Can Pressure**: -6.2

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

Comments: Temp / Pressure DATA FROM LAMBERT INTL.

ROB HESSE [Signature] 5/29/13
 Sampler Name (Print) Sampler Signature/Date

LABORATORY INFORMATION

Canister Size: 6-Liter Sampling Period: 4-Hour
 Canister Serial No.: 575 Flow Controller Serial No.: 804
 Initial Pressure: 4.5 Certified Flow Rate: 18.0
 Return Pressure: 593.4 Certified By/Date: [Signature] 5/23/2013
 Final Pressure: 1032.5 Flow Rate upon Return: 22.0

Date Shipped From Lab: 5/20/2013 Shipped By: [Signature]
 Date Returned to Lab: 5/31/2013 Received By: [Signature]
 Flow Controller Certification File ID: MS03/052013/05231305
 Canister Certification File ID: MS03/04051313
 Certification Type: SIM SCAN NJLL PAMS Other

[Signature] 6/5/13
 Chemist Signature/Date 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.: SWAPE Project: 601 Bridgeton Landfill
 Site Address and/or ID No.: Bridgeton Sanitary Landfill, Bridgeton, Missouri
 Sample Name and/or ID No.: BZ-2-CANISTER
 AAC Batch ID: 130650 AAC Sample ID: 63209

SAMPLING INFORMATION

Start Date/Time: 5/29/13 12:09 Stop Date/Time: 5/29/13 16:22
 Start Temp/Pressure*: 30.00 / 29.90 Stop Temp/Pressure*: 30.00°C / 29.85
 Initial Can Pressure**: -29 Final Can Pressure**: -9

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

Comments: THIS CANISTER / FLOW CONTROLLER MARKED "D-2" BY JOHN BLANK

TEMP / PRESSURE DATA FROM LAMBERT INTL.
ROB HESSE [Signature] 5/29/13
 Sampler Name (Print) Sampler Signature/Date

LABORATORY INFORMATION

Canister Size: 6-Liter Sampling Period: 4-Hour
 Canister Serial No.: 576 Flow Controller Serial No.: 694
 Initial Pressure: 4.2 Certified Flow Rate: 18.0
 Return Pressure: 485.6 Certified By/Date: [Signature] 5/20/2013
 Final Pressure: 1017.2 Flow Rate upon Return: 18.8

Date Shipped From Lab: 5/16/2013 Shipped By: [Signature]
 Date Returned to Lab: 5/31/2013 Received By: [Signature]
 Flow Controller Certification File ID: 1603052013104
 Canister Certification File ID: 160305151318
 Certification Type: SIM SCAN NJLL PAMS Other

[Signature] 06/03/13 [Signature] 6/5/13
 Chemist Signature/Date 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.: SWAPE Project: 601 Bridgeton Landfill
 Site Address and/or ID No.: Bridgeton Sanitary Landfill, Bridgeton, Missouri
 Sample Name and/or ID No.: U-1-CANISTER
 AAC Batch ID: 130650 AAC Sample ID: 63218

SAMPLING INFORMATION

Start Date/Time: 5/29/13 8:10 Stop Date/Time: 5/29/13 12:46
 Start Temp/Pressure*: 25°C | 29.96 Stop Temp/Pressure*: 30.6°C | 29.90
 Initial Can Pressure**: -78 Final Can Pressure**: -4

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

Comments: TEMP/PRESSURE DATA FROM LAMBERTS INTL.

ROB HESSE [Signature] 5/29/13
 Sampler Name (Print) Sampler Signature/Date

LABORATORY INFORMATION

Canister Size: 6-Liter Sampling Period: 4-Hour
 Canister Serial No.: 777 Flow Controller Serial No.: 803
 Initial Pressure: 4.7 Certified Flow Rate: 18.0
 Return Pressure: 618.8 Certified By/Date: [Signature] 5/23/2013
 Final Pressure: 1019.4 Flow Rate upon Return: 19.0

Date Shipped From Lab: 5/20/2013 Shipped By: [Signature]
 Date Returned to Lab: 5/31/2013 Received By: [Signature]
 Flow Controller Certification File ID: M603/05201305
 Canister Certification File ID: M603/04051314
 Certification Type: SIM SCAN NJLL PAMS Other

[Signature] 06/03/13 [Signature]
 Chemist Signature/Date 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.: SWAPE Project: 601 Bridgeton Landfill
 Site Address and/or ID No.: Bridgeton Sanitary Landfill, Bridgeton, Missouri
 Sample Name and/or ID No.: V-2-Canister
 AAC Batch ID: 130650 AAC Sample ID: 63227

SAMPLING INFORMATION

Start Date/Time: 5/29/13 12:57 Stop Date/Time: 5/29/13 16:46
 Start Temp/Pressure*: 30.6°C/29.70 Stop Temp/Pressure*: 30.0°C/29.86
 Initial Can Pressure**: -30 Final Can Pressure**: -6.6

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

Comments: Temp/pressure DATA FROM LAMBERT INTL.

ROB HESSE [Signature] 5/29/13
 Sampler Name (Print) Sampler Signature/Date

LABORATORY INFORMATION

Canister Size: 6-Liter Sampling Period: 4-Hour
 Canister Serial No.: 716 Flow Controller Serial No.: 709
 Initial Pressure: 4.6 Certified Flow Rate: 18.0
 Return Pressure: 602.7 Certified By/Date: JJ 5/23/2013
 Final Pressure: 1030.6 Flow Rate upon Return: 26.0

Date Shipped From Lab: 5/20/2013 Shipped By: JJ
 Date Returned to Lab: 5/31/2013 Received By: JJ
 Flow Controller Certification File ID: 1403/0523/305
 Canister Certification File ID: 1403/0510/319
 Certification Type: SIM SCAN NJLL PAMS Other

[Signature] 06/05/13 [Signature] 6/5/13
 Chemist Signature/Date 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.: SWAPE Project: 601 Bridgeton Landfill
 Site Address and/or ID No.: Bridgeton Sanitary Landfill, Bridgeton, Missouri
 Sample Name and/or ID No.: D-1-Canister
 AAC Batch ID: 130650 AAC Sample ID: 63236

SAMPLING INFORMATION

Start Date/Time: 5/29/13 8:43 Stop Date/Time: 5/29/13 13:28
 Start Temp/Pressure*: 25.0 / 29.96 Stop Temp/Pressure*: 30.6 / 29.90
 Initial Can Pressure**: -29.5 Final Can Pressure**: -10

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

Comments: Temp / Pressure Data from Lambert Intl.

Rob Hesse [Signature] 5/29/13
 Sampler Name (Print) Sampler Signature/Date

LABORATORY INFORMATION

Canister Size: 6-Liter Sampling Period: 4-Hour
 Canister Serial No.: 695 Flow Controller Serial No.: 805
 Initial Pressure: 4.6 Certified Flow Rate: 18.0
 Return Pressure: 526.9 Certified By/Date: [Signature] 5/23/2013
 Final Pressure: 1023.3 Flow Rate upon Return: 20.0

Date Shipped From Lab: 5/20/2013 Shipped By: [Signature]
 Date Returned to Lab: 5/31/2013 Received By: [Signature]
 Flow Controller Certification File ID: MS03/0523305
 Canister Certification File ID: MS03/05151326
 Certification Type: SIM SCAN NJLL PAMS Other

[Signature] [Signature]
 Chemist Signature/Date 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.: SWAPE Project: 601 Bridgeton Landfill
 Site Address and/or ID No.: Bridgeton Sanitary Landfill, Bridgeton, Missouri
 Sample Name and/or ID No.: D-2-Canister
 AAC Batch ID: 130650 AAC Sample ID: 63245

SAMPLING INFORMATION

Start Date/Time: 5/29/13 13:45 Stop Date/Time: 5/29/13 17:20
 Start Temp/Pressure*: 30.6°C/29.90 Stop Temp/Pressure*: 30.0°C/29.86
 Initial Can Pressure**: -28.5 Final Can Pressure**: -10

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

Comments: THIS CANISTER / FLOW CONTROLLER MARKED "D-1"
BY JOHN BLANK

TEMP / PRESSURE DATA FROM LAMBERT INTL.
Rob Hesse [Signature] 5/29/13
 Sampler Name (Print) Sampler Signature/Date

LABORATORY INFORMATION

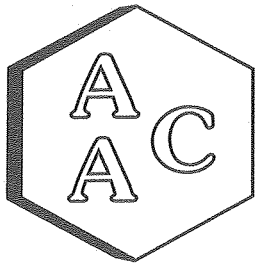
Canister Size: 6-Liter Sampling Period: 4-Hour
 Canister Serial No.: 653 Flow Controller Serial No.: 715
 Initial Pressure: 4.3 Certified Flow Rate: 18.0
 Return Pressure: 487.2 Certified By/Date: [Signature] 5/20/2013
 Final Pressure: 1024.5 Flow Rate upon Return: 18.4

Date Shipped From Lab: 5/16/2013 Shipped By: [Signature]
 Date Returned to Lab: 5/31/2013 Received By: [Signature]
 Flow Controller Certification File ID: M503/0580310
 Canister Certification File ID: M503/0515321
 Certification Type: SIM SCAN NJLL PAMS Other

[Signature] 6/5/13
 Chemist Signature/Date 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

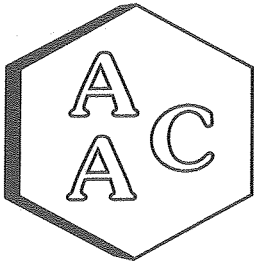
CLIENT : Soil Water Air Protection Enterprise
PROJECT NO : 130650
MATRIX : AIR
UNITS : PPB (v/v)

DATE RECEIVED : 05/31/2013
DATE REPORTED : 06/03/2013

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

Client ID AAC ID	BZ-1-Canister			Sample Reporting Limit (SRL) (MRLxDF's)	BZ-2-Canister			Sample Reporting Limit (SRL) (MRLxDF's)	Method Reporting Limit (MRL)
	130650-63200	05/29/2013			130650-63209	05/29/2013			
Date Analyzed	05/31/2013			2.09	05/31/2013			2.09	
Can Dilution Factor	1.74				Result	Qualifier	Analysis DF		
Chlorodifluoromethane	0.38	J	1.0	0.87	0.31	J	1.0	1.05	0.5
Propene	1.81		1.0	1.74	2.56		1.0	2.09	1.0
Dichlorodifluoromethane	0.59	J	1.0	0.87	0.54	J	1.0	1.05	0.5
Chloromethane	0.47	J	1.0	0.87	0.59	J	1.0	1.05	0.5
Dichlorotetrafluoroethane	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Vinyl Chloride	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Methanol	143		10.0	87.0	298		10.0	105	5.0
1,3-Butadiene	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Bromomethane	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Chloroethane	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Dichlorofluoromethane	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Ethanol	19.0		1.0	3.48	35.3		1.0	4.19	2.0
Vinyl Bromide	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Acetone	14.0		1.0	3.48	24.0		1.0	4.19	2.0
Trichlorofluoromethane	0.33	J	1.0	0.87	0.27	J	1.0	1.05	0.5
2-Propanol (IPA)	3.39	J	1.0	3.48	5.61		1.0	4.19	2.0
Acrylonitrile	<SRL	U	1.0	1.74	<SRL	U	1.0	2.09	1.0
1,1-Dichloroethene	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Methylene Chloride (DCM)	<SRL	U	1.0	1.74	<SRL	U	1.0	2.09	1.0
Allyl Chloride	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Carbon Disulfide	NR	U	1.0	0.87	NR	U	1.0	1.05	0.5
Trichlorotrifluoroethane	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
trans-1,2-Dichloroethene	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
1,1-Dichloroethane	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Methyl Tert Butyl Ether (MTBE)	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Vinyl Acetate	<SRL	U	1.0	1.74	<SRL	U	1.0	2.09	1.0
2-Butanone (MEK)	6.40		1.0	1.74	11.5		1.0	2.09	1.0
cis-1,2-Dichloroethene	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Hexane	0.66	J	1.0	0.87	<SRL	U	1.0	1.05	0.5
Chloroform	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Ethyl Acetate	0.50	J	1.0	0.87	1.13		1.0	1.05	0.5
Tetrahydrofuran	3.20		1.0	0.87	6.14		1.0	1.05	0.5
1,2-Dichloroethane	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
1,1,1-Trichloroethane	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5





Atmospheric Analysis & Consulting, Inc.

Laboratory Analysis Report

CLIENT : Soil Water Air Protection Enterprise
PROJECT NO : 130650
MATRIX : AIR
UNITS : PPB (v/v)

DATE RECEIVED : 05/31/2013
DATE REPORTED : 06/03/2013

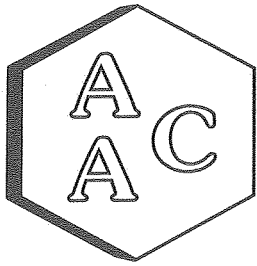
VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

Client ID AAC ID	BZ-1-Canister			Sample Reporting Limit (SRL) (MRLxDF's)	BZ-2-Canister			Sample Reporting Limit (SRL) (MRLxDF's)	Method Reporting Limit (MRL)
	130650-63200				130650-63209				
Date Sampled	05/29/2013				05/29/2013				
Date Analyzed	05/31/2013				05/31/2013				
Can Dilution Factor	1.74				2.09				
	Result	Qualifier	Analysis DF		Result	Qualifier	Analysis DF		
Benzene	3.86		1.0	0.87	7.27		1.0	1.05	0.5
Carbon Tetrachloride	0.09	J	1.0	0.87	<SRL	U	1.0	1.05	0.5
Cyclohexane	0.12	J	1.0	0.87	<SRL	U	1.0	1.05	0.5
1,2-Dichloropropane	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Bromodichloromethane	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
1,4-Dioxane	<SRL	U	1.0	0.87	0.44	J	1.0	1.05	0.5
Trichloroethene (TCE)	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
2,2,4-Trimethylpentane	0.56	J	1.0	0.87	0.17	J	1.0	1.05	0.5
Heptane	0.30	J	1.0	0.87	0.15	J	1.0	1.05	0.5
cis-1,3-Dichloropropene	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
4-Methyl-2-pentanone (MiBK)	0.17	J	1.0	0.87	0.29	J	1.0	1.05	0.5
trans-1,3-Dichloropropene	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
1,1,2-Trichloroethane	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Toluene	1.69	J	1.0	0.87	1.95	J	1.0	1.05	0.5
2-Hexanone (MBK)	<SRL	U	1.0	0.87	0.27	J	1.0	1.05	0.5
Dibromochloromethane	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
1,2-Dibromoethane	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Tetrachloroethene (PCE)	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Chlorobenzene	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Ethylbenzene	0.57	J	1.0	0.87	0.92	J	1.0	1.05	0.5
m & p-Xylenes	1.53	J	1.0	1.74	2.18	J	1.0	2.09	1.0
Bromoform	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Styrene	<SRL	U	1.0	0.87	0.13	J	1.0	1.05	0.5
1,1,2,2-Tetrachloroethane	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
o-Xylene	0.63	J	1.0	0.87	0.96	J	1.0	1.05	0.5
4-Ethyltoluene	0.19	J	1.0	0.87	0.34	J	1.0	1.05	0.5
1,3,5-Trimethylbenzene	0.21	J	1.0	0.87	0.38	J	1.0	1.05	0.5
1,2,4-Trimethylbenzene	0.66	J	1.0	0.87	1.38	J	1.0	1.05	0.5
Benzyl Chloride (a-Chlorotoluene)	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
1,3-Dichlorobenzene	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
1,4-Dichlorobenzene	0.16	J	1.0	0.87	0.80	J	1.0	1.05	0.5
1,2-Dichlorobenzene	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
1,2,4-Trichlorobenzene	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
Hexachlorobutadiene	<SRL	U	1.0	0.87	<SRL	U	1.0	1.05	0.5
BFB-Surrogate Std. % Recovery	104%				105%				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

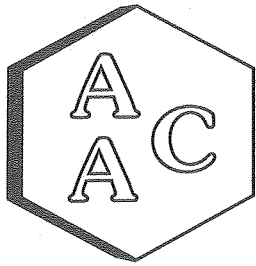
CLIENT : Soil Water Air Protection Enterprise
PROJECT NO : 130650
MATRIX : AIR
UNITS : PPB (v/v)

DATE RECEIVED : 05/31/2013
DATE REPORTED : 06/03/2013

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

Client ID	U-1-Canister			Sample Reporting Limit (SRL) (MRLxDF's)	U-2-Canister			Sample Reporting Limit (SRL) (MRLxDF's)	Method Reporting Limit (MRL)
	AAC ID	Result	Qualifier		Analysis DF	Result	Qualifier		
	130650-63218				130650-63227				
Date Sampled	05/29/2013				05/29/2013				
Date Analyzed	05/31/2013				05/31/2013				
Can Dilution Factor	1.65				1.71				
	Result	Qualifier	Analysis DF		Result	Qualifier	Analysis DF		
Chlorodifluoromethane	0.35	J	1.0	0.82	0.31	J	1.0	0.85	0.5
Propene	0.54	J	1.0	1.65	0.53	J	1.0	1.71	1.0
Dichlorodifluoromethane	0.56	J	1.0	0.82	0.55	J	1.0	0.85	0.5
Chloromethane	0.48	J	1.0	0.82	0.50	J	1.0	0.85	0.5
Dichlorotetrafluoroethane	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Vinyl Chloride	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Methanol	10.6		1.0	8.24	11.1		1.0	8.55	5.0
1,3-Butadiene	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Bromomethane	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Chloroethane	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Dichlorofluoromethane	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Ethanol	3.59		1.0	3.29	4.36		1.0	3.42	2.0
Vinyl Bromide	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Acetone	3.72		1.0	3.29	4.55		1.0	3.42	2.0
Trichlorofluoromethane	0.26	J	1.0	0.82	0.27	J	1.0	0.85	0.5
2-Propanol (IPA)	0.41	J	1.0	3.29	0.65	J	1.0	3.42	2.0
Acrylonitrile	<SRL	U	1.0	1.65	<SRL	U	1.0	1.71	1.0
1,1-Dichloroethene	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Methylene Chloride (DCM)	<SRL	U	1.0	1.65	<SRL	U	1.0	1.71	1.0
Allyl Chloride	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Carbon Disulfide	NR	U	1.0	0.82	NR	U	1.0	0.85	0.5
Trichlorotrifluoroethane	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
trans-1,2-Dichloroethene	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
1,1-Dichloroethane	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Methyl Tert Butyl Ether (MTBE)	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Vinyl Acetate	<SRL	U	1.0	1.65	<SRL	U	1.0	1.71	1.0
2-Butanone (MEK)	0.49	J	1.0	1.65	0.44	J	1.0	1.71	1.0
cis-1,2-Dichloroethene	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Hexane	0.69	J	1.0	0.82	0.43	J	1.0	0.85	0.5
Chloroform	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Ethyl Acetate	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Tetrahydrofuran	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
1,2-Dichloroethane	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
1,1,1-Trichloroethane	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5





Atmospheric Analysis & Consulting, Inc.

Laboratory Analysis Report

CLIENT : Soil Water Air Protection Enterprise
PROJECT NO : 130650
MATRIX : AIR
UNITS : PPB (v/v)

DATE RECEIVED : 05/31/2013
DATE REPORTED : 06/03/2013

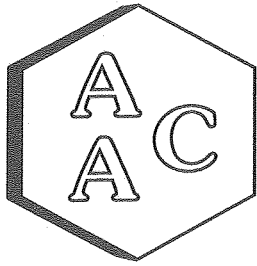
VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

Client ID	U-1-Canister			Sample Reporting Limit (SRL) (MRLxDF's)	U-2-Canister			Sample Reporting Limit (SRL) (MRLxDF's)	Method Reporting Limit (MRL)
	AAC ID	Date Sampled	Date Analyzed		AAC ID	Date Sampled	Date Analyzed		
	130650-63218	05/29/2013	05/31/2013		130650-63227	05/29/2013	05/31/2013		
Can Dilution Factor	1.65				1.71				
	Result	Qualifier	Analysis DF		Result	Qualifier	Analysis DF		
Benzene	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Carbon Tetrachloride	0.08	J	1.0	0.82	0.09	J	1.0	0.85	0.5
Cyclohexane	0.10	J	1.0	0.82	<SRL	U	1.0	0.85	0.5
1,2-Dichloropropane	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Bromodichloromethane	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
1,4-Dioxane	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Trichloroethene (TCE)	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
2,2,4-Trimethylpentane	0.56	J	1.0	0.82	0.46	J	1.0	0.85	0.5
Heptane	0.21	J	1.0	0.82	0.24	J	1.0	0.85	0.5
cis-1,3-Dichloropropene	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
4-Methyl-2-pentanone (MiBK)	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
trans-1,3-Dichloropropene	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
1,1,2-Trichloroethane	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Toluene	0.92	J	1.0	0.82	0.89	J	1.0	0.85	0.5
2-Hexanone (MBK)	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Dibromochloromethane	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
1,2-Dibromoethane	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Tetrachloroethene (PCE)	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Chlorobenzene	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Ethylbenzene	0.40	J	1.0	0.82	0.39	J	1.0	0.85	0.5
m & p-Xylenes	1.42	J	1.0	1.65	1.44	J	1.0	1.71	1.0
Bromoform	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Styrene	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
1,1,2,2-Tetrachloroethane	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
o-Xylene	0.56	J	1.0	0.82	0.56	J	1.0	0.85	0.5
4-Ethyltoluene	0.13	J	1.0	0.82	0.12	J	1.0	0.85	0.5
1,3,5-Trimethylbenzene	0.12	J	1.0	0.82	0.12	J	1.0	0.85	0.5
1,2,4-Trimethylbenzene	0.40	J	1.0	0.82	0.36	J	1.0	0.85	0.5
Benzyl Chloride (a-Chlorotoluene)	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
1,3-Dichlorobenzene	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
1,4-Dichlorobenzene	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
1,2-Dichlorobenzene	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
1,2,4-Trichlorobenzene	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
Hexachlorobutadiene	<SRL	U	1.0	0.82	<SRL	U	1.0	0.85	0.5
BFB-Surrogate Std. % Recovery	105%				105%				70-130%

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





Atmospheric Analysis & Consulting, Inc.

Laboratory Analysis Report

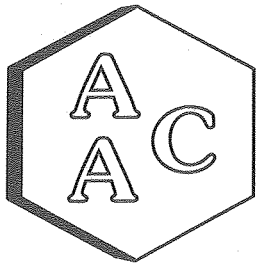
CLIENT : Soil Water Air Protection Enterprise
PROJECT NO : 130650
MATRIX : AIR
UNITS : PPB (v/v)

DATE RECEIVED : 05/31/2013
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VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

Client ID	D-1-Canister			Sample Reporting Limit (SRL) (MRLxDF's)	D-2-Canister			Sample Reporting Limit (SRL) (MRLxDF's)	Method Reporting Limit (MRL)
	AAC ID	Result	Qualifier		Analysis DF	Result	Qualifier		
	130650-63236				130650-63245				
Date Sampled	05/23/2013				05/23/2013				
Date Analyzed	05/31/2013				05/31/2013				
Can Dilution Factor	1.94				2.10				
Chlorodifluoromethane	0.39	J	1.0	0.97	0.32	J	1.0	1.05	0.5
Propene	1.18	J	1.0	1.94	1.32	J	1.0	2.10	1.0
Dichlorodifluoromethane	0.58	J	1.0	0.97	0.57	J	1.0	1.05	0.5
Chloromethane	0.45	J	1.0	0.97	0.53	J	1.0	1.05	0.5
Dichlorotetrafluoroethane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Vinyl Chloride	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Methanol	68.2		1.0	9.71	52.6		1.0	10.5	5.0
1,3-Butadiene	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Bromomethane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Chloroethane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Dichlorofluoromethane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Ethanol	8.37		1.0	3.88	6.31		1.0	4.21	2.0
Vinyl Bromide	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Acetone	8.80		1.0	3.88	7.40		1.0	4.21	2.0
Trichlorofluoromethane	0.31	J	1.0	0.97	0.27	J	1.0	1.05	0.5
2-Propanol (IPA)	1.98	J	1.0	3.88	1.62	J	1.0	4.21	2.0
Acrylonitrile	<SRL	U	1.0	1.94	<SRL	U	1.0	2.10	1.0
1,1-Dichloroethene	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Methylene Chloride (DCM)	<SRL	U	1.0	1.94	<SRL	U	1.0	2.10	1.0
Allyl Chloride	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Carbon Disulfide	NR	U	1.0	0.97	NR		1.0	1.05	0.5
Trichlorotrifluoroethane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
trans-1,2-Dichloroethene	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
1,1-Dichloroethane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Methyl Tert Butyl Ether (MTBE)	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Vinyl Acetate	<SRL	U	1.0	1.94	<SRL	U	1.0	2.10	1.0
2-Butanone (MEK)	2.39		1.0	1.94	2.08	J	1.0	2.10	1.0
cis-1,2-Dichloroethene	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Hexane	0.76	J	1.0	0.97	<SRL	U	1.0	1.05	0.5
Chloroform	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Ethyl Acetate	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Tetrahydrofuran	0.97		1.0	0.97	0.95	J	1.0	1.05	0.5
1,2-Dichloroethane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
1,1,1-Trichloroethane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5





Atmospheric Analysis & Consulting, Inc.

Laboratory Analysis Report

CLIENT : Soil Water Air Protection Enterprise
PROJECT NO : 130650
MATRIX : AIR
UNITS : PPB (v/v)

DATE RECEIVED : 05/31/2013
DATE REPORTED : 06/03/2013

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

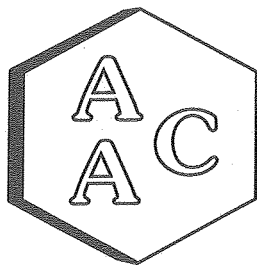
Client ID	D-1-Canister			Sample Reporting Limit (SRL) (MRLxDF's)	D-2-Canister			Sample Reporting Limit (SRL) (MRLxDF's)	Method Reporting Limit (MRL)
	AAC ID	Result	Qualifier		Analysis DF	Result	Qualifier		
	130650-63236				130650-63245				
Date Sampled	05/23/2013				05/23/2013				
Date Analyzed	05/31/2013				05/31/2013				
Can Dilution Factor	1.94				2.10				
Benzene	1.55		1.0	0.97	1.56		1.0	1.05	0.5
Carbon Tetrachloride	<SRL	U	1.0	0.97	0.11	J	1.0	1.05	0.5
Cyclohexane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
1,2-Dichloropropane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Bromodichloromethane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
1,4-Dioxane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Trichloroethene (TCE)	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
2,2,4-Trimethylpentane	0.62	J	1.0	0.97	0.17	J	1.0	1.05	0.5
Heptane	0.31	J	1.0	0.97	0.17	J	1.0	1.05	0.5
cis-1,3-Dichloropropene	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
4-Methyl-2-pentanone (MiBK)	<SRL	U	1.0	0.97	0.08	J	1.0	1.05	0.5
trans-1,3-Dichloropropene	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
1,1,2-Trichloroethane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Toluene	1.40	J	1.0	0.97	0.76	J	1.0	1.05	0.5
2-Hexanone (MBK)	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Dibromochloromethane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
1,2-Dibromoethane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Tetrachloroethene (PCE)	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Chlorobenzene	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Ethylbenzene	0.43	J	1.0	0.97	0.23	J	1.0	1.05	0.5
m & p-Xylenes	1.11	J	1.0	1.94	0.67	J	1.0	2.10	1.0
Bromoform	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Styrene	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
1,1,2,2-Tetrachloroethane	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
o-Xylene	0.43	J	1.0	0.97	0.27	J	1.0	1.05	0.5
4-Ethyltoluene	0.16	J	1.0	0.97	<SRL	U	1.0	1.05	0.5
1,3,5-Trimethylbenzene	0.14	J	1.0	0.97	<SRL	U	1.0	1.05	0.5
1,2,4-Trimethylbenzene	0.47	J	1.0	0.97	0.29	J	1.0	1.05	0.5
Benzyl Chloride (a-Chlorotoluene)	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
1,3-Dichlorobenzene	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
1,4-Dichlorobenzene	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
1,2-Dichlorobenzene	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
1,2,4-Trichlorobenzene	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	0.5
Hexachlorobutadiene	<SRL	U	1.0	0.97	<SRL	U	1.0	1.05	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.


