

ANALYTICAL REPORT

TestAmerica Laboratories, Inc.

TestAmerica Canton
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Tel: (330)497-9396

TestAmerica Job ID: 240-109102-1

Client Project/Site: Ford LTP Livonia MI - E203631

For:

ARCADIS U.S., Inc.
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Suite 500
Novi, Michigan 48377

Attn: Kristoffer Hinskey



Authorized for release by:
3/22/2019 9:38:24 AM

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Results relate only to the items tested and the sample(s) as received by the laboratory.



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Definitions/Glossary

Client: ARCADIS U.S., Inc.
Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

Qualifiers

GC/MS VOA

Qualifier	Qualifier Description
X	Surrogate is outside control limits
F1	MS and/or MSD Recovery is outside acceptance limits.
J	Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value.
U	Indicates the analyte was analyzed for but not detected.
B	Compound was found in the blank and sample.
*	LCS or LCSD is outside acceptance limits.
F2	MS/MSD RPD exceeds control limits
^	ICV,CCV,ICB,CCB, ISA, ISB, CRI, CRA, DLCK or MRL standard: Instrument related QC is outside acceptance limits.

Glossary

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

Case Narrative

Client: ARCADIS U.S., Inc.
Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

Job ID: 240-109102-1

Laboratory: TestAmerica Canton

Narrative

CASE NARRATIVE

Client: ARCADIS U.S., Inc.

Project: Ford LTP Livonia MI - E203631

Report Number: 240-109102-1

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

TestAmerica Canton attests to the validity of the laboratory data generated by TestAmerica facilities reported herein. All analyses performed by TestAmerica facilities were done using established laboratory SOPs that incorporate QA/QC procedures described in the application methods. TestAmerica's operations groups have reviewed the data for compliance with the laboratory QA/QC plan, and data have been found to be compliant with laboratory protocols unless otherwise noted below.

The test results in this report meet all NELAP requirements for parameters for which accreditation is required or available. Any exceptions to NELAP requirements are noted in this report. Pursuant to NELAP, this report may not be reproduced, except in full, without the written approval of the laboratory.

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

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

All solid sample results are reported on an "as received" basis unless otherwise indicated by the presence of a % solids value in the method header.

This laboratory report is confidential and is intended for the sole use of TestAmerica and its client.

RECEIPT

The samples were received on 3/8/2019 8:00 AM; the samples arrived in good condition, properly preserved and, where required, on ice. The temperature of the cooler at receipt was 3.2° C.

VOLATILE ORGANIC COMPOUNDS (GCMS)

Samples MW-72_030519 (240-109102-1), MW-72S_030519 (240-109102-2), MW-83_030519 (240-109102-3), MW-83S_030519 (240-109102-4), MW-82SR_030519 (240-109102-5) and TRIP BLANK (240-109102-6) were analyzed for volatile organic compounds (GCMS) in accordance with EPA SW-846 Method 8260B. The samples were analyzed on 03/14/2019 and 03/18/2019.

cis-1,2-Dichloroethene and Trichloroethene were detected in method blank MB 240-371553/6 at levels that were above the method detection limit but below the reporting limit. The values should be considered estimates, and have been flagged. If the associated sample reported a result above the MDL and/or RL, the result has been flagged. Refer to the QC report for details.

Vinyl chloride failed the recovery criteria high for LCS 240-371553/4. Vinyl chloride failed the recovery criteria high for MRL 240-371553/5. Refer to the QC report for details.

The continuing calibration verification (CCV) associated with batch 371553 recovered above the upper control limit for Vinyl Chloride and/or 1,2-Dichloropropane. The samples associated with this CCV were non-detects for the affected analytes; therefore, the data have been reported. The following samples are impacted: MW-72S_030519 (240-109102-2), MW-83_030519 (240-109102-3), MW-83S_030519 (240-109102-4), MW-82SR_030519 (240-109102-5) and TRIP BLANK (240-109102-6).