 Marcus Hueppe
 Laboratory Director



TO-15
QC
REPORT



Atmospheric Analysis & Consulting, Inc.

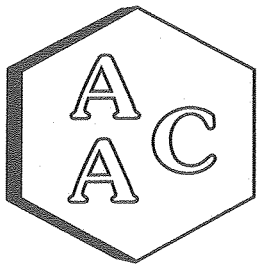
ANALYSIS DATE : 05/31/2013
 ANALYST : JJG

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	9.62	96
Chlorodifluoromethane	10.10	9.72	96
Propene	11.00	10.37	94
Dichlorodifluoromethane	9.80	9.67	99
Chloromethane	10.10	10.04	99
Dichlorotetrafluoroethane	10.10	10.14	100
Vinyl Chloride	10.20	9.66	95
Methanol	4.90	4.94	101
1,3-Butadiene	10.50	9.53	91
Bromomethane	10.20	8.75	86
Chloroethane	10.00	9.67	97
Dichlorofluoromethane	10.00	10.24	102
Ethanol	9.80	10.10	103
Vinyl Bromide	10.20	10.40	102
Acetone	10.80	9.57	89
Trichlorofluoromethane	10.10	10.69	106
2-Propanol (IPA)	11.00	10.12	92
Acrylonitrile	10.50	10.83	103
1,1-Dichloroethene	10.50	10.20	97
Methylene Chloride (DCM)	10.40	9.74	94
Allyl Chloride	11.00	10.88	99
Carbon Disulfide	10.50	9.63	92
Trichlorotrifluoroethane	10.40	10.25	99
trans-1,2-Dichloroethene	10.40	10.15	98
1,1-Dichloroethane	10.40	9.86	95
Methyl Tert Butyl Ether (MTBE)	10.60	10.57	100
Vinyl Acetate	9.70	9.68	100
2-Butanone (MEK)	10.60	10.85	102
cis-1,2-Dichloroethene	10.60	10.22	96
Hexane	10.70	9.70	91
Chloroform	10.60	10.56	100
Ethyl Acetate	11.00	11.00	100
Tetrahydrofuran	10.80	10.41	96
1,2-Dichloroethane	10.40	10.70	103
1,1,1-Trichloroethane	10.50	10.83	103





Atmospheric Analysis & Consulting, Inc.


ANALYSIS DATE : 05/31/2013
ANALYST : JJG

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

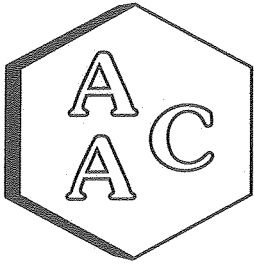
VOLATILE ORGANIC COMPOUNDS BY EPA METHOD TO-15
Continuing Calibration Verification of the 05/15/2013 Calibration

Compounds	Conc	Daily Conc	%REC*
Benzene	10.50	9.97	95
Carbon Tetrachloride	10.10	10.65	105
Cyclohexane	10.50	9.92	94
1,2-Dichloropropane	10.50	10.24	98
Bromodichloromethane	10.30	10.68	104
1,4-Dioxane	10.30	10.25	100
Trichloroethene (TCE)	10.30	10.38	101
2,2,4-Trimethylpentane	10.90	10.62	97
Heptane	10.70	10.55	99
cis-1,3-Dichloropropene	11.00	10.95	100
4-Methyl-2-pentanone (MiBK)	10.30	10.36	101
trans-1,3-Dichloropropene	9.80	9.96	102
1,1,2-Trichloroethane	10.60	10.83	102
Toluene	10.60	10.46	99
2-Hexanone (MBK)	10.80	10.89	101
Dibromochloromethane	11.00	11.62	106
1,2-Dibromoethane	10.40	10.42	100
Tetrachloroethene (PCE)	10.40	10.54	101
Chlorobenzene	10.60	10.05	95
Ethylbenzene	10.50	9.94	95
m & p-Xylenes	20.60	18.61	90
Bromoform	10.30	10.01	97
Styrene	10.40	9.76	94
1,1,2,2-Tetrachloroethane	10.60	9.68	91
o-Xylene	10.60	9.52	90
4-Ethyltoluene	10.40	9.88	95
1,3,5-Trimethylbenzene	10.20	9.40	92
1,2,4-Trimethylbenzene	10.20	9.80	96
Benzyl Chloride (a-Chlorotoluene)	10.00	10.07	101
1,3-Dichlorobenzene	10.00	9.57	96
1,4-Dichlorobenzene	10.00	9.31	93
1,2-Dichlorobenzene	10.00	9.41	94
1,2,4-Trichlorobenzene	9.30	9.03	97
Hexachlorobutadiene	9.80	9.40	96

* - %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 : 05/31/2013
AAC ID : LCS/LCSD DATE REPORTED : 05/31/2013
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	10.20	10.00	97	95	2.0
Methylene Chloride (DCM)	0.0	10.40	9.74	9.75	94	94	0.1
Benzene	0.0	10.50	9.97	9.80	95	93	1.7
Trichloroethene (TCE)	0.0	10.30	10.38	10.12	101	98	2.5
Toluene	0.0	10.60	10.46	10.34	99	98	1.2
Tetrachloroethene (PCE)	0.0	10.40	10.54	10.24	101	98	2.9
Chlorobenzene	0.0	10.60	10.05	10.37	95	98	3.1
Ethylbenzene	0.0	10.50	9.94	10.11	95	96	1.7
m & p-Xylenes	0.0	20.60	18.61	19.96	90	97	7.0
o-Xylene	0.0	10.60	9.52	9.87	90	93	3.6

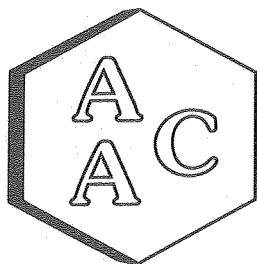
* Must be 70-130%

** Must be < 25%



Marcus Hueppe
Laboratory Director





Atmospheric Analysis & Consulting, Inc.

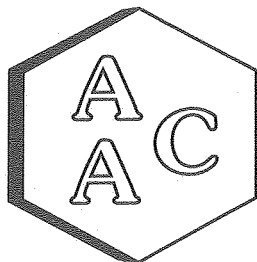
Method Blank Analysis Report

MATRIX : AIR **ANALYSIS DATE** : 05/31/2013
UNITS : ppbv **REPORT DATE** : 05/31/2013

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

<i>Client ID</i>	Method Blank	RL
<i>AAC ID</i>	MB 053113	
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 : 05/31/2013
 UNITS : ppbv REPORT DATE : 05/31/2013

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

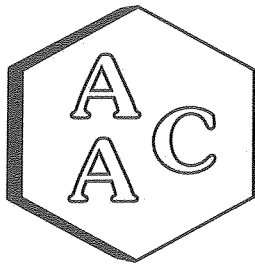
<i>Client ID</i> <i>AAC ID</i>	Method Blank MB 053113	RL
cis-1,3-Dichloropropene	<RL	0.5
4-Methyl-2-pentanone (MiBK)	<RL	0.5
trans-1,3-Dichloropropene	<RL	0.5
1,1,2-Trichloroethane	<RL	0.5
Toluene	<RL	0.5
2-Hexanone (MBK)	<RL	0.5
Dibromochloromethane	<RL	0.5
1,2-Dibromoethane	<RL	0.5
Tetrachloroethene (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.

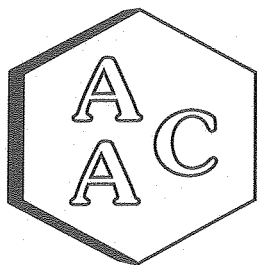
Quality Control/Quality Assurance Report

AAC ID	: 130647-63190	DATE ANALYZED	: 05/31/2013
MATRIX	: Air	DATE REPORTED	: 05/31/2013
		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	14.4	14.7	2.1
1,3-Butadiene	<SRL	<SRL	0.0
Bromomethane	<SRL	<SRL	0.0
Chloroethane	<SRL	<SRL	0.0
Dichlorofluoromethane	<SRL	<SRL	0.0
Ethanol	5.80	5.59	3.7
Vinyl Bromide	<SRL	<SRL	0.0
Acetone	5.12	5.02	2.0
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	<SRL	<SRL	0.0
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.

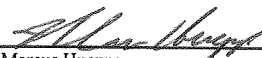
Quality Control/Quality Assurance Report

AAC ID : 130647-63190 **DATE ANALYZED** : 05/31/2013
MATRIX : Air **DATE REPORTED** : 05/31/2013
 UNITS : ppbv

TO-15 Duplicate Analysis

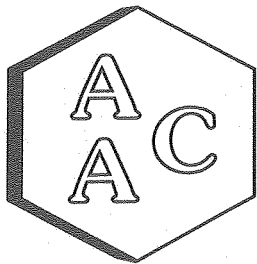
Compound	Sample Conc	Duplicate Conc	% RPD
Cyclohexane	<SRL	<SRL	0.0
1,2-Dichloropropane	<SRL	<SRL	0.0
Bromodichloromethane	<SRL	<SRL	0.0
1,4-Dioxane	<SRL	<SRL	0.0
Trichloroethene (TCE)	<SRL	<SRL	0.0
2,2,4-Trimethylpentane	<SRL	<SRL	0.0
Heptane	<SRL	<SRL	0.0
cis-1,3-Dichloropropene	<SRL	<SRL	0.0
4-Methyl-2-pentanone (MiBK)	<SRL	<SRL	0.0
trans-1,3-Dichloropropene	<SRL	<SRL	0.0
1,1,2-Trichloroethane	<SRL	<SRL	0.0
Toluene	<SRL	<SRL	0.0
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%	105%	2.8

SRL - Sample Reporting Limit



 Marcus Hueppe
 Laboratory Director





Atmospheric Analysis & Consulting, Inc.

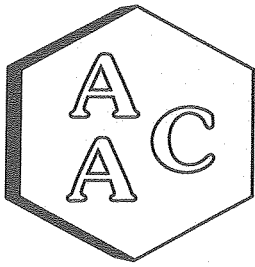
ANALYSIS DATE : 06/03/2013
 ANALYST : JJG

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	10.05	101
Chlorodifluoromethane	10.10	9.47	94
Propene	11.00	9.92	90
Dichlorodifluoromethane	9.80	9.57	98
Chloromethane	10.10	10.06	100
Dichlorotetrafluoroethane	10.10	10.14	100
Vinyl Chloride	10.20	10.07	99
Methanol	4.90	5.21	106
1,3-Butadiene	10.50	9.50	90
Bromomethane	10.20	8.28	81
Chloroethane	10.00	9.63	96
Dichlorofluoromethane	10.00	10.06	101
Ethanol	9.80	10.44	107
Vinyl Bromide	10.20	10.33	101
Acetone	10.80	9.50	88
Trichlorofluoromethane	10.10	10.68	106
2-Propanol (IPA)	11.00	10.27	93
Acrylonitrile	10.50	10.19	97
1,1-Dichloroethene	10.50	10.18	97
Methylene Chloride (DCM)	10.40	9.95	96
Allyl Chloride	11.00	11.09	101
Carbon Disulfide	10.50	9.76	93
Trichlorotrifluoroethane	10.40	10.15	98
trans-1,2-Dichloroethene	10.40	10.03	96
1,1-Dichloroethane	10.40	10.15	98
Methyl Tert Butyl Ether (MTBE)	10.60	10.78	102
Vinyl Acetate	9.70	10.02	103
2-Butanone (MEK)	10.60	10.87	103
cis-1,2-Dichloroethene	10.60	10.37	98
Hexane	10.70	9.94	93
Chloroform	10.60	10.65	100
Ethyl Acetate	11.00	11.09	101
Tetrahydrofuran	10.80	10.41	96
1,2-Dichloroethane	10.40	10.81	104
1,1,1-Trichloroethane	10.50	10.89	104





Atmospheric Analysis & Consulting, Inc.

ANALYSIS DATE : 06/03/2013
ANALYST : JJG

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

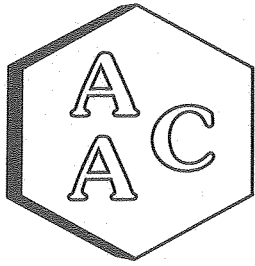
VOLATILE ORGANIC COMPOUNDS BY EPA METHOD TO-15
Continuing Calibration Verification of the 05/15/2013 Calibration

Compounds	Conc	Daily Conc	%REC*
Benzene	10.50	9.97	95
Carbon Tetrachloride	10.10	10.48	104
Cyclohexane	10.50	10.08	96
1,2-Dichloropropane	10.50	10.01	95
Bromodichloromethane	10.30	10.26	100
1,4-Dioxane	10.30	9.91	96
Trichloroethene (TCE)	10.30	10.18	99
2,2,4-Trimethylpentane	10.90	10.61	97
Heptane	10.70	10.20	95
cis-1,3-Dichloropropene	11.00	10.96	100
4-Methyl-2-pentanone (MiBK)	10.30	10.30	100
trans-1,3-Dichloropropene	9.80	9.83	100
1,1,2-Trichloroethane	10.60	10.46	99
Toluene	10.60	10.37	98
2-Hexanone (MBK)	10.80	10.33	96
Dibromochloromethane	11.00	11.43	104
1,2-Dibromoethane	10.40	10.33	99
Tetrachloroethene (PCE)	10.40	10.25	99
Chlorobenzene	10.60	10.52	99
Ethylbenzene	10.50	10.65	101
m & p-Xylenes	20.60	20.34	99
Bromoform	10.30	10.46	102
Styrene	10.40	10.11	97
1,1,2,2-Tetrachloroethane	10.60	10.31	97
o-Xylene	10.60	10.12	95
4-Ethyltoluene	10.40	10.62	102
1,3,5-Trimethylbenzene	10.20	9.91	97
1,2,4-Trimethylbenzene	10.20	10.36	102
Benzyl Chloride (a-Chlorotoluene)	10.00	10.98	110
1,3-Dichlorobenzene	10.00	10.24	102
1,4-Dichlorobenzene	10.00	9.74	97
1,2-Dichlorobenzene	10.00	9.86	99
1,2,4-Trichlorobenzene	9.30	9.46	102
Hexachlorobutadiene	9.80	9.67	99

* - %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 : 06/03/2013
AAC ID : LCS/LCSD DATE REPORTED : 06/03/2013
MEDIA : Air UNITS : ppbv

TO-15 Laboratory Control Spike Recovery

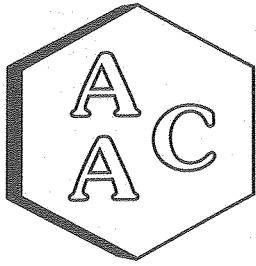
Compound	Sample Conc.	Spike Added	Spike Res	Dup Spike Res	Spike % Rec *	Spike Dup % Rec *	RPD** %
1,1-Dichloroethene	0.0	10.50	10.18	10.08	97	96	1.0
Methylene Chloride (DCM)	0.0	10.40	9.95	10.06	96	97	1.1
Benzene	0.0	10.50	9.97	9.78	95	93	1.9
Trichloroethene (TCE)	0.0	10.30	10.18	10.05	99	98	1.3
Toluene	0.0	10.60	10.37	9.99	98	94	3.7
Tetrachloroethene (PCE)	0.0	10.40	10.25	10.17	99	98	0.8
Chlorobenzene	0.0	10.60	10.52	10.31	99	97	2.0
Ethylbenzene	0.0	10.50	10.65	10.34	101	98	3.0
m & p-Xylenes	0.0	20.60	20.34	19.51	99	95	4.2
o-Xylene	0.0	10.60	10.12	9.92	95	94	2.0

* Must be 70-130%

** Must be < 25%


Marcus Hueppe
Laboratory Director





Atmospheric Analysis & Consulting, Inc.

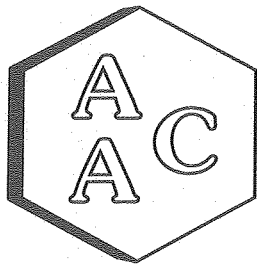
Method Blank Analysis Report

MATRIX : AIR ANALYSIS DATE : 06/03/2013
 UNITS : ppbv REPORT DATE : 06/03/2013

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

<i>Client ID</i> <i>AAC ID</i>	Method Blank MB 060313	RL
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 : 06/03/2013
UNITS : ppbv REPORT DATE : 06/03/2013

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

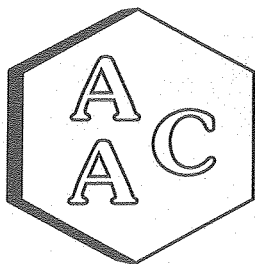
<i>Client ID</i> <i>AAC ID</i>	Method Blank MB 060313	RL
cis-1,3-Dichloropropene	<RL	0.5
4-Methyl-2-pentanone (MiBK)	<RL	0.5
trans-1,3-Dichloropropene	<RL	0.5
1,1,2-Trichloroethane	<RL	0.5
Toluene	<RL	0.5
2-Hexanone (MBK)	<RL	0.5
Dibromochloromethane	<RL	0.5
1,2-Dibromoethane	<RL	0.5
Tetrachloroethene (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	101%	--

RL - Reporting Limit



Marcus Hueppe
Laboratory Director





Atmospheric Analysis & Consulting, Inc.

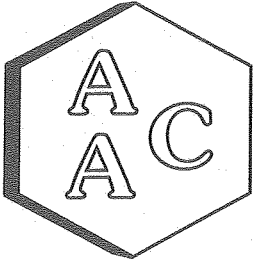
Quality Control/Quality Assurance Report

AAC ID : 130650-63200 DATE ANALYZED : 06/03/2013
MATRIX : Air DATE REPORTED : 06/03/2013
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	143	142	0.7
1,3-Butadiene	<SRL	<SRL	0.0
Bromomethane	<SRL	<SRL	0.0
Chloroethane	<SRL	<SRL	0.0
Dichlorofluoromethane	<SRL	<SRL	0.0
Ethanol	<SRL	<SRL	0.0
Vinyl Bromide	<SRL	<SRL	0.0
Acetone	<SRL	<SRL	0.0
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	<SRL	<SRL	0.0
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 : 130650-63200 DATE ANALYZED : 06/03/2013
MATRIX : Air DATE REPORTED : 06/03/2013
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	<SRL	<SRL	0.0
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%	103%	1.2
SRL - Sample Reporting Limit			


Marcus Hueppe
Laboratory Director



TO-15
RAW
DATA

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311314.D
 Acq On : 31 May 2013 18:53
 Operator : JJG
 Sample : 130650-63200 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 10 Sample Multiplier: 1

Quant Time: Jun 03 09:36:05 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	132947	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	772161	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	728009	10.00	ppbv	0.00

System Monitoring Compounds						
63) 4-Bromofluorobenzene (BFB)	22.710	174	474547	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	5754	0.22	ppbv #		92
3) Propene	4.799	42	7176	1.04	ppbv #		78
4) Dichlorodifluoromethane	4.908	85	15434	0.34	ppbv		97
5) Chloromethane	5.288	52	1144	0.27	ppbv #		1
6) Dichlorotetrafluoroethane	5.324	135	121	N.D.			
7) VinylChloride	0.000		0	N.D.			
8) Methanol	5.758	31	339308	Below Cal			
9) 1,3-Butadiene	5.867	54	260	N.D.			
10) Bromomethane	0.000		0	N.D.	dev		0.00
11) Chloroethane	0.000		0	N.D.	dev		0.00
12) Dichlorofluoromethane	0.000		0	N.D.	dev		0.00
13) Ethanol	7.043	45	61871	10.91	ppbv		
14) VinylBromide	0.000		0	N.D.			
15) Acetone	7.966	58	58139	8.07	ppbv		0.00
16) Trichlorofluoromethane	7.659	103	5142	0.19	ppbv #		89
17) 2-Propanol (IPA)	8.183	45	48117	1.95	ppbv		
18) Acrylonitrile	0.000		0	N.D.			
19) 1,1-Dichloroethene	0.000		0	N.D.			
20) MethyleneChloride (DCM)	0.000		0	N.D.	dev #		92
21) AllylChloride	0.000		0	N.D.	dev #		78
22) CarbonDisulfide	0.000		0	N.D.	dev		97
23) Trichlorotrifluoroethane	0.000		0	N.D.	dev #		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	0.000		0	N.D.	d		
28) 2-Butanone (MEK)	11.423	72	27106	3.68	ppbv		82
29) cis-1,2-Dichloroethene	0.000		0	N.D.			
30) Hexane	11.458	86	1347	0.38	ppbv #		28
31) Chloroform	12.510	83	870	N.D.			
32) EthylAcetate	12.047	43	10909	0.29	ppbv #		94

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311314.D
 Acq On : 31 May 2013 18:53
 Operator : JJG
 Sample : 130650-63200 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 10 Sample Multiplier: 1

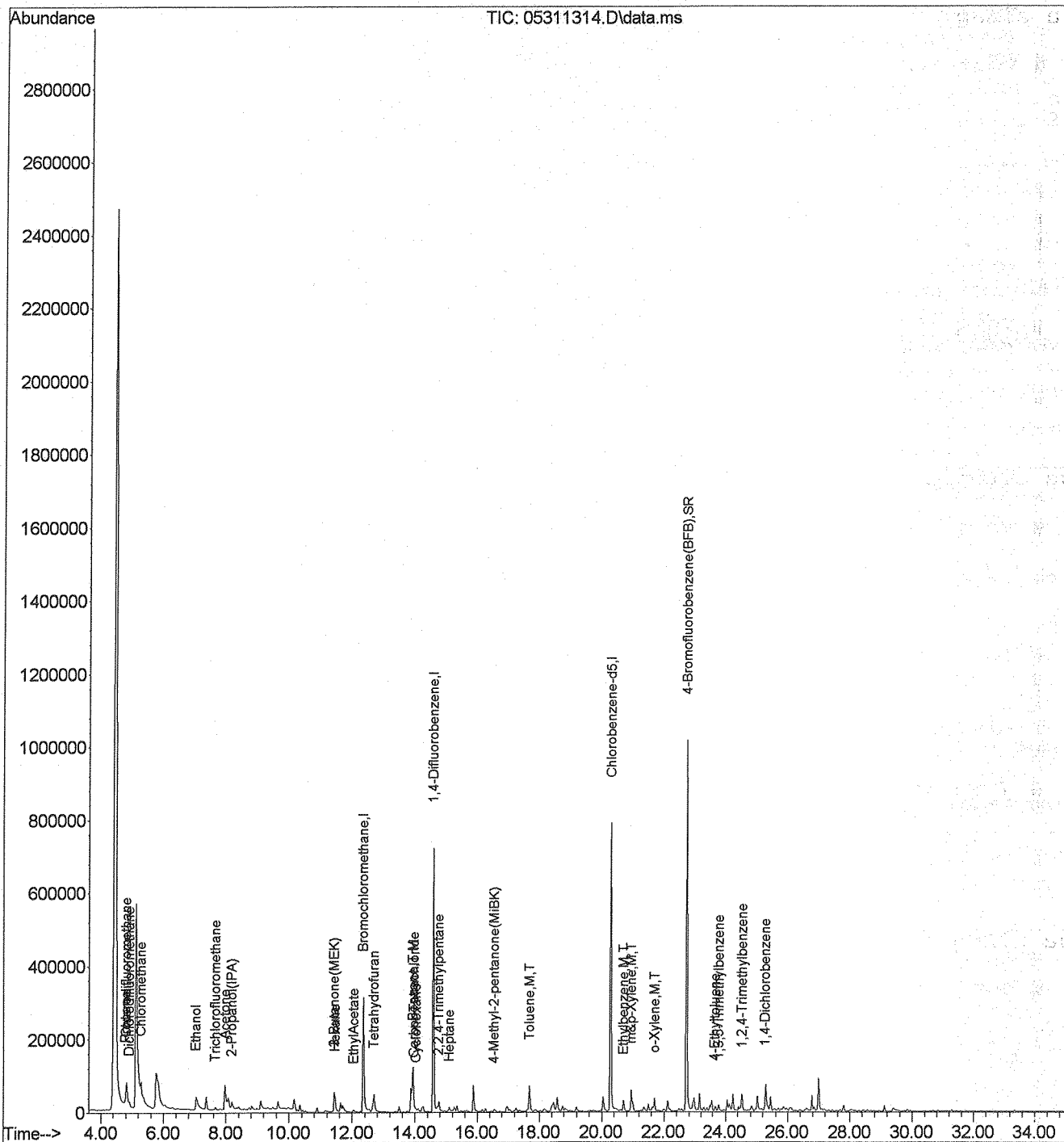
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 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
33) Tetrahydrofuran	12.689	72	13672	1.84	ppbv #	65
34) 1,2-Dichloroethane	13.616	62	124	N.D.		
35) 1,1,1-Trichloroethane	0.000		0	N.D.		
37) Benzene	13.937	78	126242	2.22	ppbv	99
38) CarbonTetrachloride	13.973	117	2528	0.05	ppbv #	97
39) Cyclohexane	14.026	69	597	0.07	ppbv #	1
40) 1,2-Dichloropropane	15.399	63	140	N.D.		
41) Bromodichloromethane	0.000		0	N.D.		
42) 1,4-Dioxane	0.000		0	N.D.	d	
43) Trichloroethene (TCE)	0.000		0	N.D.		
44) 2,2,4-Trimethylpentane	14.757	57	32849	0.32	ppbv #	92
45) Heptane	15.096	71	3016	0.17	ppbv #	46
46) cis-1,3-Dichloropropene	0.000		0	N.D.		
47) 4-Methyl-2-pentanone (M...)	16.540	58	2142	0.10	ppbv	79
48) trans-1,3-Dichloropropene	17.682	75	635	N.D.		
49) 1,1,2-Trichloroethane	0.000		0	N.D.	d	
50) Toluene	17.682	91	70041(m)	0.97	ppbv Dev (Min)	
51) 2-Hexanone (MBK)	0.000		0	N.D.	d	
52) Dibromochloromethane	19.019	129	270	N.D.		65
53) 1,2-Dibromoethane	0.000		0	N.D.		
54) Tetrachloroethene (PCE)	19.019	166	335	N.D.		
56) Chlorobenzene	20.285	114	143	N.D.		99
57) Ethylbenzene	20.695	91	31700	0.33	ppbv #	97
58) m&p-Xylene	20.945	106	33324	0.88	ppbv #	92
59) Bromoform	21.837	173	474	N.D.		
60) Styrene	21.658	104	1985	N.D.		
61) 1,1,2,2-Tetrachloroethane	0.000		0	N.D.		
62) o-Xylene	21.694	91	27460	0.36	ppbv	95
64) 4-Ethyltoluene	23.673	120	3609	0.11	ppbv #	94
65) 1,3,5-Trimethylbenzene	23.780	120	5334	0.12	ppbv #	93
66) 1,2,4-Trimethylbenzene	24.529	120	17181	0.38	ppbv #	95
67) BenzylChloride (a-Chlor...)	25.100	91	1066	N.D.		79
68) 1,3-Dichlorobenzene	0.000		0	N.D.		
69) 1,4-Dichlorobenzene	25.278	146	6233	0.09	ppbv #	90
70) 1,2-Dichlorobenzene	0.000		0	N.D.		
71) 1,2,4-Trichlorobenzene	29.451	180	274	N.D.		
72) Hexachlorobutadiene	0.000		0	N.D.		

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

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311314.D
 Acq On : 31 May 2013 18:53
 Operator : JJG
 Sample : 130650-63200 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 10 Sample Multiplier: 1

Quant Time: Jun 03 09:36:05 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\060313\
 Data File : 06031305.D
 Acq On : 3 Jun 2013 11:37
 Operator : JJG
 Sample : 130650-63200 x10
 Misc : IS/Surr: PS082712-02 + 50mL
 ALS Vial : 2 Sample Multiplier: 10

Quant Time: Jun 03 12:38:11 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	145756	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	793394	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	753276	10.00	ppbv	0.00

System Monitoring Compounds						
63) 4-Bromofluorobenzene (BFB)	22.710	174	480401	10.19	ppbv	0.00

Spiked Amount 10.000 Recovery = 101.90%

Target Compounds	R.T.	QIon	Response	Conc	Units	Dev (Min)	Qvalue
2) Chlorodifluoromethane	4.835	51	566	N.D.			
3) Propene	4.817	42	841	N.D.			
4) Dichlorodifluoromethane	4.926	85	1412	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	5.867	31	383060	82.26	ppbv		
9) 1,3-Butadiene	0.000		0	N.D.			
10) Bromomethane	6.446	96	701	N.D.			
11) Chloroethane	0.000		0	N.D.			
12) Dichlorofluoromethane	0.000		0	N.D.			
13) Ethanol	7.152	45	5921	N.D.			
14) VinylBromide	0.000		0	N.D.			
15) Acetone	8.057	58	6425	N.D.			
16) Trichlorofluoromethane	7.658	103	339	N.D.			
17) 2-Propanol (IPA)	8.256	45	6108	N.D.			
18) Acrylonitrile	0.000		0	N.D.			
19) 1,1-Dichloroethene	0.000		0	N.D.			
20) MethyleneChloride (DCM)	9.341	84	1281	N.D.			
21) AllylChloride	0.000		0	N.D.			
22) CarbonDisulfide	9.504	76	3786	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.			
26) MethylTertButylEther (M...)	0.000		0	N.D.			
27) VinylAcetate	10.888	43	120	N.D.			
28) 2-Butanone (MEK)	11.512	72	1776	N.D.			
29) cis-1,2-Dichloroethene	0.000		0	N.D.			
30) Hexane	0.000		0	N.D.			
31) Chloroform	0.000		0	N.D.			
32) EthylAcetate	12.118	43	1289	N.D.			

JJG
06/03/13

Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031305.D
 Acq On : 3 Jun 2013 11:37
 Operator : JJG
 Sample : 130650-63200 x10
 Misc : IS/Surr: PS082712-02 + 50mL
 ALS Vial : 2 Sample Multiplier: 10

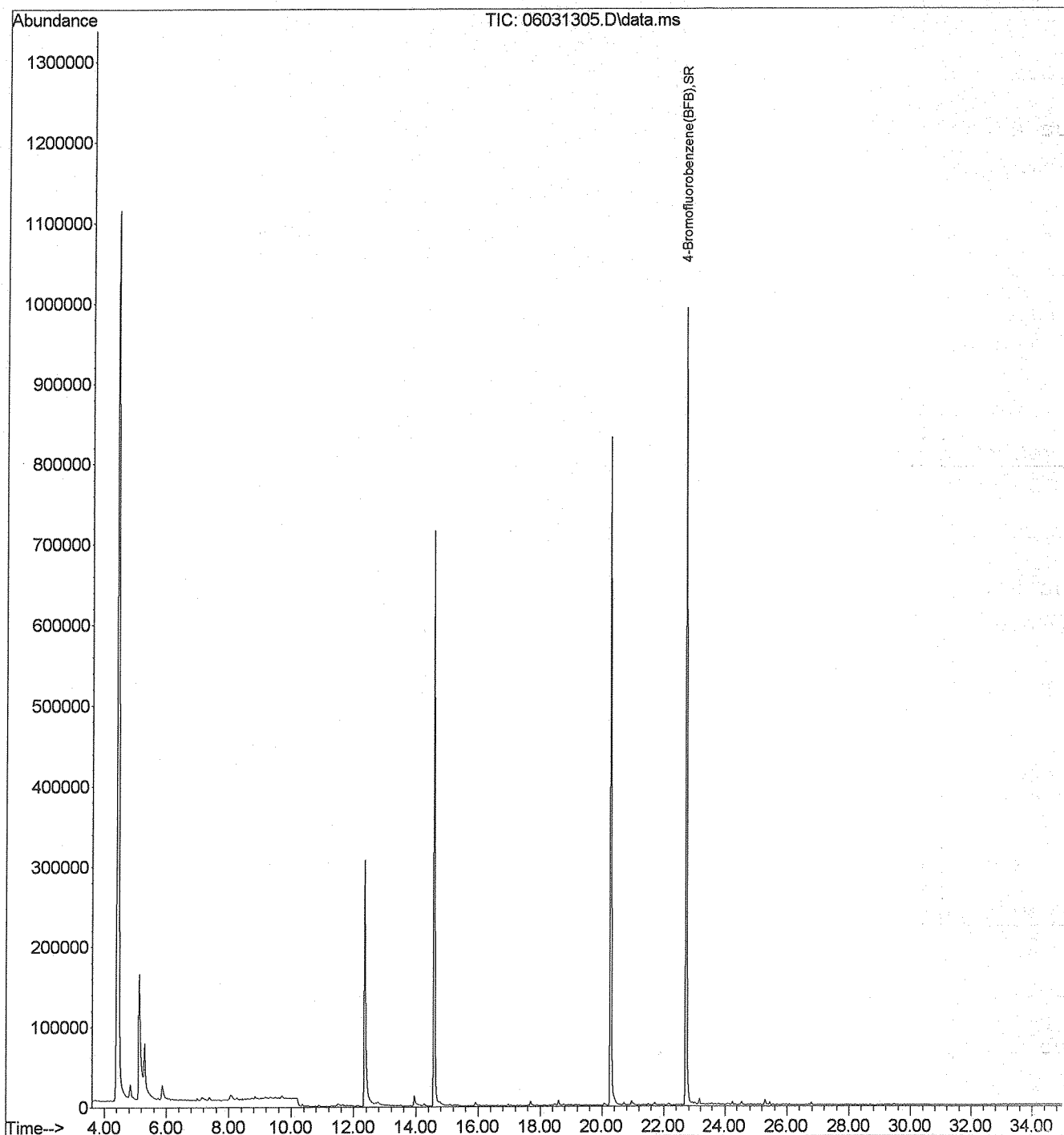
Quant Time: Jun 03 12:38:11 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
33) Tetrahydrofuran	12.760	72	1372		N.D.	
34) 1,2-Dichloroethane	0.000		0		N.D.	
35) 1,1,1-Trichloroethane	0.000		0		N.D.	
37) Benzene	13.937	78	15871		N.D.	
38) CarbonTetrachloride	0.000		0		N.D.	
39) Cyclohexane	0.000		0		N.D.	
40) 1,2-Dichloropropane	15.275	63	113		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	14.775	57	4185		N.D.	
45) Heptane	15.114	71	125		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	17.878	97	295		N.D.	
50) Toluene	17.700	91	7109		N.D.	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	20.267	114	234		N.D.	
57) Ethylbenzene	20.713	91	3358		N.D.	
58) m&p-Xylene	20.963	106	3210		N.D.	
59) Bromoform	0.000		0		N.D.	
60) Styrene	21.694	104	109		N.D.	
61) 1,1,2,2-Tetrachloroethane	0.000		0		N.D.	
62) o-Xylene	21.712	91	2881		N.D.	
64) 4-Ethyltoluene	23.691	120	354		N.D.	
65) 1,3,5-Trimethylbenzene	23.798	120	657		N.D.	
66) 1,2,4-Trimethylbenzene	24.547	120	1686		N.D.	
67) BenzylChloride (a-Chlor...)	25.225	91	110		N.D.	
68) 1,3-Dichlorobenzene	25.064	146	372		N.D.	
69) 1,4-Dichlorobenzene	25.296	146	1229		N.D.	
70) 1,2-Dichlorobenzene	25.849	146	353		N.D.	Dev (min)
71) 1,2,4-Trichlorobenzene	29.469	180	1019		N.D.	
72) Hexachlorobutadiene	30.075	225	358		N.D.	

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

Data Path : C:\msdchem\1\MS03\2013\060313\
Data File : 06031305.D
Acq On : 3 Jun 2013 11:37
Operator : JJG
Sample : 130650-63200 x10
Misc : IS/Surr: PS082712-02 + 50mL
ALS Vial : 2 Sample Multiplier: 10

Quant Time: Jun 03 12:38:11 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\053113\
 Data File : 05311315.D
 Acq On : 31 May 2013 19:41
 Operator : JJG
 Sample : 130650-63209 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Jun 03 09:38:53 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	139879	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	760226	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	730853	10.00	ppbv	0.00

System Monitoring Compounds						
63) 4-Bromofluorobenzene (BFB)	22.711	174	478599	10.46	ppbv	0.00

Spiked Amount 10.000 Recovery = 104.60%

Target Compounds	R.T.	QIon	Response	Conc	Units	Dev (Min)	Qvalue
2) Chlorodifluoromethane	4.817	51	4274	0.15	ppbv	#	93
3) Propene	4.781	42	8895	1.22	ppbv		91
4) Dichlorodifluoromethane	4.908	85	12519	0.26	ppbv		98
5) Chloromethane	5.288	52	1276	0.28	ppbv	#	24
6) Dichlorotetrafluoroethane	0.000		0	N.D.			
7) VinylChloride	0.000		0	N.D.			
8) Methanol	5.740	31	563474m	Below Cal			
9) 1,3-Butadiene	5.722	54	113	N.D.			
10) Bromomethane	0.000		0	N.D.	d		0.00
11) Chloroethane	0.000		0	N.D.	ppbv		0.00
12) Dichlorofluoromethane	0.000		0	N.D.	ppbv		0.00
13) Ethanol	7.025	45	100556m	16.86	ppbv		
14) VinylBromide	0.000		0	N.D.			
15) Acetone	7.948	58	86664m	11.44	ppbv		0.00
16) Trichlorofluoromethane	7.659	103	3717	0.13	ppbv		97
17) 2-Propanol (IPA)	8.183	45	69481m	2.68	ppbv		98
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	#	93
21) AllylChloride	9.251	39	404	N.D.	ppbv		91
22) CarbonDisulfide	0.000		0	N.D.	d		98
23) Trichlorotrifluoroethane	0.000		0	N.D.	d	#	24
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	0.000		0	N.D.	d		
28) 2-Butanone (MEK)	11.423	72	42666	5.51	ppbv		89
29) cis-1,2-Dichloroethene	0.000		0	N.D.			
30) Hexane	0.000		0	N.D.	d		00
31) Chloroform	12.511	83	234	N.D.			
32) EthylAcetate	12.047	43	21125	0.54	ppbv		96

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311315.D
 Acq On : 31 May 2013 19:41
 Operator : JJG
 Sample : 130650-63209 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 11 Sample Multiplier: 1

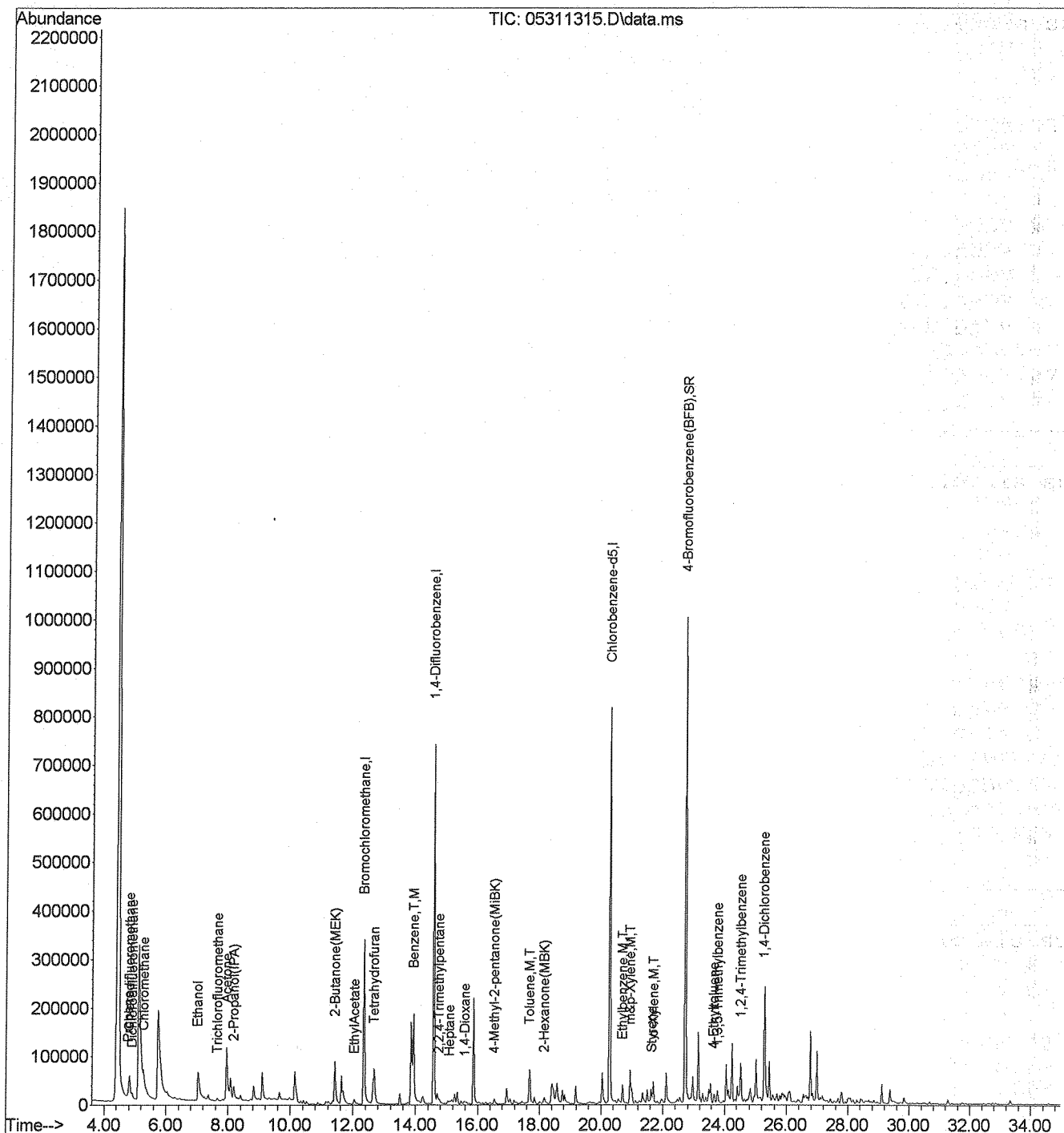
Quant Time: Jun 03 09:38:53 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
33) Tetrahydrofuran	12.671	72	22908	2.93	ppbv #	76
34) 1,2-Dichloroethane	13.616	62	117	N.D.		
35) 1,1,1-Trichloroethane	0.000		0	N.D.		
37) Benzene	13.937	78	194352	3.47	ppbv	99
38) CarbonTetrachloride	0.000		0	N.D.	d	
39) Cyclohexane	0.000		0	N.D.		
40) 1,2-Dichloropropane	15.346	63	564	N.D.		
41) Bromodichloromethane	0.000		0	N.D.		
42) 1,4-Dioxane	15.596	88	2718	0.21	ppbv	88
43) Trichloroethene (TCE)	0.000		0	N.D.		
44) 2,2,4-Trimethylpentane	14.758	57	8416	0.08	ppbv #	69
45) Heptane	15.096	71	1213	0.07	ppbv #	1
46) cis-1,3-Dichloropropene	0.000		0	N.D.		
47) 4-Methyl-2-pentanone (M...)	16.541	58	2890	0.14	ppbv #	78
48) trans-1,3-Dichloropropene	17.700	75	827	N.D.		
49) 1,1,2-Trichloroethane	0.000		0	N.D.	d	
50) Toluene	17.682	91	65825	0.93	ppbv	98
51) 2-Hexanone (MBK)	18.163	58	3181	0.13	ppbv #	71
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	143	N.D.		
57) Ethylbenzene	20.696	91	42704	0.44	ppbv #	96
58) m&p-Xylene	20.945	106	39478	1.04	ppbv #	89
59) Bromoform	21.837	173	1870	N.D.		
60) Styrene	21.659	104	3560	0.06	ppbv #	85
61) 1,1,2,2-Tetrachloroethane	0.000		0	N.D.		
62) o-Xylene	21.694	91	35323	0.46	ppbv	97
64) 4-Ethyltoluene	23.673	120	4955	0.16	ppbv #	94
65) 1,3,5-Trimethylbenzene	23.780	120	8281	0.18	ppbv #	91
66) 1,2,4-Trimethylbenzene	24.529	120	29728	0.66	ppbv	96
67) BenzylChloride (a-Chlor...)	25.118	91	2480	N.D.		
68) 1,3-Dichlorobenzene	25.064	146	571	N.D.		
69) 1,4-Dichlorobenzene	25.278	146	27148	0.38	ppbv	97
70) 1,2-Dichlorobenzene	25.849	146	580	N.D.		
71) 1,2,4-Trichlorobenzene	29.451	180	919	N.D.		
72) Hexachlorobutadiene	0.000		0	N.D.		

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

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311315.D
 Acq On : 31 May 2013 19:41
 Operator : JJG
 Sample : 130650-63209 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Jun 03 09:38:53 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



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Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031307.D
 Acq On : 3 Jun 2013 13:10
 Operator : JJG
 Sample : 130650-63209 x10
 Misc : IS/Surr: PS082712-02 + 50mL
 ALS Vial : 3 Sample Multiplier: 10

Quant Time: Jun 03 13:43:10 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	146559	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	803893	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	733552	10.00	ppbv	0.00

System Monitoring Compounds						
63) 4-Bromofluorobenzene (BFB)	22.710	174	475994	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	4.835	51	587	N.D.			
3) Propene	4.817	42	1136	N.D.			
4) Dichlorodifluoromethane	4.908	85	1251	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	5.849	31	652550	142.21	ppbv		
9) 1,3-Butadiene	0.000		0	N.D.			
10) Bromomethane	6.446	96	579	N.D.			0.00
11) Chloroethane	6.627	66	117	N.D.			0.00
12) Dichlorofluoromethane	0.000		0	N.D.			0.00
13) Ethanol	7.116	45	10595	16.95	ppbv		85
14) VinylBromide	0.000		0	N.D.			
15) Acetone	8.039	58	9542	12.02	ppbv		0.91
16) Trichlorofluoromethane	7.659	103	311	N.D.			
17) 2-Propanol (IPA)	8.238	45	9200	N.D.			103.70%
18) Acrylonitrile	0.000		0	N.D.			
19) 1,1-Dichloroethene	0.000		0	N.D.			Qvalue
20) MethyleneChloride (DCM)	9.323	84	1398	N.D.			
21) AllylChloride	0.000		0	N.D.			
22) CarbonDisulfide	9.486	76	3220	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.			
27) VinylAcetate	0.000		0	N.D.			
28) 2-Butanone (MEK)	11.494	72	3134	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.			85
32) EthylAcetate	12.100	43	2246	N.D.			0.91

Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031307.D
 Acq On : 3 Jun 2013 13:10
 Operator : JJG
 Sample : 130650-63209 x10
 Misc : IS/Surr: PS082712-02 + 50mL
 ALS Vial : 3 Sample Multiplier: 10

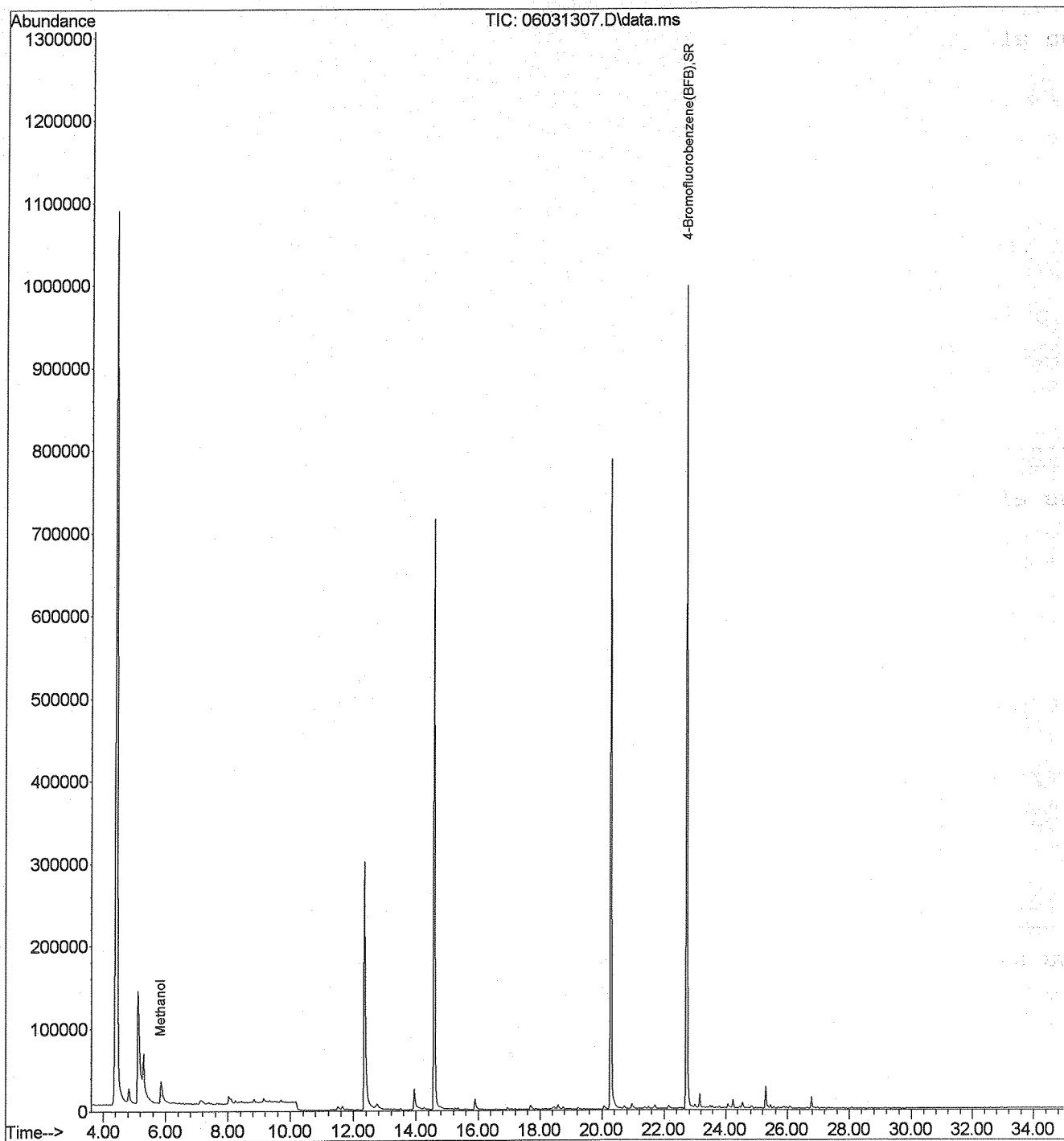
Quant Time: Jun 03 13:43:10 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) Tetrahydrofuran	12.742	72	2193		N.D.	
34) 1,2-Dichloroethane	0.000		0		N.D.	
35) 1,1,1-Trichloroethane	0.000		0		N.D.	
37) Benzene	13.937	78	23831		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	14.757	57	1665		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	17.860	97	1301		N.D.	
50) Toluene	17.700	91	6875		N.D.	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	20.374	114	111		N.D.	
57) Ethylbenzene	20.713	91	4297		N.D.	
58) m&p-Xylene	20.963	106	3844		N.D.	
59) Bromoform	21.836	173	380		N.D.	
60) Styrene	0.000		0		N.D.	
61) 1,1,2,2-Tetrachloroethane	0.000		0		N.D.	
62) o-Xylene	21.694	91	3373		N.D.	
64) 4-Ethyltoluene	23.691	120	325		N.D.	
65) 1,3,5-Trimethylbenzene	23.798	120	632		N.D.	
66) 1,2,4-Trimethylbenzene	24.547	120	2861		N.D.	
67) BenzylChloride (a-Chlor...)	25.296	91	4680		N.D.	
68) 1,3-Dichlorobenzene	0.000		0		N.D.	
69) 1,4-Dichlorobenzene	25.296	146	2612		N.D.	
70) 1,2-Dichlorobenzene	25.866	146	128		N.D.	
71) 1,2,4-Trichlorobenzene	29.469	180	278		N.D.	
72) Hexachlorobutadiene	0.000		0		N.D.	