Case Narrative

Client: ARCADIS U.S., Inc.
Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

Job ID: 240-109102-1 (Continued)

Laboratory: TestAmerica Canton (Continued)

The laboratory control sample (LCS) for 371553 recovered outside control limits one or multiple analytes. These analytes were biased high in the LCS and were not detected in the associated samples; therefore, the data have been reported: MW-72S_030519 (240-109102-2), MW-83_030519 (240-109102-3), MW-83S_030519 (240-109102-4), MW-82SR_030519 (240-109102-5), TRIP BLANK (240-109102-6) and (LCS 240-371553/4).

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

VOLATILE ORGANIC COMPOUNDS (GCMS SIM)

Samples MW-72_030519 (240-109102-1), MW-72S_030519 (240-109102-2), MW-83_030519 (240-109102-3), MW-83S_030519 (240-109102-4) and MW-82SR_030519 (240-109102-5) were analyzed for volatile organic compounds (GCMS SIM) in accordance with EPA SW-846 Method 8260B SIM. The samples were analyzed on 03/13/2019.

The pH is greater than 2 for the following sample: MW-83_030519 (240-109102-3).

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.



Method Summary

Client: ARCADIS U.S., Inc.
Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

Method	Method Description	Protocol	Laboratory
8260B	Volatile Organic Compounds (GC/MS)	SW846	TAL CAN
8260B SIM	Volatile Organic Compounds (GC/MS)	SW846	TAL CAN
5030B	Purge and Trap	SW846	TAL CAN

Protocol References:

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

Laboratory References:

TAL CAN = TestAmerica Canton, 4101 Shuffel Street NW, North Canton, OH 44720, TEL (330)497-9396



Sample Summary

Client: ARCADIS U.S., Inc.
Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

Lab Sample ID	Client Sample ID	Matrix	Collected	Received
240-109102-1	MW-72_030519	Water	03/05/19 14:20	03/08/19 08:00
240-109102-2	MW-72S_030519	Water	03/05/19 15:50	03/08/19 08:00
240-109102-3	MW-83_030519	Water	03/05/19 12:13	03/08/19 08:00
240-109102-4	MW-83S_030519	Water	03/05/19 14:07	03/08/19 08:00
240-109102-5	MW-82SR_030519	Water	03/05/19 17:16	03/08/19 08:00
240-109102-6	TRIP BLANK	Water	03/05/19 00:00	03/08/19 08:00

- 1
- 2
- 3
- 4
- 5
- 6
- 7
- 8
- 9
- 10
- 11
- 12
- 13
- 14

Detection Summary

Client: ARCADIS U.S., Inc.
Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

Client Sample ID: MW-72_030519

Lab Sample ID: 240-109102-1

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
1,4-Dioxane	0.91	J	2.0	0.86	ug/L	1		8260B SIM	Total/NA
Vinyl chloride	1.9		1.0	0.20	ug/L	1		8260B	Total/NA

Client Sample ID: MW-72S_030519

Lab Sample ID: 240-109102-2

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Trichloroethene	0.20	J B	1.0	0.10	ug/L	1		8260B	Total/NA
Vinyl chloride	0.24	J *	1.0	0.20	ug/L	1		8260B	Total/NA

Client Sample ID: MW-83_030519

Lab Sample ID: 240-109102-3

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Trichloroethene	0.17	J B	1.0	0.10	ug/L	1		8260B	Total/NA

Client Sample ID: MW-83S_030519

Lab Sample ID: 240-109102-4

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Trichloroethene	0.20	J B	1.0	0.10	ug/L	1		8260B	Total/NA
Vinyl chloride	0.27	J *	1.0	0.20	ug/L	1		8260B	Total/NA

Client Sample ID: MW-82SR_030519

Lab Sample ID: 240-109102-5

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Trichloroethene	0.16	J B	1.0	0.10	ug/L	1		8260B	Total/NA
Vinyl chloride	0.27	J *	1.0	0.20	ug/L	1		8260B	Total/NA

Client Sample ID: TRIP BLANK

Lab Sample ID: 240-109102-6

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Trichloroethene	0.16	J B	1.0	0.10	ug/L	1		8260B	Total/NA

This Detection Summary does not include radiochemical test results.