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

Data Path : C:\msdchem\1\MS03\2013\060313\
Data File : 06031307.D
Acq On : 3 Jun 2013 13:10
Operator : JJG
Sample : 130650-63209 x10
Misc : IS/Surr: PS082712-02 + 50mL
ALS Vial : 3 Sample Multiplier: 10

Quant Time: Jun 03 13:43:10 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\053113\
 Data File : 05311316.D
 Acq On : 31 May 2013 20:29
 Operator : JJG
 Sample : 130650-63218 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 12 Sample Multiplier: 1

Quant Time: Jun 03 09:42:02 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	132307	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	779099	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	720945	10.00	ppbv	0.00

System Monitoring Compounds						
63) 4-Bromofluorobenzene(BFB)	22.710	174	473346	10.49	ppbv	0.00

Spiked Amount 10.000 Recovery = 104.90%

Target Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)	Qvalue
2) Chlorodifluoromethane	4.835	51	5572	0.21	ppbv	#	94
3) Propene	4.799	42	2239	0.33	ppbv	#	45
4) Dichlorodifluoromethane	4.908	85	15344	0.34	ppbv	#	97
5) Chloromethane	5.288	52	1245	0.29	ppbv	#	1
6) Dichlorotetrafluoroethane	5.324	135	262	N.D.			
7) VinylChloride	0.000		0	N.D.			
8) Methanol	5.849	31	27478	6.46	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.	ppbv		0.00
12) Dichlorofluoromethane	0.000		0	N.D.	ppbv		0.00
13) Ethanol	7.116	45	12277	2.18	ppbv		
14) VinylBromide	0.000		0	N.D.			
15) Acetone	8.020	58	16187	2.26	ppbv		0.00
16) Trichlorofluoromethane	7.658	103	4241	0.16	ppbv		96
17) 2-Propanol(IPA)	8.238	45	6228	0.25	ppbv	#	7
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	#	94
21) AllylChloride	0.000		0	N.D.	d	#	45
22) CarbonDisulfide	0.000		0	N.D.	d	#	97
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.			
27) VinylAcetate	0.000		0	N.D.	d		
28) 2-Butanone(MEK)	11.476	72	2222	0.30	ppbv	#	1
29) cis-1,2-Dichloroethene	0.000		0	N.D.			0.00
30) Hexane	11.458	86	1475	0.42	ppbv	#	74
31) Chloroform	12.492	83	800	N.D.			
32) EthylAcetate	0.000		0	N.D.	d		

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311316.D
 Acq On : 31 May 2013 20:29
 Operator : JJG
 Sample : 130650-63218 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 12 Sample Multiplier: 1

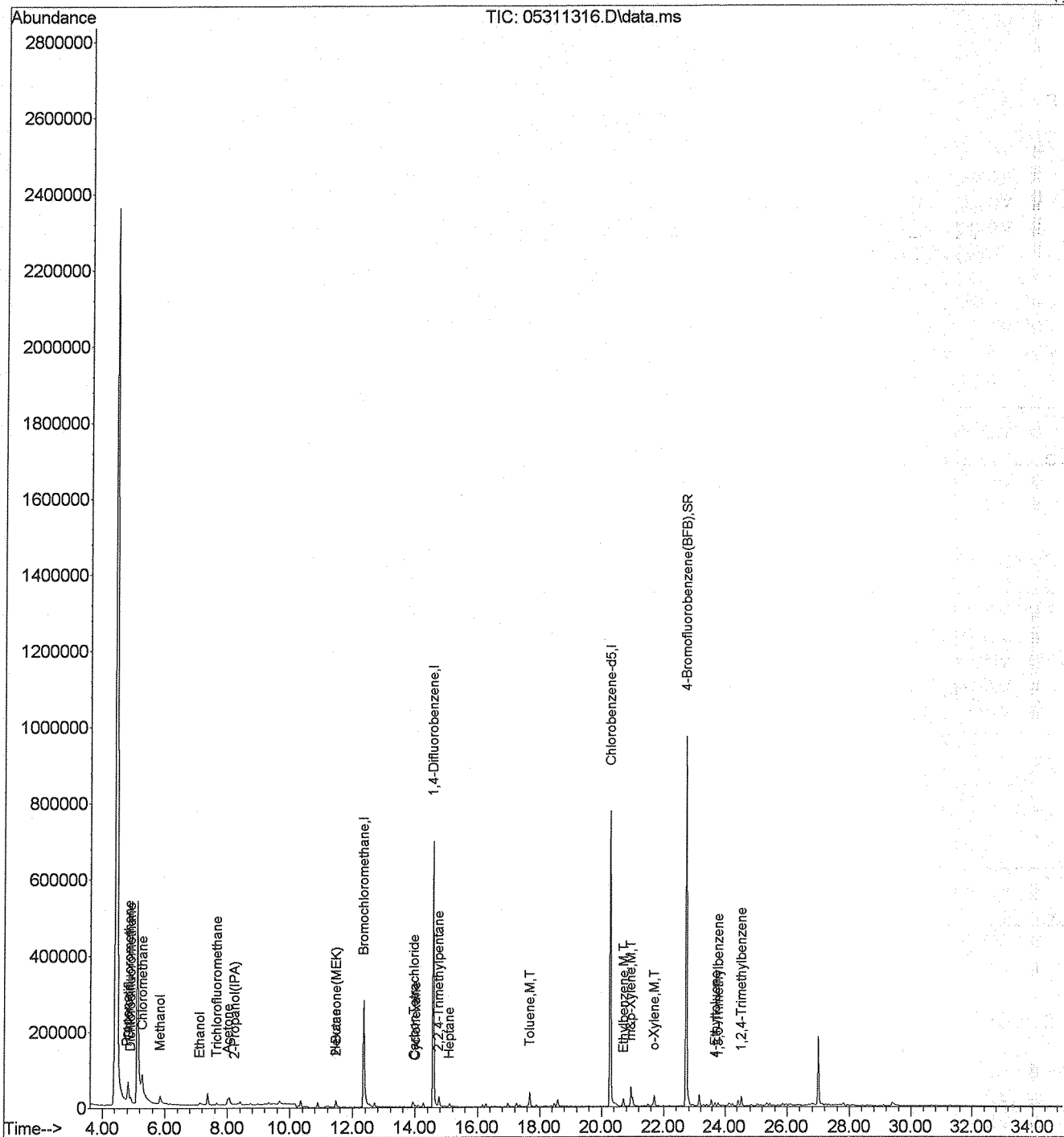
Quant Time: Jun 03 09:42:02 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
33) Tetrahydrofuran	0.000		0		N.D.	
34) 1,2-Dichloroethane	13.616	62	120		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	2506	0.05	ppbv #	97
39) Cyclohexane	14.008	69	536	0.06	ppbv #	48
40) 1,2-Dichloropropane	15.381	63	255		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	34722	0.34	ppbv	98
45) Heptane	15.096	71	2340	0.13	ppbv #	66
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	17.682	75	279		N.D.	
49) 1,1,2-Trichloroethane	0.000		0		N.D.	
50) Toluene	17.682	91	40700m	0.56	ppbv	Dev (Min)
51) 2-Hexanone (MBK)	0.000		0		N.D.	
52) Dibromochloromethane	19.019	129	285		N.D.	
53) 1,2-Dibromoethane	0.000		0		N.D.	
54) Tetrachloroethene (PCE)	19.019	166	247		N.D.	
56) Chlorobenzene	0.000		0		N.D.	
57) Ethylbenzene	20.713	91	23304	0.24	ppbv #	99
58) m&p-Xylene	20.945	106	32367	0.86	ppbv #	93
59) Bromoform	0.000		0		N.D.	
60) Styrene	21.676	104	946		N.D.	
61) 1,1,2,2-Tetrachloroethane	0.000		0		N.D.	
62) o-Xylene	21.694	91	26065	0.34	ppbv #	96
64) 4-Ethyltoluene	23.691	120	2367	0.08	ppbv #	80
65) 1,3,5-Trimethylbenzene	23.780	120	3337	0.07	ppbv #	96
66) 1,2,4-Trimethylbenzene	24.529	120	10689	0.24	ppbv #	95
67) BenzylChloride (a-Chlor...)	25.189	91	107		N.D.	
68) 1,3-Dichlorobenzene	0.000		0		N.D.	
69) 1,4-Dichlorobenzene	25.296	146	178		N.D.	
70) 1,2-Dichlorobenzene	0.000		0		N.D.	
71) 1,2,4-Trichlorobenzene	29.468	180	113		N.D.	
72) Hexachlorobutadiene	0.000		0		N.D.	

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

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311316.D
 Acq On : 31 May 2013 20:29
 Operator : JJG
 Sample : 130650-63218 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 12 Sample Multiplier: 1

Quant Time: Jun 03 09:42:02 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



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Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311317.D
 Acq On : 31 May 2013 21:17
 Operator : JJG
 Sample : 130650-63227 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Jun 03 09:49:03 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	133480	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	775162	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	723278	10.00	ppbv	0.00

System Monitoring Compounds						
63) 4-Bromofluorobenzene (BFB)	22.710	174	474297	10.48	ppbv	0.00

Spiked Amount 10.000 Recovery = 104.80%

Target Compounds	R.T.	QIon	Response	Conc	Units	Dev (Min)	Qvalue
2) Chlorodifluoromethane	4.817	51	4862	0.18	ppbv	#	98
3) Propene	4.799	42	2174	0.31	ppbv	#	74
4) Dichlorodifluoromethane	4.908	85	14418	0.32	ppbv	#	99
5) Chloromethane	5.288	52	1238	0.29	ppbv	#	1
6) Dichlorotetrafluoroethane	5.324	135	249	N.D.			
7) VinylChloride	0.000		0	N.D.		Dev (Min)	
8) Methanol	5.867	31	27813	6.48	ppbv		
9) 1,3-Butadiene	0.000		0	N.D.			
10) Bromomethane	6.428	96	396	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	7.116	45	14505	2.55	ppbv		
14) VinylBromide	0.000		0	N.D.			
15) Acetone	8.002	58	19208	2.66	ppbv		0.00
16) Trichlorofluoromethane	7.659	103	4373	0.16	ppbv	#	94
17) 2-Propanol (IPA)	8.238	45	9367	0.38	ppbv		80%
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	#	98
21) AllylChloride	0.000		0	N.D.	d	#	74
22) CarbonDisulfide	0.000		0	N.D.	d	#	99
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)	11.494	72	1928	0.26	ppbv	#	58
29) cis-1,2-Dichloroethene	0.000		0	N.D.			0.00
30) Hexane	11.458	86	877	0.25	ppbv	#	32
31) Chloroform	12.493	83	553	N.D.			
32) EthylAcetate	0.000		0	N.D.	d		

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Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311317.D
 Acq On : 31 May 2013 21:17
 Operator : JJG
 Sample : 130650-63227 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 13 Sample Multiplier: 1

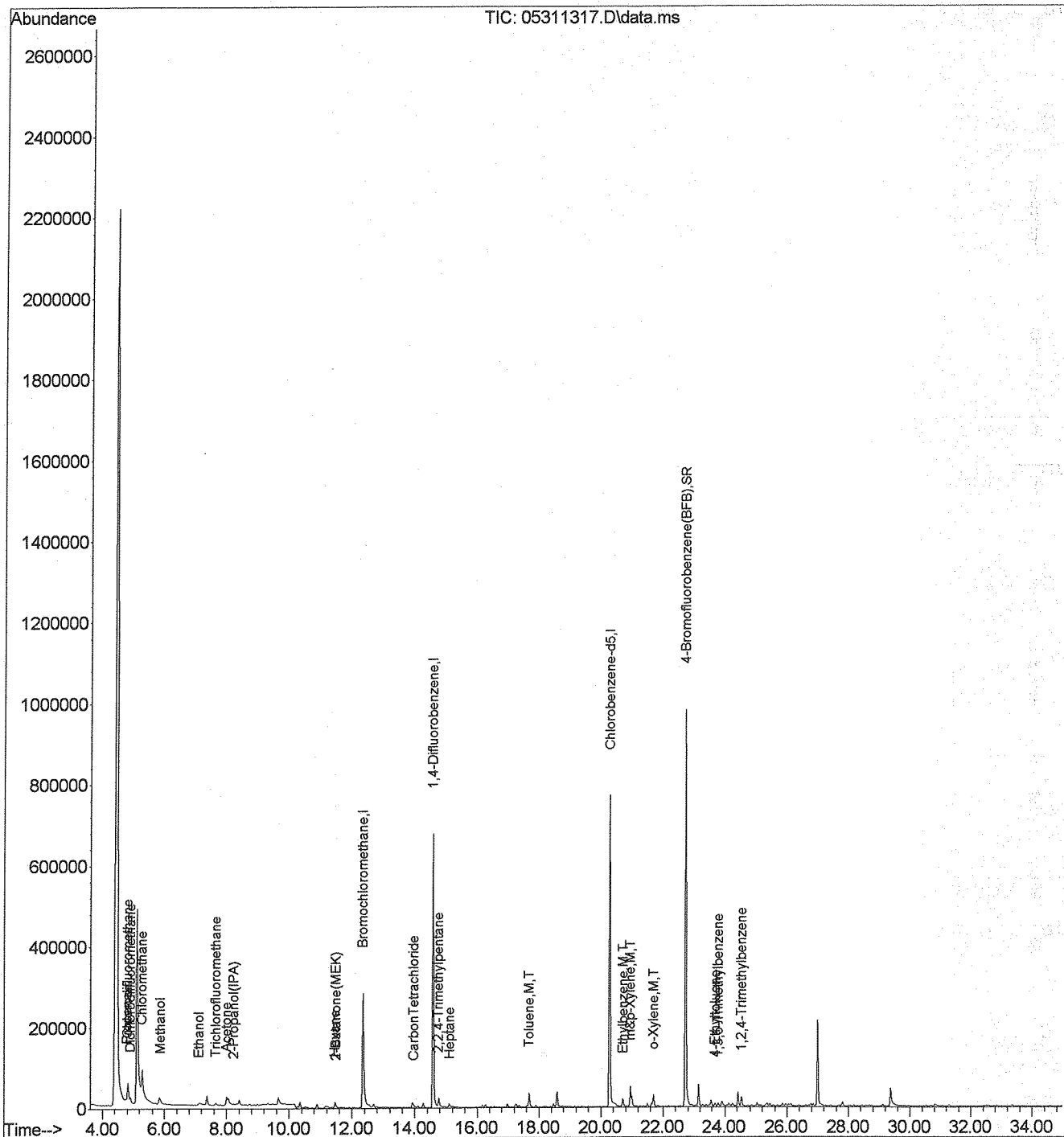
Quant Time: Jun 03 09:49:03 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
33) Tetrahydrofuran	0.000		0	N.D.		
34) 1,2-Dichloroethane	13.634	62	142	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	2429	0.05	ppbv #	94
39) Cyclohexane	14.008	69	279	N.D.		
40) 1,2-Dichloropropane	15.346	63	354	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	27391	0.27	ppbv	93
45) Heptane	15.114	71	2464	0.14	ppbv	94
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	17.682	75	291	N.D.		
49) 1,1,2-Trichloroethane	0.000		0	N.D.		
50) Toluene	17.682	91	37528	0.52	ppbv #	97
51) 2-Hexanone (MBK)	0.000		0	N.D.		
52) Dibromochloromethane	19.019	129	188	N.D.		
53) 1,2-Dibromoethane	0.000		0	N.D.		
54) Tetrachloroethene (PCE)	19.019	166	470	N.D.		
56) Chlorobenzene	20.285	114	129	N.D.		
57) Ethylbenzene	20.713	91	21480	0.23	ppbv #	99
58) m&p-Xylene	20.945	106	31629	0.84	ppbv	91
59) Bromoform	0.000		0	N.D.		
60) Styrene	21.694	104	1129	N.D.		
61) 1,1,2,2-Tetrachloroethane	0.000		0	N.D.		
62) o-Xylene	21.694	91	25314	0.33	ppbv	96
64) 4-Ethyltoluene	23.691	120	2234	0.07	ppbv #	88
65) 1,3,5-Trimethylbenzene	23.780	120	2999	0.07	ppbv #	89
66) 1,2,4-Trimethylbenzene	24.529	120	9495	0.21	ppbv #	94
67) BenzylChloride (a-Chlor...)	25.118	91	273	N.D.		
68) 1,3-Dichlorobenzene	0.000		0	N.D.		
69) 1,4-Dichlorobenzene	25.296	146	151	N.D.		
70) 1,2-Dichlorobenzene	0.000		0	N.D.		
71) 1,2,4-Trichlorobenzene	0.000		0	N.D.		
72) Hexachlorobutadiene	0.000		0	N.D.		

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

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311317.D
 Acq On : 31 May 2013 21:17
 Operator : JJG
 Sample : 130650-63227 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Jun 03 09:49:03 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



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Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311318.D
 Acq On : 31 May 2013 22:05
 Operator : JJG
 Sample : 130650-63236 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jun 03 09:51:26 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	136062	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	777462	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	738805	10.00	ppbv	0.00

System Monitoring Compounds						
63) 4-Bromofluorobenzene (BFB)	22.710	174	480469	10.39	ppbv	0.00

Spiked Amount 10.000 Recovery = 103.90%

Target Compounds	R.T.	QIon	Response	Conc	Units	Dev (Min)	Qvalue
2) Chlorodifluoromethane	4.818	51	5281	0.20	ppbv	#	96
3) Propene	4.781	42	4291	0.61	ppbv	#	69
4) Dichlorodifluoromethane	4.890	85	13727	0.30	ppbv	#	97
5) Chloromethane	5.306	52	1009	0.23	ppbv	#	6
6) Dichlorotetrafluoroethane	5.324	135	130	N.D.			
7) VinylChloride	0.000		0	N.D.			
8) Methanol	5.813	31	138931	35.13	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	7.080	45	25022	4.31	ppbv		
14) VinylBromide	0.000		0	N.D.			
15) Acetone	7.984	58	33371	4.53	ppbv		0.00
16) Trichlorofluoromethane	7.659	103	4265	0.16	ppbv	#	90
17) 2-Propanol (IPA)	8.202	45	25647	1.02	ppbv	#	90
18) Acrylonitrile	0.000		0	N.D.			
19) 1,1-Dichloroethene	0.000		0	N.D.			
20) MethyleneChloride (DCM)	0.000		0	N.D.	ppbv	#	96
21) AllylChloride	0.000		0	N.D.	ppbv	#	69
22) CarbonDisulfide	0.000		0	N.D.	ppbv	#	97
23) Trichlorotrifluoroethane	0.000		0	N.D.	ppbv	#	6
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)	11.441	72	9254	1.23	ppbv	#	59
29) cis-1,2-Dichloroethene	0.000		0	N.D.			
30) Hexane	11.458	86	1400	0.39	ppbv	#	17
31) Chloroform	12.493	83	866	N.D.			
32) EthylAcetate	0.000		0	N.D.	d		

JJG
Qvalue

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311318.D
 Acq On : 31 May 2013 22:05
 Operator : JJG
 Sample : 130650-63236 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 14 Sample Multiplier: 1

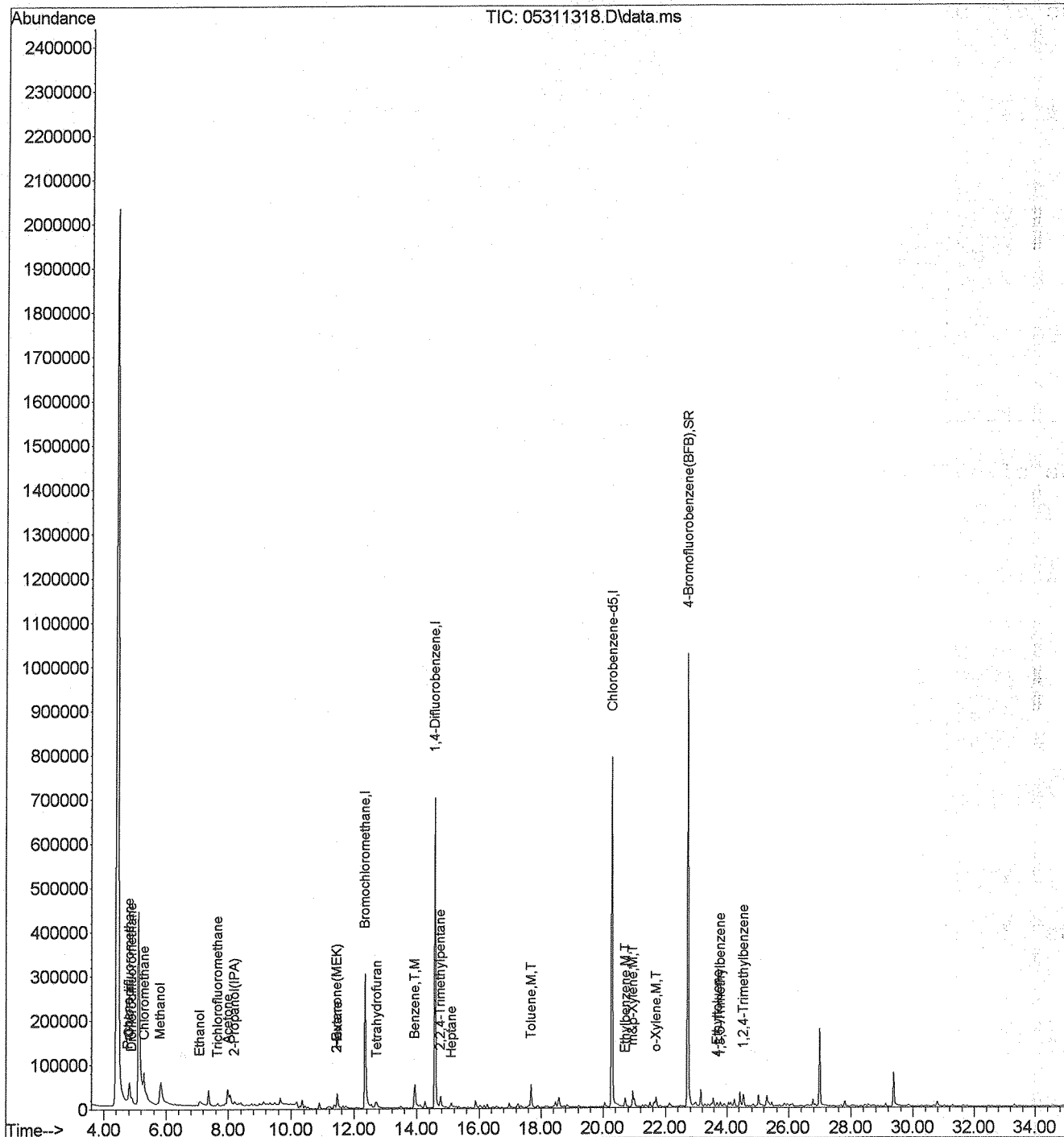
Quant Time: Jun 03 09:51:26 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) Tetrahydrofuran	12.725	72	3785	0.50	ppbv #	62
34) 1,2-Dichloroethane	0.000		0	N.D.		
35) 1,1,1-Trichloroethane	13.313	97	134	N.D.		
37) Benzene	13.937	78	46052	0.80	ppbv	99
38) CarbonTetrachloride	0.000		0	N.D.	d	
39) Cyclohexane	0.000		0	N.D.	d	
40) 1,2-Dichloropropane	15.328	63	295	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	32596	0.32	ppbv #	94
45) Heptane	15.114	71	2834	0.16	ppbv #	78
46) cis-1,3-Dichloropropene	0.000		0	N.D.		
47) 4-Methyl-2-pentanone (M...)	16.558	58	703	N.D.		
48) trans-1,3-Dichloropropene	17.664	75	388	N.D.		
49) 1,1,2-Trichloroethane	0.000		0	N.D.	d	
50) Toluene	17.682	91	52600	0.72	ppbv #	99
51) 2-Hexanone (MBK)	18.199	58	631	N.D.		
52) Dibromochloromethane	19.019	129	113	N.D.		
53) 1,2-Dibromoethane	0.000		0	N.D.		
54) Tetrachloroethene (PCE)	19.019	166	141	N.D.		
56) Chlorobenzene	20.285	114	112	N.D.		
57) Ethylbenzene	20.713	91	21044	0.22	ppbv #	97
58) m&p-Xylene	20.945	106	22001	0.57	ppbv	92
59) Bromoform	0.000		0	N.D.		
60) Styrene	21.658	104	1504	N.D.		
61) 1,1,2,2-Tetrachloroethane	0.000		0	N.D.		
62) o-Xylene	21.694	91	17449	0.22	ppbv #	97
64) 4-Ethyltoluene	23.691	120	2571	0.08	ppbv #	98
65) 1,3,5-Trimethylbenzene	23.780	120	3174	0.07	ppbv #	96
66) 1,2,4-Trimethylbenzene	24.529	120	11130	0.24	ppbv #	92
67) BenzylChloride (a-Chlor...)	25.189	91	121	N.D.		
68) 1,3-Dichlorobenzene	0.000		0	N.D.		
69) 1,4-Dichlorobenzene	0.000		0	N.D.	d	
70) 1,2-Dichlorobenzene	0.000		0	N.D.		
71) 1,2,4-Trichlorobenzene	29.451	180	131	N.D.		
72) Hexachlorobutadiene	0.000		0	N.D.		

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

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311318.D
 Acq On : 31 May 2013 22:05
 Operator : JJG
 Sample : 130650-63236 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jun 03 09:51:26 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



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Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311319.D
 Acq On : 31 May 2013 22:53
 Operator : JJG
 Sample : 130650-63245 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 15 Sample Multiplier: 1

Quant Time: Jun 03 09:53:05 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	137023	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	767232	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	731543	10.00	ppbv	0.00

System Monitoring Compounds						
63) 4-Bromofluorobenzene (BFB)	22.711	174	473330	10.34	ppbv	0.00

Spiked Amount 10.000 Recovery = 103.40%

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Chlorodifluoromethane	4.817	51	3995	0.15	ppbv	# 95
3) Propene	4.781	42	4503	0.63	ppbv	88
4) Dichlorodifluoromethane	4.890	85	12455	0.27	ppbv	98
5) Chloromethane	5.288	52	1109	0.25	ppbv	# 1
6) Dichlorotetrafluoroethane	0.000		0	N.D.		
7) VinylChloride	0.000		0	N.D.		Dev (Min)
8) Methanol	5.831	31	103282	25.00	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	7.079	45	17521	3.00	ppbv	
14) VinylBromide	0.000		0	N.D.		
15) Acetone	7.984	58	26093	3.52	ppbv	0.00
16) Trichlorofluoromethane	7.659	103	3431	0.13	ppbv	# 98
17) 2-Propanol (IPA)	8.201	45	19566	0.77	ppbv	100%
18) Acrylonitrile	9.052	52	114	N.D.		
19) 1,1-Dichloroethene	0.000		0	N.D.		Qvalue
20) MethyleneChloride (DCM)	0.000		0	N.D.	d	# 95
21) AllylChloride	9.233	39	256	N.D.		93
22) CarbonDisulfide	9.468	76	68254	1.47	ppbv	98
23) Trichlorotrifluoroethane	8.980	103	816	N.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.		
27) VinylAcetate	10.888	43	801	N.D.		
28) 2-Butanone (MEK)	11.459	72	7484	0.99	ppbv	# 66
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	267	N.D.		
32) EthylAcetate	0.000		0	N.D.	d	

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Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311319.D
 Acq On : 31 May 2013 22:53
 Operator : JJG
 Sample : 130650-63245 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 15 Sample Multiplier: 1

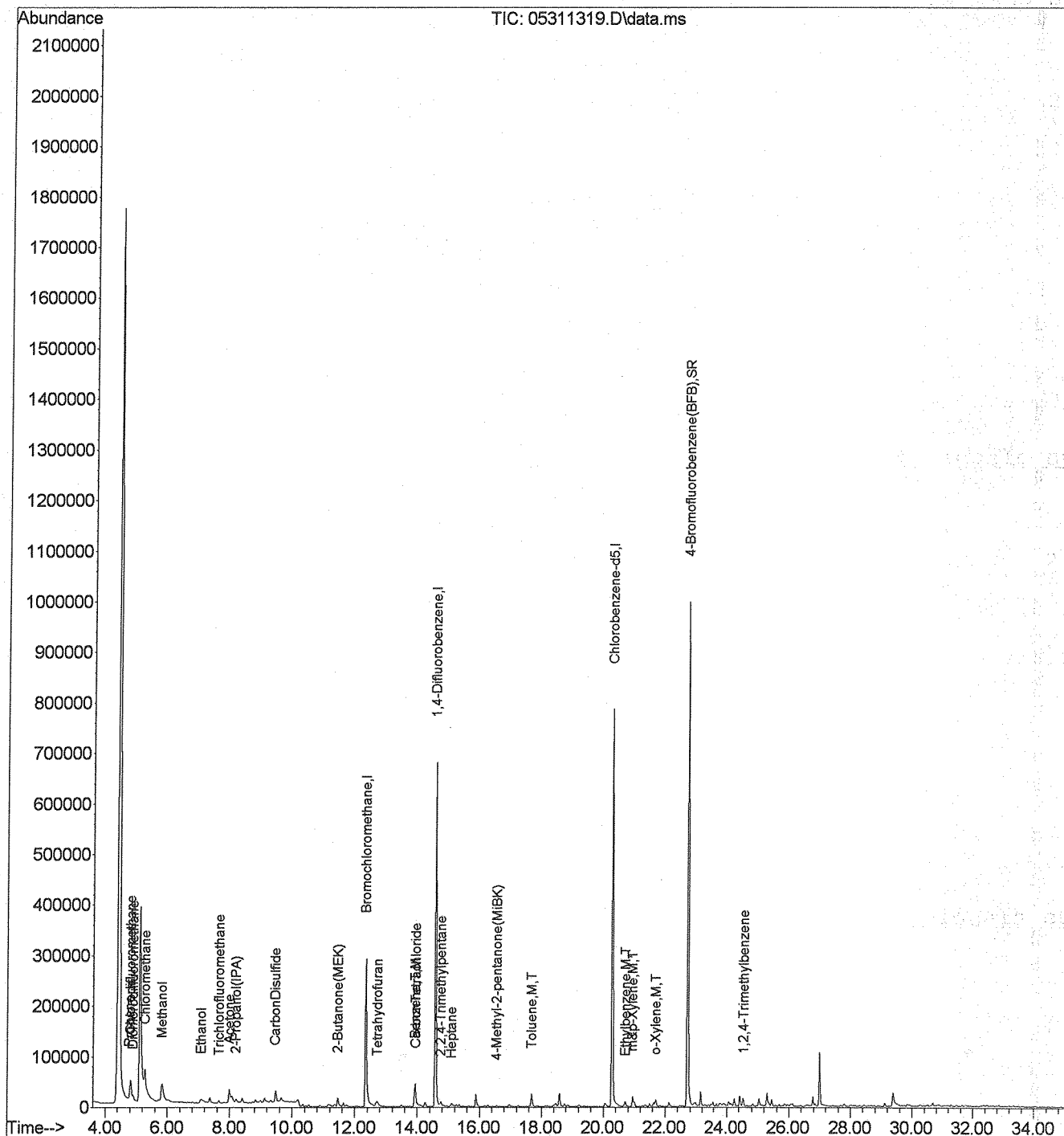
Quant Time: Jun 03 09:53:05 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
33) Tetrahydrofuran	12.725	72	3477	0.45	ppbv #	75
34) 1,2-Dichloroethane	13.598	62	113	N.D.		
35) 1,1,1-Trichloroethane	0.000		0	N.D.		
37) Benzene	13.937	78	41784	0.74	ppbv	98
38) CarbonTetrachloride	13.955	117	2113	0.05	ppbv #	96
39) Cyclohexane	0.000		0	N.D.		
40) 1,2-Dichloropropane	15.382	63	117	N.D.		
41) Bromodichloromethane	0.000		0	N.D.		
42) 1,4-Dioxane	15.667	88	126	N.D.		
43) Trichloroethene (TCE)	0.000		0	N.D.		
44) 2,2,4-Trimethylpentane	14.758	57	8049	0.08	ppbv #	90
45) Heptane	15.096	71	1477	0.08	ppbv #	67
46) cis-1,3-Dichloropropene	0.000		0	N.D.		
47) 4-Methyl-2-pentanone (M...)	16.559	58	867	0.04	ppbv #	81
48) trans-1,3-Dichloropropene	17.682	75	137	N.D.		
49) 1,1,2-Trichloroethane	0.000		0	N.D.	d	
50) Toluene	17.682	91	26159	0.36	ppbv #	97
51) 2-Hexanone (MBK)	18.199	58	108	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	112	N.D.		
57) Ethylbenzene	20.713	91	10892	0.11	ppbv #	98
58) m&p-Xylene	20.945	106	12304	0.32	ppbv #	84
59) Bromoform	0.000		0	N.D.		
60) Styrene	21.658	104	1109	N.D.		
61) 1,1,2,2-Tetrachloroethane	0.000		0	N.D.		
62) o-Xylene	21.694	91	10398	0.13	ppbv #	96
64) 4-Ethyltoluene	23.691	120	1195	N.D.		
65) 1,3,5-Trimethylbenzene	23.780	120	1816	N.D.		
66) 1,2,4-Trimethylbenzene	24.547	120	6156	0.14	ppbv	93
67) BenzylChloride (a-Chlor...)	25.189	91	108	N.D.		
68) 1,3-Dichlorobenzene	0.000		0	N.D.		
69) 1,4-Dichlorobenzene	25.278	146	2399	N.D.		
70) 1,2-Dichlorobenzene	0.000		0	N.D.		
71) 1,2,4-Trichlorobenzene	29.469	180	108	N.D.		
72) Hexachlorobutadiene	0.000		0	N.D.		

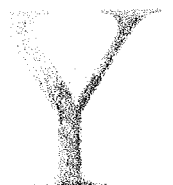
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\msdchem\1\MS03\2013\053113\
Data File : 05311319.D
Acq On : 31 May 2013 22:53
Operator : JJG
Sample : 130650-63245 x1
Misc : IS/Surr: PS082712-02 + 500mL
ALS Vial : 15 Sample Multiplier: 1

Quant Time: Jun 03 09:53:05 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

Sequence Name: C:\msdchem\1\sequence\2013\053113.S
Comment: GCMS-03
Operator: JJG
Data Path: C:\MSDCHEM\1\MS03\2013\053113\
Instrument Control Pre-Seq Cmd:
Data Analysis Pre-Seq Cmd:
Instrument Control Post-Seq Cmd:
Data Analysis Post-Seq Cmd:

Method Sections To Run On A Barcode Mismatch
(X) Full Method (X) Inject Anyway
() Reprocessing Only () Don't Inject

Line	Sample Name/Misc Info
1) Sample	1 05311301 TO15-5MS TO15 BFB 053113
2) Sample	1 05311302 TO15-5MS TO15 CCV 053113
3) Sample	1 05311303 TO15-5MS TO15 LCSD 053113
4) Sample	1 05311304 TO15-5MS TO15 MB 053113
5) Sample	2 05311305 TO15-5MS 130647-63190 x1
6) Sample	2 05311306 TO15-5MS 130647-63190 x1 dp
7) Sample	3 05311307 TO15-5MS 130647-63191 x1
8) Sample	4 05311308 TO15-5MS 130647-63192 x1
9) Sample	5 05311309 TO15-5MS 130647-63193 x1
10) Sample	6 05311310 TO15-5MS 130647-63194 x1
11) Sample	7 05311311 TO15-5MS 130647-63195 x1
12) Sample	8 05311312 TO15-5MS 130647-63196 x1
13) Sample	9 05311313 TO15-5MS 130647-63197 x1
14) Sample	10 05311314 TO15-5MS 130650-63200 x1
15) Sample	11 05311315 TO15-5MS 130650-63209 x1
16) Sample	12 05311316 TO15-5MS 130650-63218 x1
17) Sample	13 05311317 TO15-5MS 130650-63227 x1
18) Sample	14 05311318 TO15-5MS 130650-63236 x1
19) Sample	15 05311319 TO15-5MS 130650-63245 x1
20) Sample	16 05311320 TO15-5MS 130653-63265 x1
21) Sample	2 05311321 TO15-5MS 130653-63266 x1
22) Sample	3 05311322 TO15-5MS 130653-63267 x1
23) Sample	4 05311323 TO15-5MS 130653-63268 x1
24) Sample	5 05311324 TO15-5MS 130647-63193 x2
25) Sample	7 05311325 TO15-5MS 130647-63195 x5
26) Sample	6 05311326 TO15-5MS Flow Check#053113-01
27) Sample	1 05311327 TO15-5MS Can Check#000529

Comments: _____

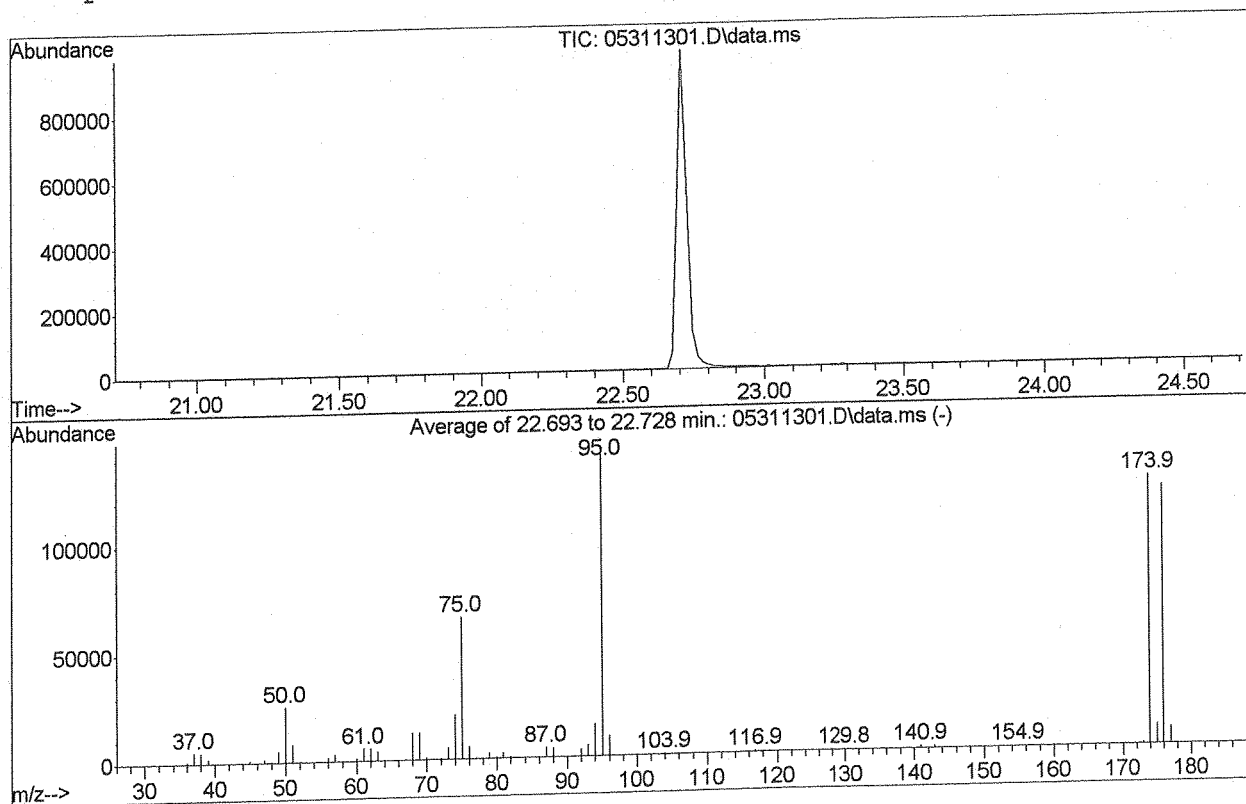
Analyst: *JJG*

Date: 06/03/13

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311301.D
 Acq On : 31 May 2013 8:32 am
 Operator : JJG
 Sample : TO15 BFB 053113
 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	18.1	25488	PASS
75	95	30	60	46.4	65440	PASS
95	95	100	100	100.0	141003	PASS
96	95	5	9	6.6	9311	PASS
173	174	0.00	2	0.9	1137	PASS
174	95	50	100	88.2	124371	PASS
175	174	5	9	7.3	9107	PASS
176	174	95	101	96.6	120152	PASS
177	176	5	9	6.8	8113	PASS

JJG
05/31/13

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311302.D
 Acq On : 31 May 2013 9:19
 Operator : JJG
 Sample : TO15 CCV 053113
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 31 15:58:54 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	146685	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	752301	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	756486	10.00	ppbv	0.00

System Monitoring Compounds						
63) 4-Bromofluorobenzene(BFB)	22.711	174	455206	9.62	ppbv	0.00

Spiked Amount 10.000 Recovery = 96.20%

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Chlorodifluoromethane	4.817	51	282455	9.72	ppbv	100
3) Propene	4.781	42	79110m	10.37	ppbv	
4) Dichlorodifluoromethane	4.908	85	485056	9.67	ppbv	99
5) Chloromethane	5.288	52	47233m	10.04	ppbv	
6) Dichlorotetrafluoroethane	5.324	135	346658	10.14	ppbv	89
7) VinylChloride	5.668	62	161122m	9.66	ppbv	
8) Methanol	5.849	31	23428m	4.94	ppbv	
9) 1,3-Butadiene	5.849	54	100966m	9.53	ppbv	
10) Bromomethane	6.446	96	105092m	8.75	ppbv	100
11) Chloroethane	6.736	66	26251	9.67	ppbv	98
12) Dichlorofluoromethane	7.007	67	367549m	10.24	ppbv	100
13) Ethanol	7.043	45	63181m	10.10	ppbv	
14) VinylBromide	7.261	108	146782m	10.40	ppbv	
15) Acetone	7.966	58	76006m	9.57	ppbv	100
16) Trichlorofluoromethane	7.677	103	313168	10.69	ppbv	99
17) 2-Propanol (IPA)	8.147	45	275368m	10.12	ppbv	
18) Acrylonitrile	8.962	52	128832m	10.83	ppbv	
19) 1,1-Dichloroethene	8.726	96	170377	10.20	ppbv	96
20) MethyleneChloride (DCM)	9.323	84	149386m	9.74	ppbv	
21) AllylChloride	9.305	39	146219m	10.88	ppbv	
22) CarbonDisulfide	9.486	76	477553m	9.63	ppbv	
23) Trichlorotrifluoroethane	8.998	103	246483	10.25	ppbv	97
24) trans-1,2-Dichloroethene	10.424	96	182437m	10.15	ppbv	
25) 1,1-Dichloroethane	10.906	63	361184	9.86	ppbv	100
26) MethylTertButylEther (M...)	10.442	73	514412	10.57	ppbv	99
27) VinylAcetate	10.888	43	438278m	9.68	ppbv	
28) 2-Butanone (MEK)	11.423	72	88194m	10.85	ppbv	
29) cis-1,2-Dichloroethene	11.905	96	197672	10.22	ppbv	99
30) Hexane	11.477	86	37597	9.70	ppbv	84
31) Chloroform	12.493	83	429780	10.56	ppbv	97
32) EthylAcetate	12.011	43	450464	11.00	ppbv	97

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Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311302.D
 Acq On : 31 May 2013 9:19
 Operator : JJG
 Sample : TO15 CCV 053113
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 31 15:58:54 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

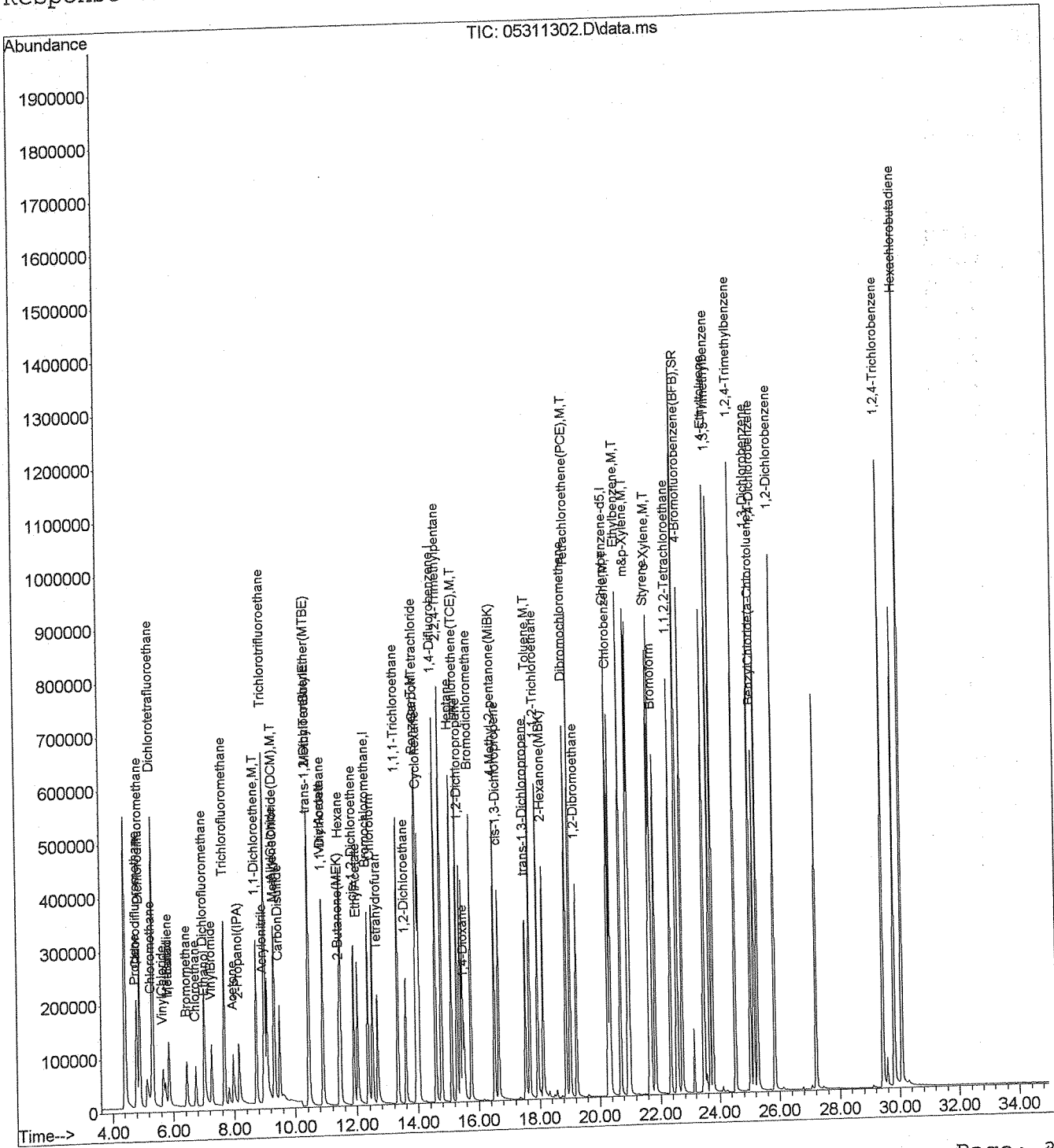
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) Tetrahydrofuran	12.671	72	852310	10.41	ppbv	
34) 1,2-Dichloroethane	13.599	62	315891	10.70	ppbv	97
35) 1,1,1-Trichloroethane	13.331	97	487892	10.83	ppbv	99
37) Benzene	13.937	78	552734	9.97	ppbv	99
38) CarbonTetrachloride	13.973	117	487943	10.65	ppbv	99
39) Cyclohexane	14.026	69	81482	9.92	ppbv	95
40) 1,2-Dichloropropane	15.400	63	228251	10.24	ppbv	96
41) Bromodichloromethane	15.756	85	305276	10.68	ppbv	99
42) 1,4-Dioxane	15.524	88	134023m	10.25	ppbv	
43) Trichloroethene (TCE)	15.293	130	278715	10.38	ppbv	99
44) 2,2,4-Trimethylpentane	14.775	57	1047326	10.62	ppbv	99
45) Heptane	15.114	71	185741	10.55	ppbv	99
46) cis-1,3-Dichloropropene	16.648	75	347585	10.95	ppbv	98
47) 4-Methyl-2-pentanone (M...)	16.523	58	209434	10.36	ppbv	98
48) trans-1,3-Dichloropropene	17.539	75	322018	9.96	ppbv	99
49) 1,1,2-Trichloroethane	17.932	97	268071	10.83	ppbv	99
50) Toluene	17.682	91	735162	10.46	ppbv	99
51) 2-Hexanone (MBK)	18.128	58	272743	10.89	ppbv	97
52) Dibromochloromethane	18.877	129	552160	11.62	ppbv	99
53) 1,2-Dibromoethane	19.233	107	422672	10.42	ppbv	99
54) Tetrachloroethene (PCE)	19.019	166	415184	10.54	ppbv	99
56) Chlorobenzene	20.357	114	197607	10.05	ppbv	99
57) Ethylbenzene	20.696	91	992383	9.94	ppbv	100
58) m&p-Xylene	20.999	106	731019	18.61	ppbv	98
59) Bromoform	21.819	173	527500	10.01	ppbv #	96
60) Styrene	21.641	104	619961	9.76	ppbv	100
61) 1,1,2,2-Tetrachloroethane	22.336	83	550882	9.68	ppbv	99
62) o-Xylene	21.694	91	760234	9.52	ppbv	99
64) 4-Ethyltoluene	23.674	120	324083	9.88	ppbv	99
65) 1,3,5-Trimethylbenzene	23.781	120	450315	9.40	ppbv	99
66) 1,2,4-Trimethylbenzene	24.529	120	457427	9.80	ppbv	99
67) BenzylChloride (a-Chlor...)	25.154	91	707251	10.07	ppbv	100
68) 1,3-Dichlorobenzene	25.047	146	703212	9.57	ppbv	99
69) 1,4-Dichlorobenzene	25.261	146	687499m	9.31	ppbv	99
70) 1,2-Dichlorobenzene	25.831	146	722735m	9.41	ppbv	99
71) 1,2,4-Trichlorobenzene	29.433	180	686176m	9.03	ppbv	97
72) Hexachlorobutadiene	30.075	225	578384m	9.40	ppbv	99

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

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Data Path : C:\msdchem\1\MS03\2013\053113\
Data File : 05311302.D
Acq On : 31 May 2013 9:19
Operator : JJJG
Sample : TO15 CCV 053113
Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 31 15:58:54 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\053113\
 Data File : 05311303.D
 Acq On : 31 May 2013 10:05
 Operator : JJG
 Sample : TO15 LCSD 053113
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 31 16:01:04 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	148225	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	769764	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	730036	10.00	ppbv	0.00
System Monitoring Compounds						
63) 4-Bromofluorobenzene (BFB)	22.710	174	462255	10.12	ppbv	0.00

Spiked Amount 10.000

Recovery = 101.20%

Target Compounds	R.T.	QIon	Response	Conc	Units	Dev (Min)	Qvalue
2) Chlorodifluoromethane	4.817	51	277483m	9.45	ppbv		
3) Propene	4.781	42	78550m	10.19	ppbv		
4) Dichlorodifluoromethane	4.908	85	473004	9.33	ppbv		99
5) Chloromethane	5.288	52	460730	9.69	ppbv		
6) Dichlorotetrafluoroethane	5.324	135	340828	9.87	ppbv		92
7) VinylChloride	5.650	62	165177m	9.80	ppbv		
8) Methanol	5.849	31	22878m	4.78	ppbv		
9) 1,3-Butadiene	5.849	54	105277m	9.84	ppbv		
10) Bromomethane	6.446	96	111831m	9.21	ppbv		0.00
11) Chloroethane	6.736	66	26950	9.83	ppbv		0.93
12) Dichlorofluoromethane	7.007	67	350375	9.66	ppbv		0.99
13) Ethanol	7.043	45	59358m	9.39	ppbv		
14) VinylBromide	7.260	108	148341m	10.40	ppbv		
15) Acetone	7.966	58	77767m	9.69	ppbv		0.00
16) Trichlorofluoromethane	7.677	103	307607	10.39	ppbv		99
17) 2-Propanol (IPA)	8.147	45	272676m	9.92	ppbv		
18) Acrylonitrile	8.961	52	126782m	10.55	ppbv		
19) 1,1-Dichloroethene	8.726	96	168791	10.00	ppbv		97
20) MethyleneChloride (DCM)	9.323	84	151145m	9.75	ppbv		
21) AllylChloride	9.305	39	148303m	10.92	ppbv		
22) CarbonDisulfide	9.486	76	486256m	9.70	ppbv		99
23) Trichlorotrifluoroethane	8.998	103	248277	10.22	ppbv		97
24) trans-1,2-Dichloroethene	10.424	96	185793m	10.23	ppbv		92
25) 1,1-Dichloroethane	10.906	63	360683	9.74	ppbv		100
26) MethylTertButylEther (M...)	10.442	73	515449m	10.48	ppbv		
27) VinylAcetate	10.888	43	436529m	9.54	ppbv		
28) 2-Butanone (MEK)	11.423	72	87320m	10.63	ppbv		
29) cis-1,2-Dichloroethene	11.904	96	200309	10.25	ppbv		99
30) Hexane	11.476	86	38543	9.84	ppbv		87
31) Chloroform	12.493	83	433724	10.54	ppbv		97
32) EthylAcetate	12.011	43	444177	10.73	ppbv		97

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311303.D
 Acq On : 31 May 2013 10:05
 Operator : JJG
 Sample : TO15 LCSD 053113
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 31 16:01:04 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

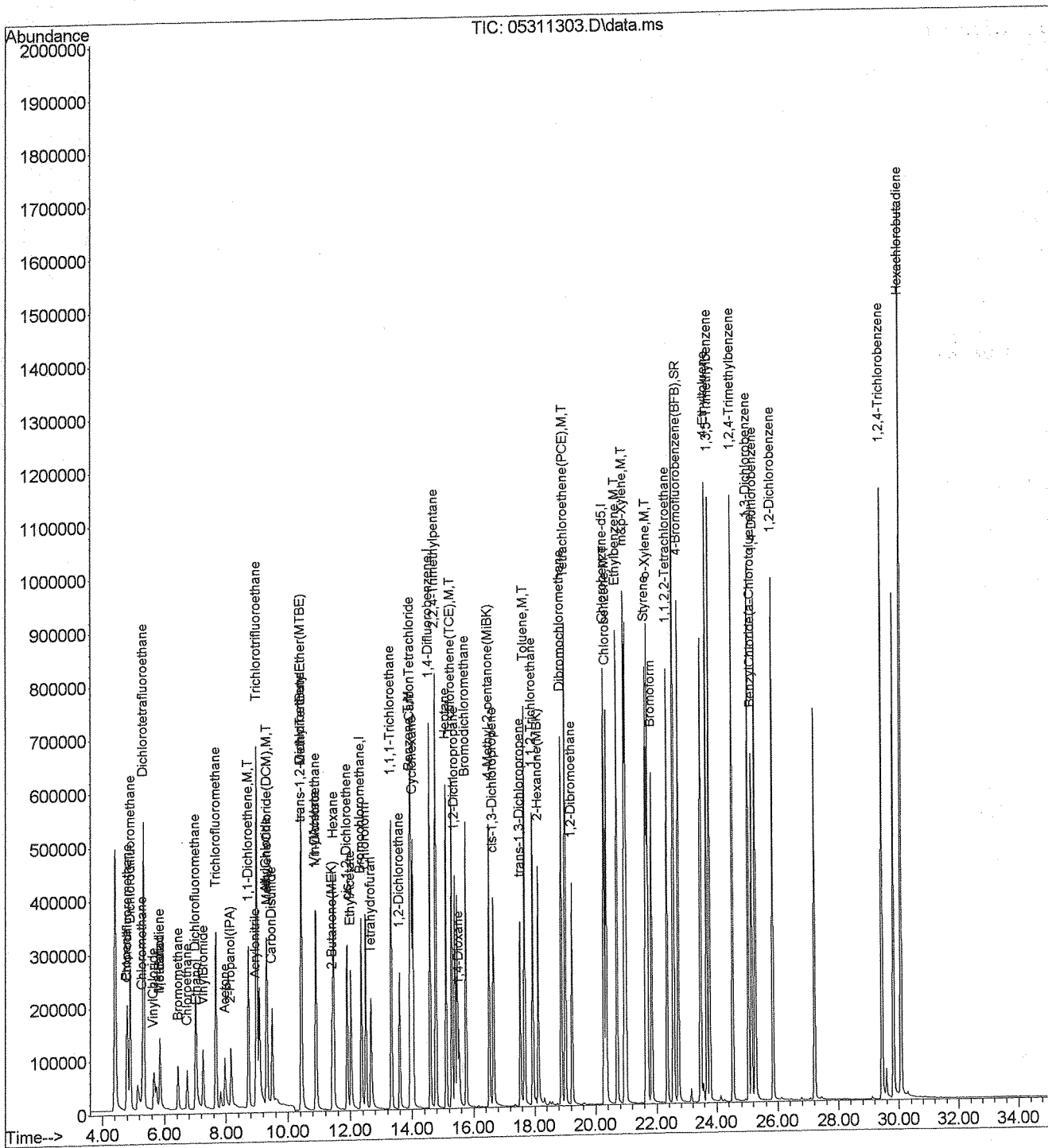
Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
33) Tetrahydrofuran	12.671	72	86404m	10.44	ppbv	
34) 1,2-Dichloroethane	13.598	62	316194	10.60	ppbv	99
35) 1,1,1-Trichloroethane	13.331	97	486237	10.68	ppbv	99
37) Benzene	13.937	78	555998	9.80	ppbv	99
38) CarbonTetrachloride	13.973	117	490779	10.47	ppbv	99
39) Cyclohexane	14.026	69	83350	9.92	ppbv	94
40) 1,2-Dichloropropane	15.399	63	225436	9.89	ppbv	97
41) Bromodichloromethane	15.756	85	306686	10.48	ppbv	98
42) 1,4-Dioxane	15.524	88	134724m	10.07	ppbv	
43) Trichloroethene (TCE)	15.292	130	278045	10.12	ppbv	98
44) 2,2,4-Trimethylpentane	14.775	57	1051308	10.41	ppbv	99
45) Heptane	15.114	71	179210	9.95	ppbv	97
46) cis-1,3-Dichloropropene	16.647	75	355245	10.94	ppbv	98
47) 4-Methyl-2-pentanone (M...)	16.523	58	209891	10.15	ppbv	98
48) trans-1,3-Dichloropropene	17.539	75	321126	9.71	ppbv	99
49) 1,1,2-Trichloroethane	17.931	97	266677	10.53	ppbv	99
50) Toluene	17.682	91	743636	10.34	ppbv	100
51) 2-Hexanone (MBK)	18.127	58	268499	10.48	ppbv	97
52) Dibromochloromethane	18.876	129	558658	11.49	ppbv	99
53) 1,2-Dibromoethane	19.233	107	428542	10.33	ppbv	99
54) Tetrachloroethene (PCE)	19.019	166	413007	10.24	ppbv	99
56) Chlorobenzene	20.356	114	196745	10.37	ppbv	98
57) Ethylbenzene	20.695	91	974128	10.11	ppbv	100
58) m&p-Xylene	20.945	106	756674	19.96	ppbv	95
59) Bromoform	21.836	173	518783	10.20	ppbv	100
60) Styrene	21.640	104	623918	10.17	ppbv	100
61) 1,1,2,2-Tetrachloroethane	22.336	83	563905	10.26	ppbv	98
62) o-Xylene	21.694	91	760674	9.87	ppbv	100
64) 4-Ethyltoluene	23.673	120	329642	10.42	ppbv	100
65) 1,3,5-Trimethylbenzene	23.780	120	456676	9.88	ppbv	99
66) 1,2,4-Trimethylbenzene	24.529	120	454079	10.08	ppbv	98
67) BenzylChloride (a-Chlor...)	25.153	91	721619	10.65	ppbv	99
68) 1,3-Dichlorobenzene	25.046	146	714472	10.08	ppbv	99
69) 1,4-Dichlorobenzene	25.260	146	708397m	9.94	ppbv	99
70) 1,2-Dichlorobenzene	25.831	146	723618m	9.76	ppbv	100
71) 1,2,4-Trichlorobenzene	29.433	180	694134m	9.47	ppbv	97
72) Hexachlorobutadiene	30.075	225	576795m	9.72	ppbv	99

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

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311303.D
 Acq On : 31 May 2013 10:05
 Operator : JJG
 Sample : TO15 LCSO 053113
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 31 16:01:04 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\053113\
 Data File : 05311304.D
 Acq On : 31 May 2013 10:53
 Operator : JJG
 Sample : TO15 MB 053113
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 31 16:01:41 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	138528	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	768901	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	714667	10.00	ppbv	0.00
System Monitoring Compounds						
63) 4-Bromofluorobenzene (BFB)	22.711	174	457690	10.23	ppbv	0.00
Spiked Amount	10.000			Recovery	=	102.30%

Target Compounds					Qvalue
2) Chlorodifluoromethane	0.000		0		N.D.
3) Propene	4.836	42	129		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. d
8) Methanol	0.000		0		N.D.
9) 1,3-Butadiene	0.000		0		N.D. d
10) Bromomethane	0.000		0		N.D.
11) Chloroethane	0.000		0		N.D.
12) Dichlorofluoromethane	0.000		0		N.D.
13) Ethanol	0.000		0		N.D.
14) VinylBromide	0.000		0		N.D. d
15) Acetone	0.000		0		N.D.
16) Trichlorofluoromethane	0.000		0		N.D.
17) 2-Propanol (IPA)	8.328	45	241		N.D.
18) Acrylonitrile	0.000		0		N.D.
19) 1,1-Dichloroethene	0.000		0		N.D. d
20) MethyleneChloride (DCM)	0.000		0		N.D.
21) AllylChloride	0.000		0		N.D. d
22) CarbonDisulfide	0.000		0		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.
26) MethylTertButylEther (M...)	0.000		0		N.D.
27) VinylAcetate	0.000		0		N.D.
28) 2-Butanone (MEK)	0.000		0		N.D.
29) cis-1,2-Dichloroethene	0.000		0		N.D.
30) Hexane	0.000		0		N.D.
31) Chloroform	0.000		0		N.D.
32) EthylAcetate	0.000		0		N.D.

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311304.D
 Acq On : 31 May 2013 10:53
 Operator : JJG
 Sample : TO15 MB 053113
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 1 Sample Multiplier: 1

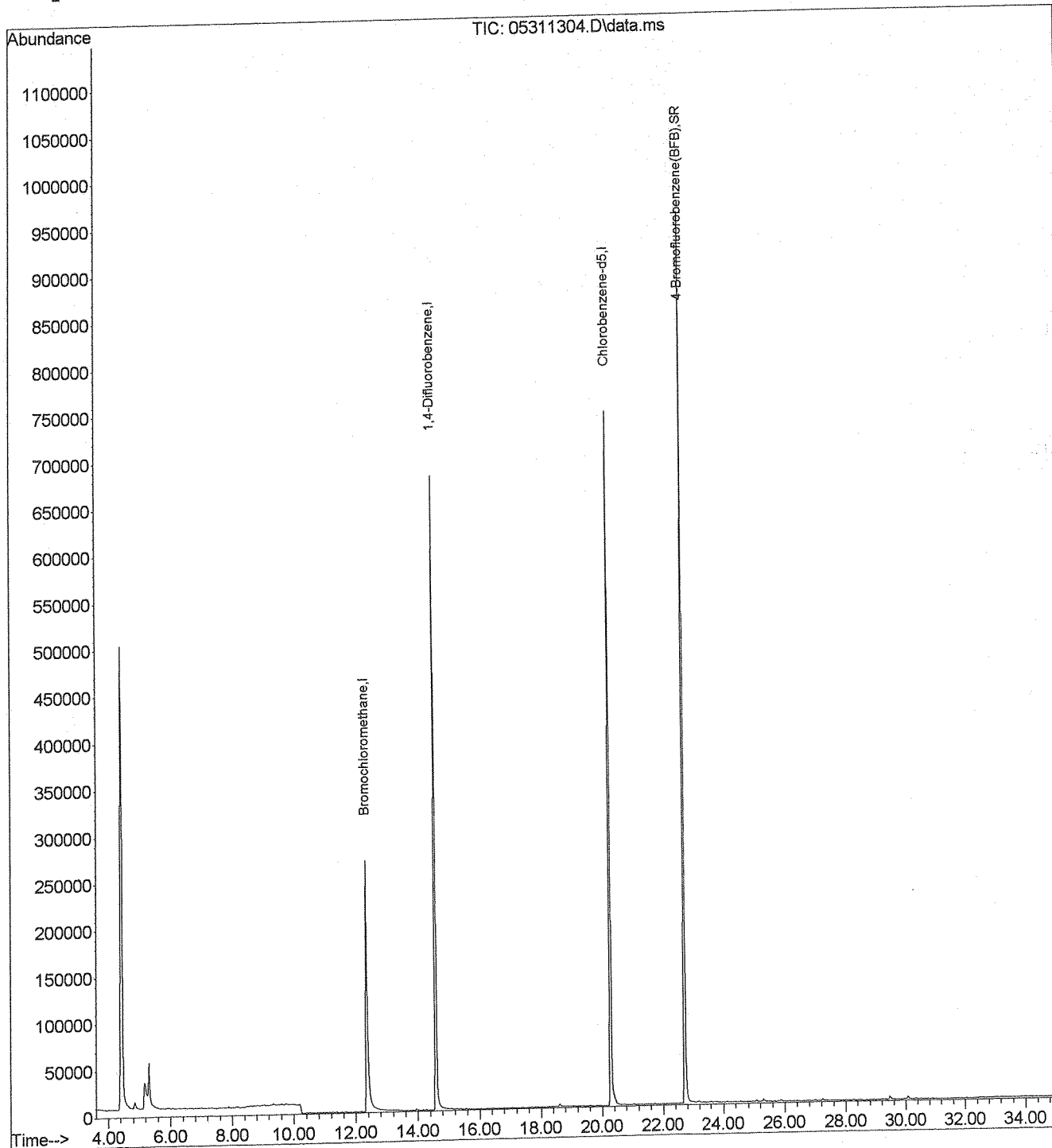
Quant Time: May 31 16:01:41 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

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	13.937	78	2127		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	769		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	123		N.D.	
57) Ethylbenzene	20.731	91	740		N.D.	
58) m&p-Xylene	21.017	106	120		N.D.	
59) Bromoform	0.000		0		N.D.	
60) Styrene	21.694	104	296		N.D.	
61) 1,1,2,2-Tetrachloroethane	22.354	83	175		N.D.	
62) o-Xylene	21.712	91	581		N.D.	
64) 4-Ethyltoluene	23.691	120	254		N.D.	
65) 1,3,5-Trimethylbenzene	23.798	120	381		N.D.	
66) 1,2,4-Trimethylbenzene	24.565	120	306		N.D.	
67) BenzylChloride (a-Chlor...)	25.207	91	770		N.D.	
68) 1,3-Dichlorobenzene	25.064	146	2005		N.D.	
69) 1,4-Dichlorobenzene	25.296	146	2609		N.D.	
70) 1,2-Dichlorobenzene	25.867	146	1342		N.D.	
71) 1,2,4-Trichlorobenzene	0.000		0		N.D.	d
72) Hexachlorobutadiene	30.075	225	1188		N.D.	

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

Data Path : C:\msdchem\1\MS03\2013\053113\
Data File : 05311304.D
Acq On : 31 May 2013 10:53
Operator : JJG
Sample : TO15 MB 053113
Misc : IS/Surr: PS082712-02 + 500mL
ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 31 16:01:41 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\053113\
 Data File : 05311305.D
 Acq On : 31 May 2013 11:41
 Operator : JJG
 Sample : 130647-63190 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 31 16:03:57 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	141154	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	774304	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	740604	10.00	ppbv	0.00
System Monitoring Compounds						
63) 4-Bromofluorobenzene (BFB)	22.710	174	470787	10.16	ppbv	0.00
Spiked Amount	10.000		Recovery	=	101.60%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue	Dev(Min)
2) Chlorodifluoromethane	4.835	51	4423	0.16	ppbv	#	96
3) Propene	0.000		0	N.D.	d		
4) Dichlorodifluoromethane	4.908	85	13542	0.28	ppbv		98
5) Chloromethane	5.306	52	1042	0.23	ppbv	#	6
6) Dichlorotetrafluoroethane	5.324	135	228	N.D.			
7) VinylChloride	0.000		0	N.D.			
8) Methanol	5.867	31	33313m	7.37	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.			0.00
13) Ethanol	7.134	45	17815m	2.96	ppbv		
14) VinylBromide	0.000		0	N.D.			
15) Acetone	8.002	58	19956m	2.61	ppbv		0.00
16) Trichlorofluoromethane	7.658	103	3831	0.14	ppbv		97
17) 2-Propanol (IPA)	8.219	45	22036m	0.84	ppbv		
18) Acrylonitrile	0.000		0	N.D.			
19) 1,1-Dichloroethene	0.000		0	N.D.			
20) MethyleneChloride (DCM)	0.000		0	N.D.	d	#	98
21) AllylChloride	9.233	39	128	N.D.			
22) CarbonDisulfide	0.000		0	N.D.	d		98
23) Trichlorotrifluoroethane	0.000		0	N.D.	d	#	6
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	10.888	43	1152	N.D.			
28) 2-Butanone (MEK)	0.000		0	N.D.	d		
29) cis-1,2-Dichloroethene	0.000		0	N.D.			
30) Hexane	0.000		0	N.D.	d		
31) Chloroform	12.493	83	396	N.D.			
32) EthylAcetate	0.000		0	N.D.	d		

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311305.D
 Acq On : 31 May 2013 11:41
 Operator : JJG
 Sample : 130647-63190 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 2 Sample Multiplier: 1

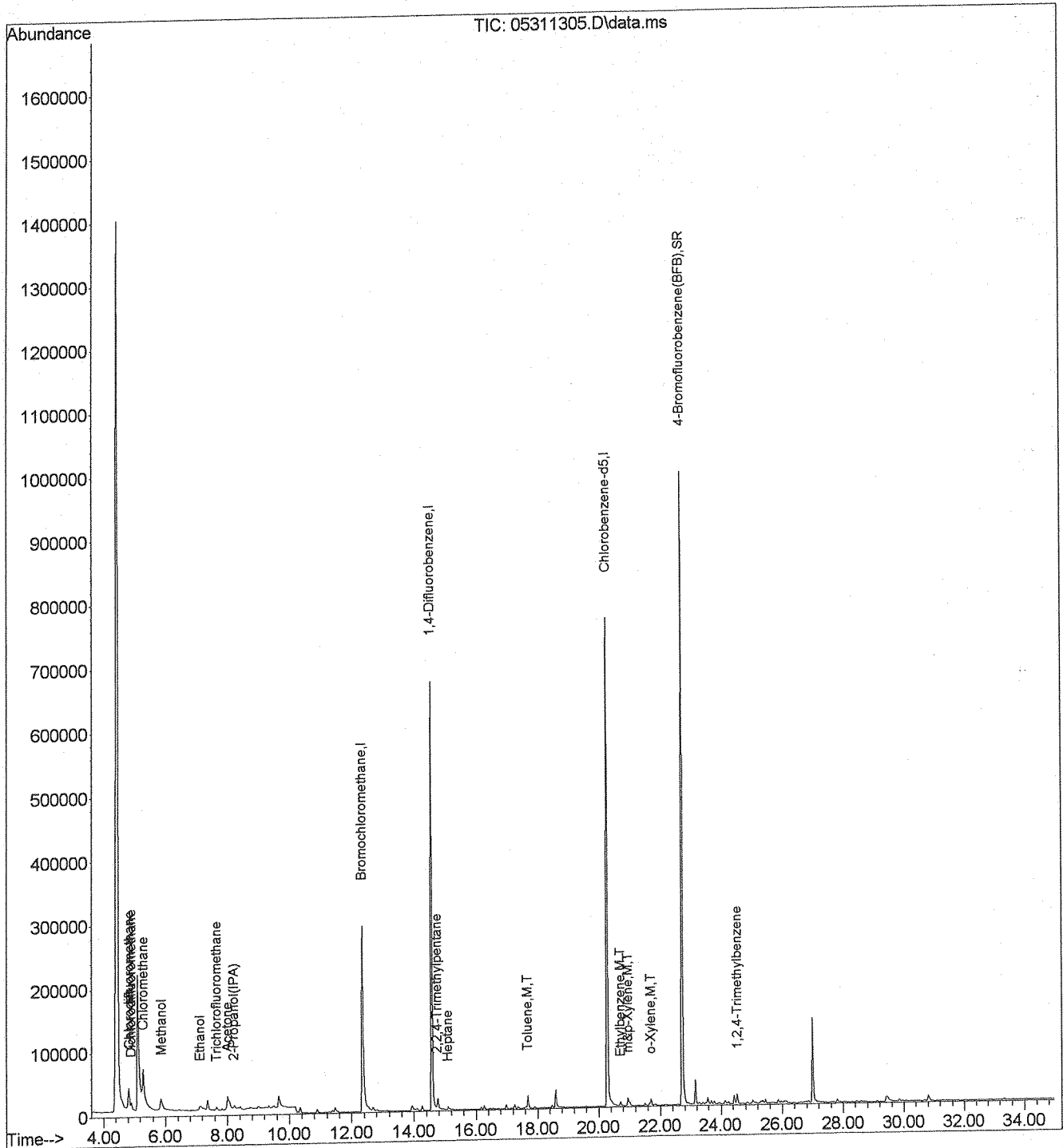
Quant Time: May 31 16:03:57 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

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	0.000		0		N.D. d	
39) Cyclohexane	14.008	69	282		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	14.757	57	22408	0.22	ppbv #	98
45) Heptane	15.096	71	1229	0.07	ppbv #	83
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	17.682	75	120		N.D.	
49) 1,1,2-Trichloroethane	0.000		0		N.D.	
50) Toluene	17.682	91	22183	0.31	ppbv #	97
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	147		N.D.	
57) Ethylbenzene	20.713	91	7119	0.07	ppbv #	96
58) m&p-Xylene	20.963	106	9179	0.24	ppbv #	97
59) Bromoform	0.000		0		N.D.	
60) Styrene	21.676	104	595		N.D.	
61) 1,1,2,2-Tetrachloroethane	0.000		0		N.D.	
62) o-Xylene	21.694	91	8412	0.11	ppbv	95
64) 4-Ethyltoluene	0.000		0		N.D. d	
65) 1,3,5-Trimethylbenzene	0.000		0		N.D. d	
66) 1,2,4-Trimethylbenzene	24.547	120	7269	0.16	ppbv #	96
67) BenzylChloride (a-Chlor...)	25.207	91	379		N.D.	
68) 1,3-Dichlorobenzene	25.064	146	586		N.D.	
69) 1,4-Dichlorobenzene	25.296	146	1078		N.D.	
70) 1,2-Dichlorobenzene	25.849	146	614		N.D.	
71) 1,2,4-Trichlorobenzene	29.451	180	1944		N.D.	
72) Hexachlorobutadiene	30.075	225	613		N.D.	

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

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311305.D
 Acq On : 31 May 2013 11:41
 Operator : JJG
 Sample : 130647-63190 x1
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 31 16:03:57 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\053113\
 Data File : 05311306.D
 Acq On : 31 May 2013 12:29
 Operator : JJG
 Sample : 130647-63190 x1 dp
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 31 16:06:29 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	139450	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	768342	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	721337	10.00	ppbv	0.00

System Monitoring Compounds						
63) 4-Bromofluorobenzene (BFB)	22.711	174	471546	10.45	ppbv	0.00

Spiked Amount 10.000

Recovery = 104.50%

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue	Dev(Min)
2) Chlorodifluoromethane	4.836	51	4269	0.15	ppbv	# 96	
3) Propene	0.000		0	N.D.	d		
4) Dichlorodifluoromethane	4.908	85	13742	0.29	ppbv	# 98	
5) Chloromethane	5.306	52	1085	0.24	ppbv	# 13	
6) Dichlorotetrafluoroethane	5.324	135	113	N.D.			
7) VinylChloride	0.000		0	N.D.			
8) Methanol	5.867	31	333940	7.48	ppbv		
9) 1,3-Butadiene	0.000		0	N.D.	d		
10) Bromomethane	0.000		0	N.D.			
11) Chloroethane	0.000		0	N.D.			
12) Dichlorofluoromethane	0.000		0	N.D.			
13) Ethanol	7.134	45	169640	2.85	ppbv		
14) VinylBromide	0.000		0	N.D.			
15) Acetone	8.002	58	193530	2.56	ppbv		
16) Trichlorofluoromethane	7.659	103	3946	0.14	ppbv	95	
17) 2-Propanol (IPA)	8.220	45	222330	0.86	ppbv		
18) Acrylonitrile	0.000		0	N.D.			
19) 1,1-Dichloroethene	0.000		0	N.D.	d	# 96	
20) MethyleneChloride (DCM)	0.000		0	N.D.			
21) AllylChloride	9.251	39	237	N.D.			
22) CarbonDisulfide	0.000		0	N.D.	d	# 98	
23) Trichlorotrifluoroethane	8.998	103	910	N.D.		# 13	
24) trans-1,2-Dichloroethene	0.000		0	N.D.			
25) 1,1-Dichloroethane	0.000		0	N.D.			
26) MethylTertButyleEther (M...)	0.000		0	N.D.			
27) VinylAcetate	10.888	43	1305	N.D.			
28) 2-Butanone (MEK)	0.000		0	N.D.	d		
29) cis-1,2-Dichloroethene	0.000		0	N.D.			
30) Hexane	0.000		0	N.D.	d		
31) Chloroform	12.493	83	322	N.D.			
32) EthylAcetate	0.000		0	N.D.	d		

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Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311306.D
 Acq On : 31 May 2013 12:29
 Operator : JJG
 Sample : 130647-63190 x1 dp
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 2 Sample Multiplier: 1

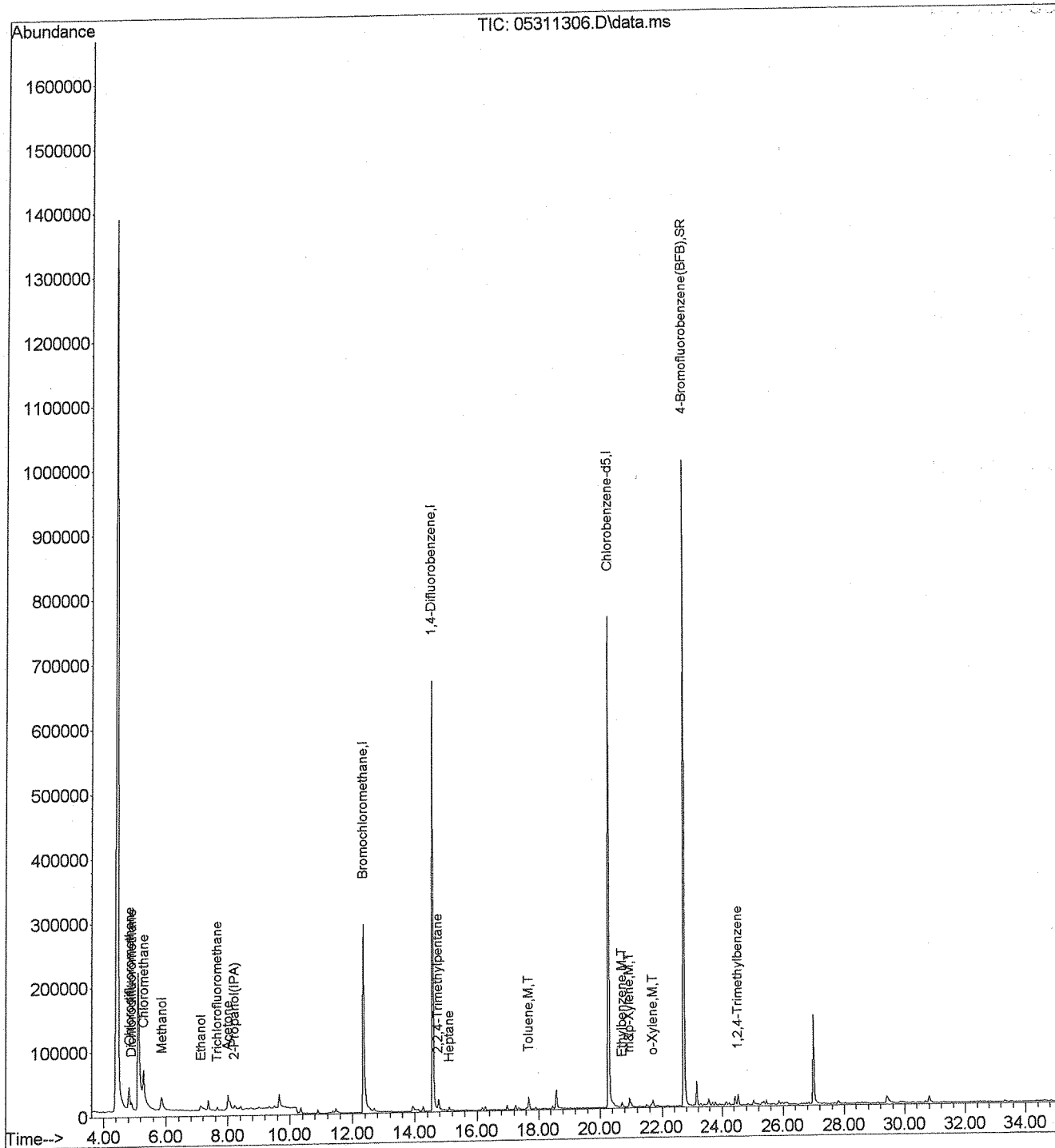
Quant Time: May 31 16:06:29 2013
 Quant Method : C:\msdchem\1\METHODS\2013\051513.M
 Quant Title : TO-15/TO-14
 QLast Update : Thu May 16 10:13:39 2013
 Response via : Initial Calibration

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	0.000		0		N.D. d	
39) Cyclohexane	14.026	69	297		N.D.	
40) 1,2-Dichloropropane	15.275	63	117		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	14.758	57	21428	0.21	ppbv	95
45) Heptane	15.114	71	1228	0.07	ppbv #	89
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	17.682	75	132		N.D.	
49) 1,1,2-Trichloroethane	0.000		0		N.D.	
50) Toluene	17.682	91	20636	0.29	ppbv	(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	20.268	114	256		N.D.	
57) Ethylbenzene	20.713	91	7287	0.08	ppbv #	96
58) m&p-Xylene	20.945	106	8927	0.24	ppbv #	94
59) Bromoform	0.000		0		N.D.	
60) Styrene	21.676	104	657		N.D.	
61) 1,1,2,2-Tetrachloroethane	0.000		0		N.D.	
62) o-Xylene	21.694	91	8082	0.11	ppbv #	92
64) 4-Ethyltoluene	0.000		0		N.D. d	
65) 1,3,5-Trimethylbenzene	0.000		0		N.D. d	
66) 1,2,4-Trimethylbenzene	24.529	120	7104	0.16	ppbv	97
67) BenzylChloride (a-Chlor...)	25.189	91	112		N.D.	
68) 1,3-Dichlorobenzene	25.064	146	149		N.D.	
69) 1,4-Dichlorobenzene	25.296	146	505		N.D.	
70) 1,2-Dichlorobenzene	25.867	146	303		N.D.	
71) 1,2,4-Trichlorobenzene	29.451	180	1067		N.D.	
72) Hexachlorobutadiene	30.075	225	267		N.D.	

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

Data Path : C:\msdchem\1\MS03\2013\053113\
 Data File : 05311306.D
 Acq On : 31 May 2013 12:29
 Operator : JJG
 Sample : 130647-63190 x1 dp
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 31 16:06:29 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



MS #3 Instrument Logbook

Sequence Name: C:\msdchem\1\sequence\2013\060313.S

Comment: GCMS-03

Operator: JJG

Data Path: C:\MSDCHEM\1\MS03\2013\060313\

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 06031301 TO15-5MS TO15 BFB 060313
2) Sample	1 06031302 TO15-5MS TO15 CCV 060313
3) Sample	1 06031303 TO15-5MS TO15 LCSD 060313
4) Sample	1 06031304 TO15-5MS TO15 MB 060313
5) Sample	2 06031305 TO15-5MS 130650-63200 x10
6) Sample	2 06031306 TO15-5MS 130650-63200 x10 dp
7) Sample	3 06031307 TO15-5MS 130650-63209 x10
8) Sample	4 06031308 TO15-5MS 130653-63267 x5
9) Sample	5 06031309 TO15-5MS 130638-63167 x10
10) Sample	6 06031310 TO15-5MS 130637-63164 x10
11) Sample	7 06031311 TO15-5MS 130638-63165 x10
12) Sample	8 06031312 TO15-5MS 130637-63162 x10
13) Sample	9 06031313 TO15-5MS 130656-63280 x20
14) Sample	9 06031314 TO15-5MS 130656-63280 x50
15) Sample	9 06031315 TO15-5MS 130656-63280 x100
16) Sample	8 06031316 TO15-5MS 130637-63162 x25
17) Sample	7 06031317 TO15-5MS 130638-63165 x25
18) Sample	6 06031318 TO15-5MS 130637-63164 x25
19) Sample	8 06031319 TO15-5MS 130637-63162 x50
20) Sample	9 06031320 TO15-5MS 130656-63280 x50 rr
21) Sample	10 06031321 TO15-5MS 130656-63280 x100
22) Sample	10 06031322 TO15-5MS 130656-63280 x100 dp
23) Sample	11 06031323 TO15-5MS 130637-63162 x100
24) Sample	12 06031324 TO15-5MS 130637-63164 x100
25) Sample	13 06031325 TO15-5MS 130638-63165 x100
26) Sample	14 06031326 TO15-5MS 130638-63167 x100
27) Sample	15 06031327 TO15-5MS Lab Air 060313 x1
28) Sample	16 06031328 TO15-5MS Lab Air 060313 x1
29) Sample	16 06031329 TO15-5MS Lab Air 060313 x1
30) Sample	16 06031330 TO15-5MS Lab Air 060313 x1
31) Sample	16 06031331 TO15-5MS Lab Air 060313 x1

Comments: _____

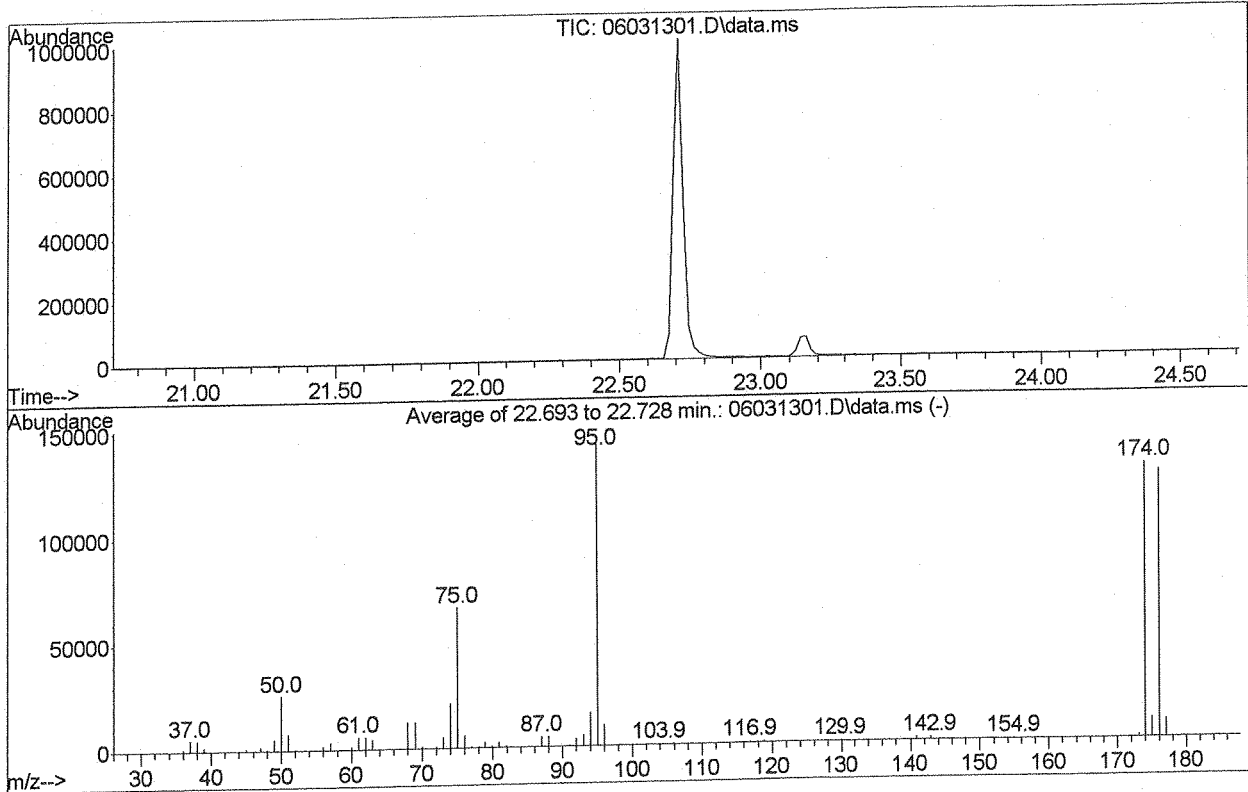
Analyst: _____

Date: 06/03/13

Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031301.D
 Acq On : 3 Jun 2013 8:31 am
 Operator : JJG
 Sample : TO15 BFB 060313
 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	17.8	25633	PASS
75	95	30	60	46.0	66285	PASS
95	95	100	100	100.0	144131	PASS
96	95	5	9	6.7	9589	PASS
173	174	0.00	2	0.9	1151	PASS
174	95	50	100	90.0	129768	PASS
175	174	5	9	7.4	9580	PASS
176	174	95	101	97.5	126579	PASS
177	176	5	9	6.7	8458	PASS

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Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031302.D
 Acq On : 3 Jun 2013 9:17
 Operator : JJG
 Sample : TO15 CCV 060313
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Jun 03 12:30:29 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	150793	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	783130	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	731958	10.00	ppbv	0.00
System Monitoring Compounds						
63) 4-Bromofluorobenzene (BFB)	22.711	174	460513	10.05	ppbv	0.00
Spiked Amount	10.000			Recovery	=	100.50%

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Chlorodifluoromethane	4.817	51	282780	9.47	ppbv	100
3) Propene	4.781	42	77829m	9.92	ppbv	
4) Dichlorodifluoromethane	4.908	85	493281	9.57	ppbv	99
5) Chloromethane	5.288	52	48654m	10.06	ppbv	
6) Dichlorotetrafluoroethane	5.342	135	356314	10.14	ppbv	98
7) VinylChloride	5.668	62	172667m	10.07	ppbv	
8) Methanol	5.849	31	25351m	5.21	ppbv	
9) 1,3-Butadiene	5.867	54	103467m	9.50	ppbv	
10) Bromomethane	6.446	96	102263m	8.28	ppbv	99
11) Chloroethane	6.736	66	26873	9.63	ppbv	98
12) Dichlorofluoromethane	7.025	67	371073	10.06	ppbv	99
13) Ethanol	7.043	45	67120m	10.44	ppbv	
14) VinylBromide	7.260	108	149796m	10.33	ppbv	
15) Acetone	7.966	58	77626m	9.50	ppbv	99
16) Trichlorofluoromethane	7.677	103	321560	10.68	ppbv	97
17) 2-Propanol (IPA)	8.165	45	287042m	10.27	ppbv	
18) Acrylonitrile	8.961	52	124545m	10.19	ppbv	
19) 1,1-Dichloroethene	8.726	96	174866	10.18	ppbv	94
20) MethyleneChloride (DCM)	9.323	84	156913m	9.95	ppbv	100
21) AllylChloride	9.305	39	153252m	11.09	ppbv	
22) CarbonDisulfide	9.486	76	497528m	9.76	ppbv	99
23) Trichlorotrifluoroethane	8.998	103	250889	10.15	ppbv	95
24) trans-1,2-Dichloroethene	10.424	96	185329	10.03	ppbv	99
25) 1,1-Dichloroethane	10.906	63	382225	10.15	ppbv	100
26) MethylTertButylEther (M...)	10.442	73	539250	10.78	ppbv	98
27) VinylAcetate	10.888	43	466423m	10.02	ppbv	
28) 2-Butanone (MEK)	11.423	72	90852m	10.87	ppbv	
29) cis-1,2-Dichloroethene	11.904	96	206277	10.37	ppbv	99
30) Hexane	11.476	86	39621	9.94	ppbv	85
31) Chloroform	12.493	83	445666	10.65	ppbv	97
32) EthylAcetate	12.011	43	466969	11.09	ppbv	97

Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031302.D
 Acq On : 3 Jun 2013 9:17
 Operator : JJG
 Sample : TO15 CCV 060313
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Jun 03 12:30:29 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

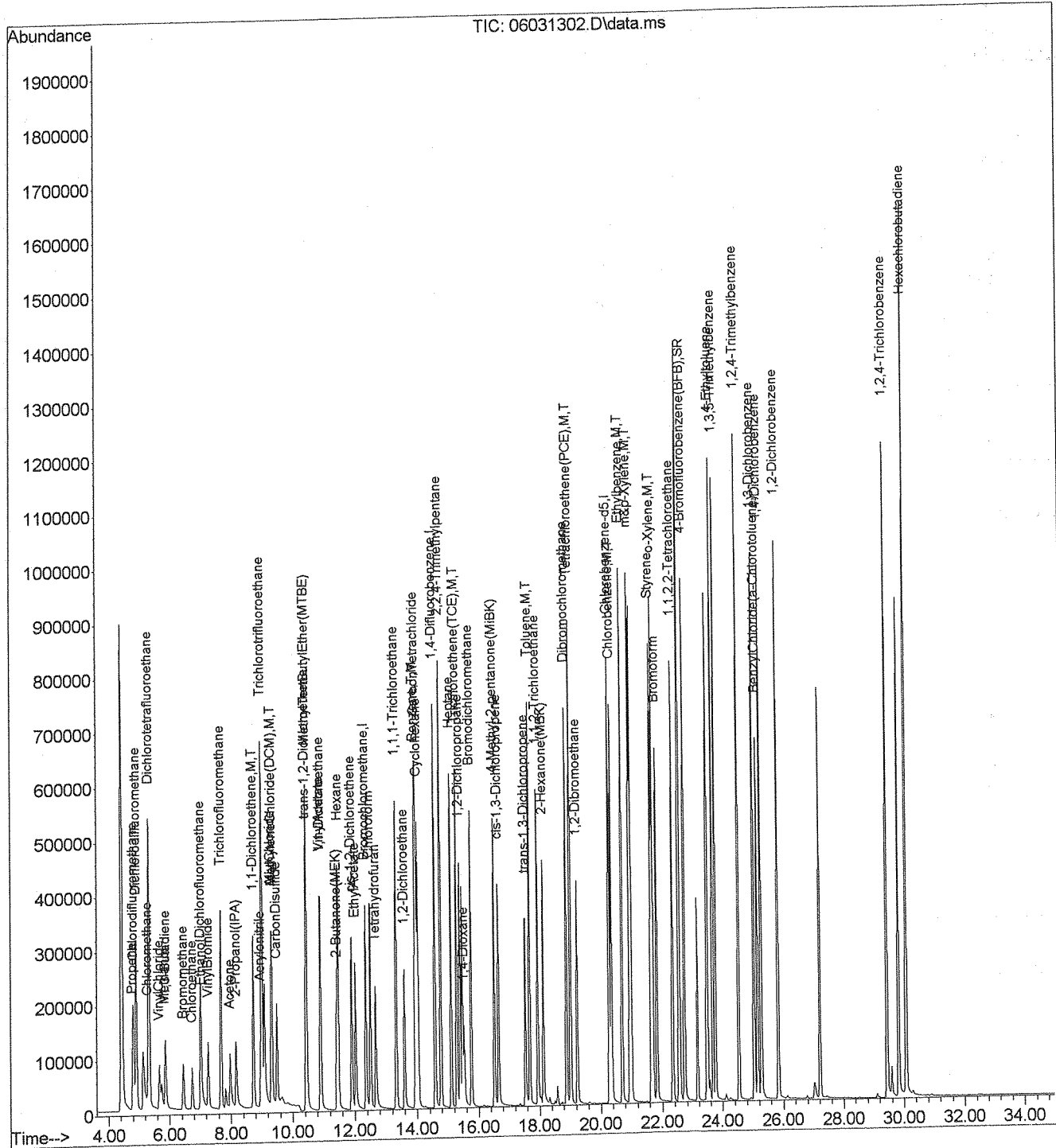
Compound	R.T.	QI on	Response	Conc	Units	Dev (Min)
33) Tetrahydrofuran	12.671	72	87643	10.41	ppbv	90
34) 1,2-Dichloroethane	13.598	62	328345	10.81	ppbv	98
35) 1,1,1-Trichloroethane	13.331	97	504547	10.89	ppbv	99
37) Benzene	13.937	78	575311	9.97	ppbv	99
38) CarbonTetrachloride	13.973	117	499891	10.48	ppbv	99
39) Cyclohexane	14.026	69	86234	10.08	ppbv	95
40) 1,2-Dichloropropane	15.399	63	232322	10.01	ppbv	96
41) Bromodichloromethane	15.756	85	305477	10.26	ppbv	99
42) 1,4-Dioxane	15.524	88	134834m	9.91	ppbv	
43) Trichloroethene (TCE)	15.292	130	284459	10.18	ppbv	99
44) 2,2,4-Trimethylpentane	14.775	57	1090124	10.61	ppbv	99
45) Heptane	15.114	71	186909	10.20	ppbv	98
46) cis-1,3-Dichloropropene	16.648	75	361949	10.96	ppbv	99
47) 4-Methyl-2-pentanone (M...)	16.523	58	216602	10.30	ppbv	96
48) trans-1,3-Dichloropropene	17.539	75	330786	9.83	ppbv	99
49) 1,1,2-Trichloroethane	17.932	97	269506	10.46	ppbv	99
50) Toluene	17.682	91	759202	10.37	ppbv	100
51) 2-Hexanone (MBK)	18.128	58	269319	10.33	ppbv	97
52) Dibromochloromethane	18.877	129	565442	11.43	ppbv	99
53) 1,2-Dibromoethane	19.233	107	435865	10.33	ppbv	100
54) Tetrachloroethene (PCE)	19.019	166	420229	10.25	ppbv	99
56) Chlorobenzene	20.357	114	200167	10.52	ppbv	99
57) Ethylbenzene	20.695	91	1028425	10.65	ppbv	100
58) m&p-Xylene	20.945	106	773279	20.34	ppbv	98
59) Bromoform	21.819	173	533326	10.46	ppbv #	95
60) Styrene	21.641	104	621590	10.11	ppbv	99
61) 1,1,2,2-Tetrachloroethane	22.336	83	568157	10.31	ppbv	99
62) o-Xylene	21.694	91	781733	10.12	ppbv	100
64) 4-Ethyltoluene	23.673	120	337038	10.62	ppbv	100
65) 1,3,5-Trimethylbenzene	23.780	120	459329	9.91	ppbv	98
66) 1,2,4-Trimethylbenzene	24.529	120	467805	10.36	ppbv	99
67) BenzylChloride (a-Chlor...)	25.153	91	746115	10.98	ppbv	99
68) 1,3-Dichlorobenzene	25.046	146	728004	10.24	ppbv	99
69) 1,4-Dichlorobenzene	25.260	146	696508m	9.74	ppbv	99
70) 1,2-Dichlorobenzene	25.831	146	732920m	9.86	ppbv	100
71) 1,2,4-Trichlorobenzene	29.433	180	695325m	9.46	ppbv	97
72) Hexachlorobutadiene	30.075	225	575537m	9.67	ppbv	99

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

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031302.D
 Acq On : 3 Jun 2013 9:17
 Operator : JJG
 Sample : TO15 CCV 060313
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Jun 03 12:30:29 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



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Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031303.D
 Acq On : 3 Jun 2013 10:03
 Operator : JJG
 Sample : TO15 LCSD 060313
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Jun 03 12:32:31 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	147894	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	793680	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	750236	10.00	ppbv	0.00
System Monitoring Compounds						
63) 4-Bromofluorobenzene (BFB)	22.711	174	459525	9.79	ppbv	0.00
Spiked Amount	10.000		Recovery	=	97.90%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Chlorodifluoromethane	4.817	51	272820	9.31	ppbv	99
3) Propene	4.781	42	76964m	10.00	ppbv	
4) Dichlorodifluoromethane	4.908	85	481319	9.52	ppbv	100
5) Chloromethane	5.288	52	46927m	9.89	ppbv	
6) Dichlorotetrafluoroethane	5.342	135	346447	10.05	ppbv	
7) VinylChloride	5.668	62	169751m	10.09	ppbv	
8) Methanol	5.849	31	23820m	4.99	ppbv	
9) 1,3-Butadiene	5.867	54	103101m	9.65	ppbv	
10) Bromomethane	6.446	96	108005m	8.92	ppbv	100.00
11) Chloroethane	6.736	66	25388	9.28	ppbv	99.99
12) Dichlorofluoromethane	7.025	67	371407m	10.26	ppbv	100.00
13) Ethanol	7.043	45	61505m	9.75	ppbv	
14) VinylBromide	7.260	108	148456m	10.43	ppbv	
15) Acetone	7.966	58	74522m	9.30	ppbv	100.00
16) Trichlorofluoromethane	7.677	103	312619	10.58	ppbv	99
17) 2-Propanol (IPA)	8.165	45	281595m	10.27	ppbv	99.99
18) Acrylonitrile	8.962	52	128892m	10.75	ppbv	
19) 1,1-Dichloroethene	8.726	96	169853	10.08	ppbv	96
20) MethyleneChloride (DCM)	9.323	84	155569m	10.06	ppbv	99
21) AllylChloride	9.305	39	147031m	10.85	ppbv	
22) CarbonDisulfide	9.486	76	485132m	9.70	ppbv	100
23) Trichlorotrifluoroethane	8.998	103	249249	10.28	ppbv	96
24) trans-1,2-Dichloroethene	10.424	96	187597m	10.36	ppbv	99
25) 1,1-Dichloroethane	10.906	63	368954	9.99	ppbv	99
26) MethylTertButylEther (M...)	10.442	73	523521	10.67	ppbv	98
27) VinylAcetate	10.888	43	455161m	9.97	ppbv	
28) 2-Butanone (MEK)	11.423	72	87961	10.73	ppbv	93
29) cis-1,2-Dichloroethene	11.904	96	204956	10.51	ppbv	99
30) Hexane	11.476	86	40772	10.43	ppbv	89
31) Chloroform	12.511	83	442508	10.78	ppbv	99
32) EthylAcetate	12.011	43	466192	11.29	ppbv	97

Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031303.D
 Acq On : 3 Jun 2013 10:03
 Operator : JJG
 Sample : TO15 LCSD 060313
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Jun 03 12:32:31 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	87286	10.57	ppbv	95
34) 1,2-Dichloroethane	13.598	62	319521	10.73	ppbv	99
35) 1,1,1-Trichloroethane	13.331	97	497848	10.96	ppbv	99
37) Benzene	13.937	78	572047	9.78	ppbv	99
38) CarbonTetrachloride	13.973	117	497821	10.30	ppbv	99
39) Cyclohexane	14.026	69	84658	9.77	ppbv	96
40) 1,2-Dichloropropane	15.400	63	234351	9.97	ppbv	97
41) Bromodichloromethane	15.756	85	312316	10.35	ppbv	99
42) 1,4-Dioxane	15.524	88	137235m	9.95	ppbv	
43) Trichloroethene (TCE)	15.293	130	284633	10.05	ppbv	99
44) 2,2,4-Trimethylpentane	14.775	57	1067684	10.26	ppbv	99
45) Heptane	15.114	71	188762	10.16	ppbv	98
46) cis-1,3-Dichloropropene	16.648	75	360975	10.78	ppbv	99
47) 4-Methyl-2-pentanone (M...)	16.523	58	214533	10.06	ppbv	97
48) trans-1,3-Dichloropropene	17.539	75	329615	9.67	ppbv	99
49) 1,1,2-Trichloroethane	17.932	97	268572	10.29	ppbv	99
50) Toluene	17.682	91	740907	9.99	ppbv	99
51) 2-Hexanone (MBK)	18.128	58	280958	10.63	ppbv	97
52) Dibromochloromethane	18.877	129	569001	11.35	ppbv	100
53) 1,2-Dibromoethane	19.233	107	438905	10.26	ppbv	100
54) Tetrachloroethene (PCE)	19.019	166	422591	10.17	ppbv	99
56) Chlorobenzene	20.357	114	201094	10.31	ppbv	99
57) Ethylbenzene	20.696	91	1023973	10.34	ppbv	100
58) m&p-Xylene	20.945	106	760135	19.51	ppbv	94
59) Bromoform	21.819	173	537006	10.27	ppbv #	96
60) Styrene	21.641	104	622726	9.88	ppbv	100
61) 1,1,2,2-Tetrachloroethane	22.336	83	575151	10.19	ppbv	99
62) o-Xylene	21.694	91	786079	9.92	ppbv	100
64) 4-Ethyltoluene	23.673	120	333591	10.26	ppbv	100
65) 1,3,5-Trimethylbenzene	23.780	120	456220	9.60	ppbv	99
66) 1,2,4-Trimethylbenzene	24.529	120	453390	9.79	ppbv	98
67) BenzylChloride (a-Chlor...)	25.154	91	749691	10.76	ppbv	99
68) 1,3-Dichlorobenzene	25.047	146	731849	10.05	ppbv	99
69) 1,4-Dichlorobenzene	25.261	146	693750m	9.47	ppbv	99
70) 1,2-Dichlorobenzene	25.831	146	734503m	9.64	ppbv	99
71) 1,2,4-Trichlorobenzene	29.433	180	713464m	9.47	ppbv	97
72) Hexachlorobutadiene	30.075	225	593007m	9.72	ppbv	99

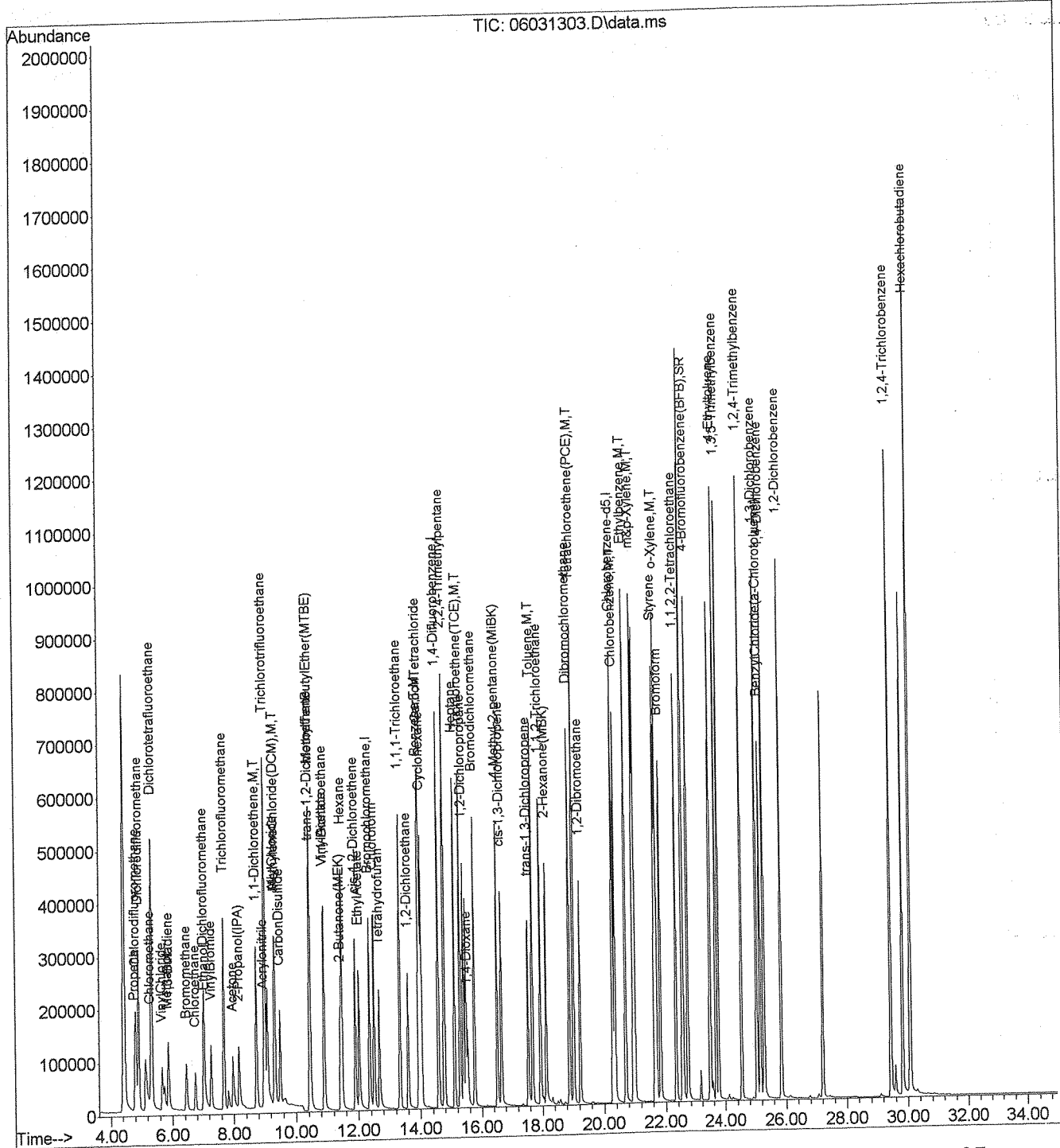
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Handwritten signature and date: JJG 6/3/13

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031303.D
 Acq On : 3 Jun 2013 10:03
 Operator : JJG
 Sample : TO15 LCSD 060313
 Misc : IS/Surr: PS082712-02 + Cal: PS040413-01
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Jun 03 12:32:31 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]

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031304.D
 Acq On : 3 Jun 2013 10:51
 Operator : JJG
 Sample : TO15 MB 060313
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Jun 03 11:50:22 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	145761	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	785623	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	751249	10.00	ppbv	0.00
System Monitoring Compounds						
63) 4-Bromofluorobenzene(BFB)	22.711	174	473021	10.06	ppbv	0.00
Spiked Amount	10.000		Recovery	= 100.60%		

Target Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)	Qvalue
2) Chlorodifluoromethane	0.000		0		N.D.		
3) Propene	4.817	42	224		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	5.903	31	1306		N.D.		
9) 1,3-Butadiene	0.000		0		N.D.		
10) Bromomethane	6.446	96	1169		N.D.	0.00	
11) Chloroethane	0.000		0		N.D.	0.00	
12) Dichlorofluoromethane	0.000		0		N.D.	0.00	
13) Ethanol	7.188	45	109		N.D.		
14) VinylBromide	0.000		0		N.D.		
15) Acetone	8.129	58	1454		N.D.	0.00	
16) Trichlorofluoromethane	0.000		0		N.D.		
17) 2-Propanol (IPA)	8.310	45	340		N.D.		
18) Acrylonitrile	0.000		0		N.D.		
19) 1,1-Dichloroethene	0.000		0		N.D.		
20) MethyleneChloride (DCM)	9.341	84	1490		N.D.		
21) AllylChloride	0.000		0		N.D.		
22) CarbonDisulfide	9.486	76	4581		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.		
26) MethylTertButylEther (M...)	0.000		0		N.D.		
27) VinylAcetate	0.000		0		N.D.		
28) 2-Butanone (MEK)	11.405	72	316		N.D.		
29) cis-1,2-Dichloroethene	0.000		0		N.D.		
30) Hexane	0.000		0		N.D.		
31) Chloroform	0.000		0		N.D.		
32) EthylAcetate	12.136	43	248		N.D.		

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031304.D
 Acq On : 3 Jun 2013 10:51
 Operator : JJG
 Sample : TO15 MB 060313
 Misc : IS/Surr: PS082712-02 + 500mL
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Jun 03 11:50:22 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

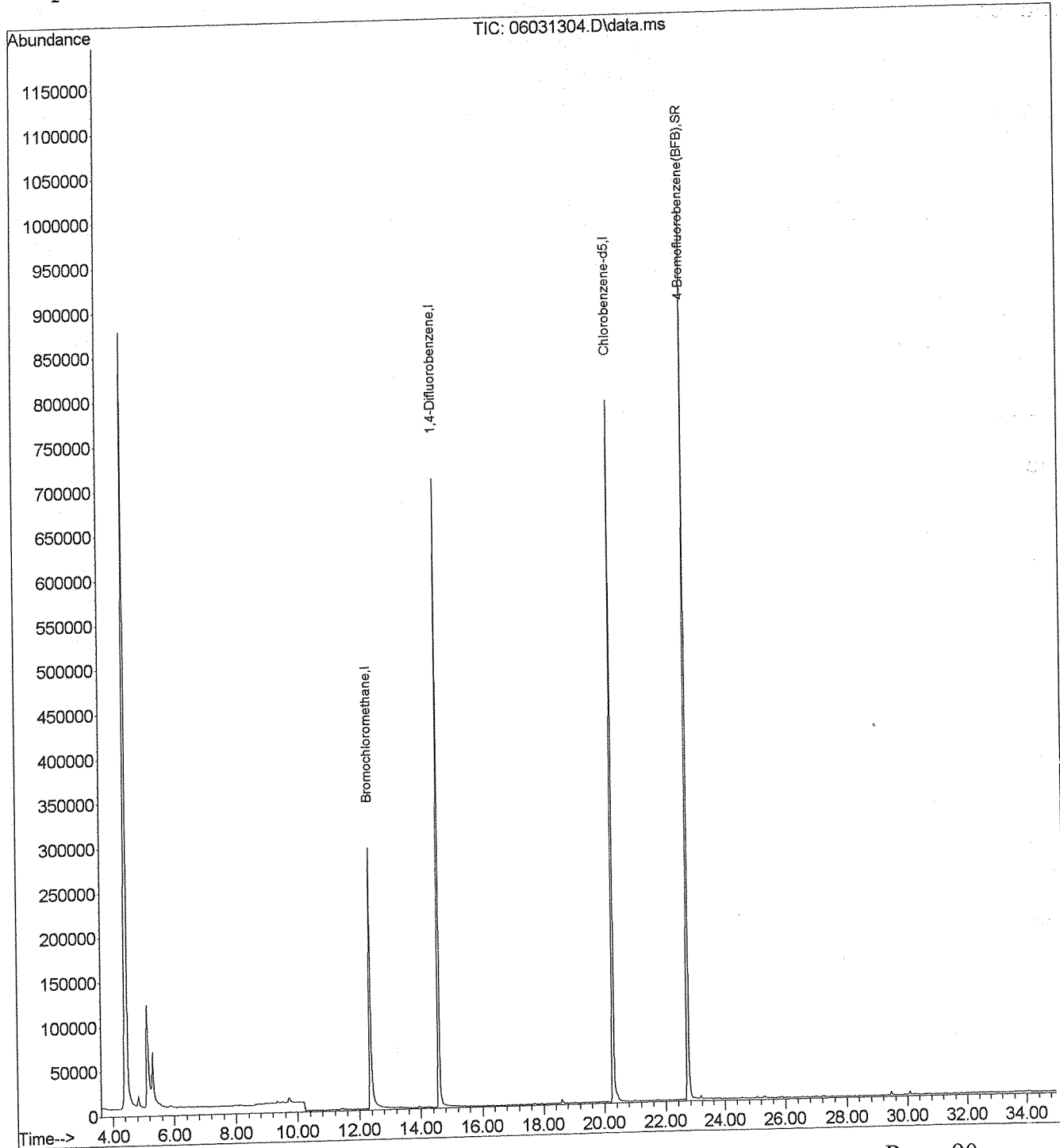
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) Tetrahydrofuran	0.000		0		N.D.	
34) 1,2-Dichloroethane	0.000		0		N.D.	
35) 1,1,1-Trichloroethane	0.000		0		N.D.	
37) Benzene	13.937	78	2965		N.D.	
38) CarbonTetrachloride	0.000		0		N.D.	
39) Cyclohexane	0.000		0		N.D.	
40) 1,2-Dichloropropane	15.292	63	113		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	17.682	91	1987		N.D.	
51) 2-Hexanone (MBK)	0.000		0		N.D.	
52) Dibromochloromethane	19.019	129	172		N.D.	
53) 1,2-Dibromoethane	0.000		0		N.D.	
54) Tetrachloroethene (PCE)	19.019	166	154		N.D.	
56) Chlorobenzene	20.375	114	112		N.D.	
57) Ethylbenzene	20.713	91	987		N.D.	
58) m&p-Xylene	21.016	106	517		N.D.	
59) Bromoform	0.000		0		N.D.	
60) Styrene	21.694	104	562		N.D.	
61) 1,1,2,2-Tetrachloroethane	0.000		0		N.D.	
62) o-Xylene	21.712	91	536		N.D.	
64) 4-Ethyltoluene	23.709	120	151		N.D.	
65) 1,3,5-Trimethylbenzene	23.798	120	181		N.D.	
66) 1,2,4-Trimethylbenzene	24.565	120	266		N.D.	
67) BenzylChloride (a-Chlor...)	25.207	91	854		N.D.	
68) 1,3-Dichlorobenzene	25.064	146	1959		N.D.	
69) 1,4-Dichlorobenzene	25.296	146	2252		N.D.	
70) 1,2-Dichlorobenzene	25.867	146	1319		N.D.	
71) 1,2,4-Trichlorobenzene	29.469	180	3546		N.D.	
72) Hexachlorobutadiene	30.075	225	1269		N.D.	

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

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\MS03\2013\060313\
Data File : 06031304.D
Acq On : 3 Jun 2013 10:51
Operator : JJG
Sample : TO15 MB 060313
Misc : IS/Surr: PS082712-02 + 500mL
ALS Vial : 1 Sample Multiplier: 1

Quant Time: Jun 03 11:50:22 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



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031305.D
 Acq On : 3 Jun 2013 11:37
 Operator : JJG
 Sample : 130650-63200 x10
 Misc : IS/Surr: PS082712-02 + 50mL
 ALS Vial : 2 Sample Multiplier: 10

Quant Time: Jun 03 12:38:11 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	145756	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	793394	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	753276	10.00	ppbv	0.00
System Monitoring Compounds						
63) 4-Bromofluorobenzene (BFB)	22.710	174	480401	10.19	ppbv	0.00
Spiked Amount	10.000		Recovery	= 101.90%		

Target Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)	Qvalue
2) Chlorodifluoromethane	4.835	51	566	N.D.			
3) Propene	4.817	42	841	N.D.			
4) Dichlorodifluoromethane	4.926	85	1412	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	5.867	31	383066	82.26	ppbv		
9) 1,3-Butadiene	0.000		0	N.D.			
10) Bromomethane	6.446	96	701	N.D.		0.00	
11) Chloroethane	0.000		0	N.D.		0.00	
12) Dichlorofluoromethane	0.000		0	N.D.		0.00	
13) Ethanol	7.152	45	5921	N.D.			
14) VinylBromide	0.000		0	N.D.			
15) Acetone	8.057	58	6425	N.D.		0.00	
16) Trichlorofluoromethane	7.658	103	339	N.D.			
17) 2-Propanol (IPA)	8.256	45	6108	N.D.			
18) Acrylonitrile	0.000		0	N.D.			
19) 1,1-Dichloroethene	0.000		0	N.D.			
20) MethyleneChloride (DCM)	9.341	84	1281	N.D.			
21) AllylChloride	0.000		0	N.D.			
22) CarbonDisulfide	9.504	76	3786	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.			
26) MethylTertButylEther (M...)	0.000		0	N.D.			
27) VinylAcetate	10.888	43	120	N.D.			
28) 2-Butanone (MEK)	11.512	72	1776	N.D.			
29) cis-1,2-Dichloroethene	0.000		0	N.D.			
30) Hexane	0.000		0	N.D.			
31) Chloroform	0.000		0	N.D.			
32) EthylAcetate	12.118	43	1289	N.D.			

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031305.D
 Acq On : 3 Jun 2013 11:37
 Operator : JJG
 Sample : 130650-63200 x10
 Misc : IS/Surr: PS082712-02 + 50mL
 ALS Vial : 2 Sample Multiplier: 10

Quant Time: Jun 03 12:38:11 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

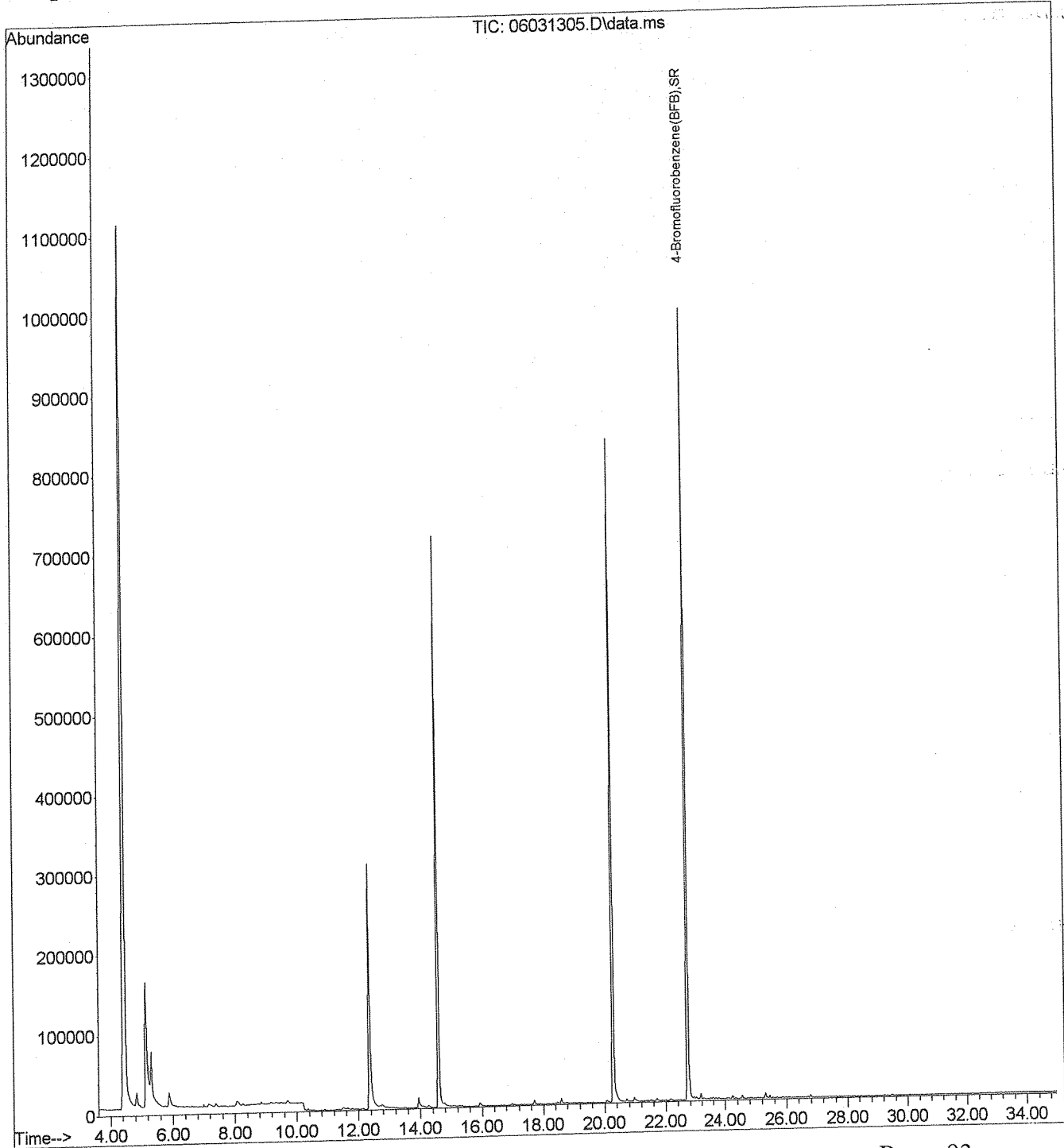
Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
33) Tetrahydrofuran	12.760	72	1372			N.D.
34) 1,2-Dichloroethane	0.000		0			N.D.
35) 1,1,1-Trichloroethane	0.000		0			N.D.
37) Benzene	13.937	78	15871			N.D.
38) CarbonTetrachloride	0.000		0			N.D.
39) Cyclohexane	0.000		0			N.D.
40) 1,2-Dichloropropane	15.275	63	113			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	14.775	57	4185			N.D.
45) Heptane	15.114	71	125			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	17.878	97	295			N.D.
50) Toluene	17.700	91	7109			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	20.267	114	234			N.D.
57) Ethylbenzene	20.713	91	3358			N.D.
58) m&p-Xylene	20.963	106	3210			N.D.
59) Bromoform	0.000		0			N.D.
60) Styrene	21.694	104	109			N.D.
61) 1,1,2,2-Tetrachloroethane	0.000		0			N.D.
62) o-Xylene	21.712	91	2881			N.D.
64) 4-Ethyltoluene	23.691	120	354			N.D.
65) 1,3,5-Trimethylbenzene	23.798	120	657			N.D.
66) 1,2,4-Trimethylbenzene	24.547	120	1686			N.D.
67) BenzylChloride (a-Chlor...)	25.225	91	110			N.D.
68) 1,3-Dichlorobenzene	25.064	146	372			N.D.
69) 1,4-Dichlorobenzene	25.296	146	1229			N.D.
70) 1,2-Dichlorobenzene	25.849	146	353			N.D.
71) 1,2,4-Trichlorobenzene	29.469	180	1019			N.D.
72) Hexachlorobutadiene	30.075	225	358			N.D.

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

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\MS03\2013\060313\
Data File : 06031305.D
Acq On : 3 Jun 2013 11:37
Operator : JJG
Sample : 130650-63200 x10
Misc : IS/Surr: PS082712-02 + 50mL
ALS Vial : 2 Sample Multiplier: 10

Quant Time: Jun 03 12:38:11 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



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031306.D
 Acq On : 3 Jun 2013 12:25
 Operator : JJG
 Sample : 130650-63200 x10 dp
 Misc : IS/Surr: PS082712-02 + 50mL
 ALS Vial : 2 Sample Multiplier: 10

Quant Time: Jun 03 12:57:49 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	146289	10.00	ppbv	0.00
36) 1,4-Difluorobenzene	14.579	114	808196	10.00	ppbv	0.00
55) Chlorobenzene-d5	20.285	117	747587	10.00	ppbv	0.00
System Monitoring Compounds						
63) 4-Bromofluorobenzene (BFB)	22.710	174	482157	10.31	ppbv	0.00
Spiked Amount	10.000		Recovery	= 103.10%		

Target Compounds	R.T.	QIon	Response	Conc	Units	Dev (Min)	Ovalue
2) Chlorodifluoromethane	4.836	51	600		N.D.		
3) Propene	4.818	42	909		N.D.		
4) Dichlorodifluoromethane	4.926	85	1602		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	5.867	31	381220	81.54	ppbv		
9) 1,3-Butadiene	0.000		0		N.D.		
10) Bromomethane	6.446	96	592		N.D.		
11) Chloroethane	0.000		0		N.D.		
12) Dichlorofluoromethane	0.000		0		N.D.		
13) Ethanol	0.000		0		N.D.		d
14) VinylBromide	0.000		0		N.D.		
15) Acetone	8.057	58	6380		N.D.		
16) Trichlorofluoromethane	7.659	103	412		N.D.		
17) 2-Propanol (IPA)	8.256	45	6139		N.D.		
18) Acrylonitrile	0.000		0		N.D.		
19) 1,1-Dichloroethene	0.000		0		N.D.		
20) MethyleneChloride (DCM)	9.324	84	1477		N.D.		
21) AllylChloride	0.000		0		N.D.		
22) CarbonDisulfide	9.504	76	3692		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.		
26) MethylTertButyleEther (M...)	0.000		0		N.D.		
27) VinylAcetate	10.906	43	118		N.D.		
28) 2-Butanone (MEK)	11.512	72	1755		N.D.		
29) cis-1,2-Dichloroethene	0.000		0		N.D.		
30) Hexane	0.000		0		N.D.		
31) Chloroform	0.000		0		N.D.		
32) EthylAcetate	12.118	43	1247		N.D.		

Data Path : C:\msdchem\1\MS03\2013\060313\
 Data File : 06031306.D
 Acq On : 3 Jun 2013 12:25
 Operator : JJG
 Sample : 130650-63200 x10 dp
 Misc : IS/Surr: PS082712-02 + 50mL
 ALS Vial : 2 Sample Multiplier: 10

Quant Time: Jun 03 12:57:49 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

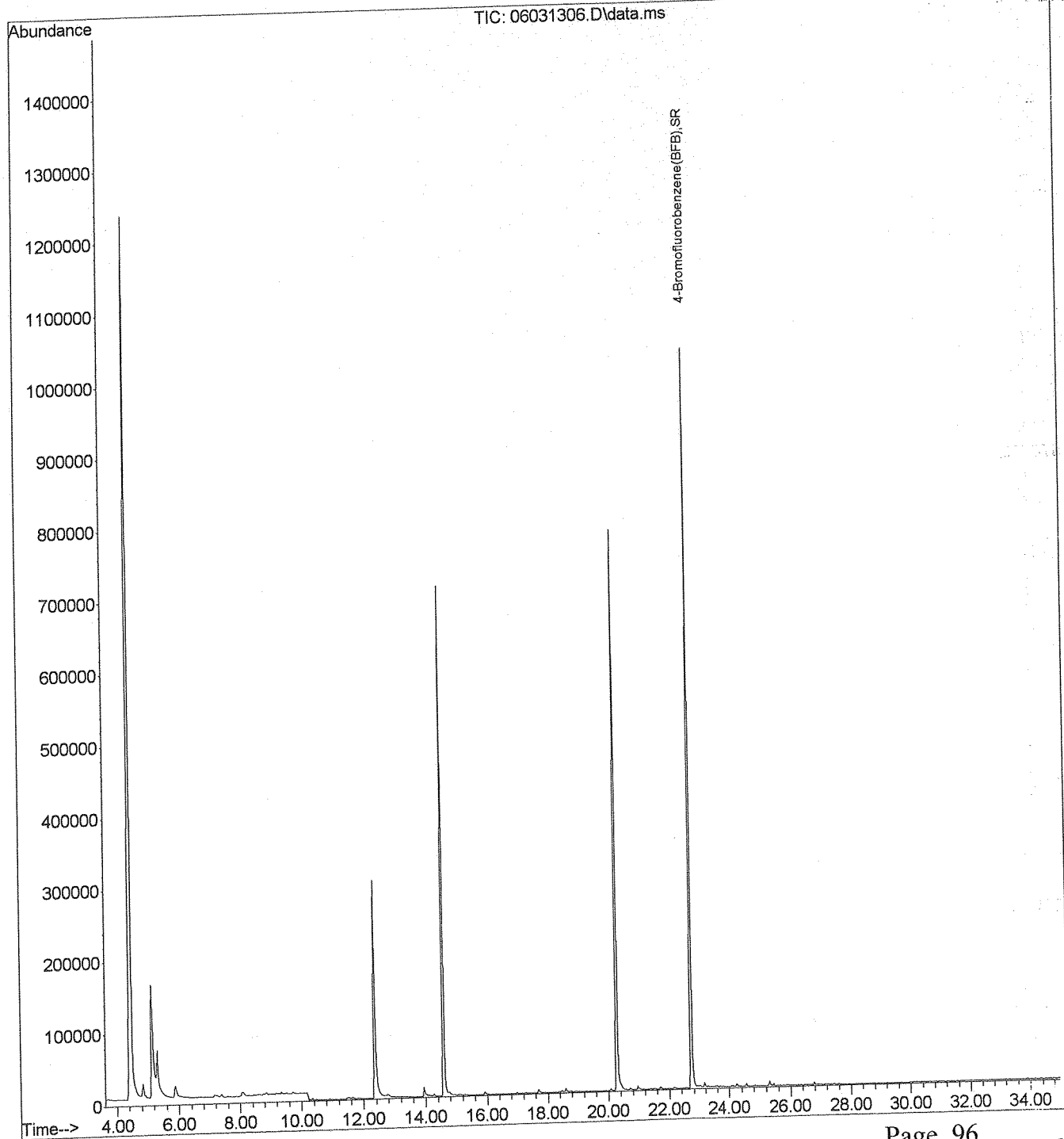
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) Tetrahydrofuran	12.760	72	1398			N.D.
34) 1,2-Dichloroethane	0.000		0			N.D.
35) 1,1,1-Trichloroethane	0.000		0			N.D.
37) Benzene	13.937	78	15751			N.D.
38) CarbonTetrachloride	0.000		0			N.D.
39) Cyclohexane	0.000		0			N.D.
40) 1,2-Dichloropropane	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	14.757	57	4392			N.D.
45) Heptane	15.114	71	262			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	17.860	97	402			N.D.
50) Toluene	17.682	91	6905			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	20.267	114	109			N.D.
57) Ethylbenzene	20.713	91	2943			N.D.
58) m&p-Xylene	20.963	106	3337			N.D.
59) Bromoform	0.000		0			N.D.
60) Styrene	0.000		0			N.D.
61) 1,1,2,2-Tetrachloroethane	0.000		0			N.D.
62) o-Xylene	21.712	91	2770			N.D.
64) 4-Ethyltoluene	23.691	120	163			N.D.
65) 1,3,5-Trimethylbenzene	23.798	120	361			N.D.
66) 1,2,4-Trimethylbenzene	24.547	120	1588			N.D.
67) BenzylChloride (a-Chlor...	25.296	91	1280			N.D.
68) 1,3-Dichlorobenzene	0.000		0			N.D.
69) 1,4-Dichlorobenzene	25.296	146	899			N.D.
70) 1,2-Dichlorobenzene	25.867	146	138			N.D.
71) 1,2,4-Trichlorobenzene	29.486	180	373			N.D.
72) Hexachlorobutadiene	30.075	225	142			N.D.

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

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\MS03\2013\060313\
Data File : 06031306.D
Acq On : 3 Jun 2013 12:25
Operator : JJG
Sample : 130650-63200 x10 dp
Misc : IS/Surr: PS082712-02 + 50mL
ALS Vial : 2 Sample Multiplier: 10

Quant Time: Jun 03 12:57:49 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

Calibration Status 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

#	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

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
-----ISTD-----									
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	7.88
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)									

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

		ISTD										
29)	cis-1,2-Dichlo...	1.339	1.387	1.385	1.382	1.308	1.264	1.167	1.319			6.17
30)	Hexane	0.270	0.298	0.295	0.272	0.256	0.239	0.219	0.264			10.89
31)	Chloroform	2.980	3.005	2.926	2.844	2.659	2.617	2.396	2.775			8.12
32)	EthylAcetate	2.836	3.052	2.959	2.914	2.819	2.582	2.380	2.792			8.36
33)	Tetrahydrofuran	0.574	0.569	0.601	0.585	0.559	0.535	0.483	0.558			7.01
34)	1,2-Dichloroet...	2.043	2.248	2.117	2.034	1.930	1.874	1.847	2.013			7.07
35)	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)	CarbonTetrachl...	0.664	0.694	0.674	0.620	0.600	0.541	0.469	0.609			13.26
39)	Cyclohexane	0.121	0.121	0.117	0.111	0.109	0.098	0.087	0.109			11.74
40)	1,2-Dichloropr...	0.324	0.333	0.328	0.306	0.291	0.265	0.226	0.296			13.14
41)	Bromodichlorom...	0.411	0.420	0.409	0.393	0.371	0.346	0.311	0.380			10.47
42)	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)	2,2,4-Trimethy...	1.494	1.513	1.438	1.352	1.312	1.136	0.935	1.312			16.01
45)	Heptane	0.250	0.247	0.252	0.241	0.240	0.218	0.189	0.234			9.72
46)	cis-1,3-Dichlo...	0.443	0.456	0.446	0.443	0.425	0.389	0.351	0.422			9.07
47)	4-Methyl-2-pen...	0.264	0.293	0.298	0.287	0.277	0.248	0.212	0.269			11.25
48)	trans-1,3-Dich...	0.406	0.438	0.451	0.457	0.441	0.423	0.391	0.430			5.59
49)	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)	2-Hexanone (MBK)	0.339	0.353	0.364	0.351	0.351	0.307	0.266	0.333			10.43
52)	Dibromochlorom...	0.667	0.684	0.678	0.663	0.647	0.577	0.504	0.632			10.56
53)	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)	Bromoform	0.762	0.809	0.762	0.719	0.685	0.627	0.514	0.697			14.37
60)	Styrene	0.895	0.947	0.887	0.890	0.834	0.768	0.660	0.840			11.58
61)	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)	4-Ethyltoluene	0.453	0.499	0.463	0.465	0.439	0.394	0.321	0.433			13.60

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