TestAmerica Canton

Client Sample Results

Client: ARCADIS U.S., Inc.
 Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

Client Sample ID: MW-72_030519

Lab Sample ID: 240-109102-1

Date Collected: 03/05/19 14:20

Matrix: Water

Date Received: 03/08/19 08:00

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

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,4-Dioxane	0.91	J	2.0	0.86	ug/L			03/13/19 12:28	1
Surrogate	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	98		63 - 125					03/13/19 12:28	1

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

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/18/19 16:27	1
cis-1,2-Dichloroethene	1.0	U	1.0	0.16	ug/L			03/18/19 16:27	1
Tetrachloroethene	1.0	U	1.0	0.15	ug/L			03/18/19 16:27	1
trans-1,2-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/18/19 16:27	1
Trichloroethene	1.0	U	1.0	0.10	ug/L			03/18/19 16:27	1
Vinyl chloride	1.9		1.0	0.20	ug/L			03/18/19 16:27	1
Surrogate	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	89		70 - 121					03/18/19 16:27	1
4-Bromofluorobenzene (Surr)	68		59 - 120					03/18/19 16:27	1
Toluene-d8 (Surr)	79		70 - 123					03/18/19 16:27	1
Dibromofluoromethane (Surr)	91		75 - 128					03/18/19 16:27	1

Client Sample Results

Client: ARCADIS U.S., Inc.
 Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

Client Sample ID: MW-72S_030519

Lab Sample ID: 240-109102-2

Date Collected: 03/05/19 15:50

Matrix: Water

Date Received: 03/08/19 08:00

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

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,4-Dioxane	2.0	U	2.0	0.86	ug/L			03/13/19 12:53	1
Surrogate	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	99		63 - 125					03/13/19 12:53	1

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

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/14/19 18:59	1
cis-1,2-Dichloroethene	1.0	U	1.0	0.16	ug/L			03/14/19 18:59	1
Tetrachloroethene	1.0	U	1.0	0.15	ug/L			03/14/19 18:59	1
trans-1,2-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/14/19 18:59	1
Trichloroethene	0.20	J B	1.0	0.10	ug/L			03/14/19 18:59	1
Vinyl chloride	0.24	J *	1.0	0.20	ug/L			03/14/19 18:59	1
Surrogate	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	113		70 - 121					03/14/19 18:59	1
4-Bromofluorobenzene (Surr)	100		59 - 120					03/14/19 18:59	1
Toluene-d8 (Surr)	114		70 - 123					03/14/19 18:59	1
Dibromofluoromethane (Surr)	98		75 - 128					03/14/19 18:59	1

Client Sample Results

Client: ARCADIS U.S., Inc.
 Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

Client Sample ID: MW-83_030519

Lab Sample ID: 240-109102-3

Date Collected: 03/05/19 12:13

Matrix: Water

Date Received: 03/08/19 08:00

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

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,4-Dioxane	2.0	U	2.0	0.86	ug/L			03/13/19 13:19	1
Surrogate	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	101		63 - 125					03/13/19 13:19	1

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

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/14/19 19:22	1
cis-1,2-Dichloroethene	1.0	U	1.0	0.16	ug/L			03/14/19 19:22	1
Tetrachloroethene	1.0	U	1.0	0.15	ug/L			03/14/19 19:22	1
trans-1,2-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/14/19 19:22	1
Trichloroethene	0.17	J B	1.0	0.10	ug/L			03/14/19 19:22	1
Vinyl chloride	1.0	U *	1.0	0.20	ug/L			03/14/19 19:22	1
Surrogate	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	109		70 - 121					03/14/19 19:22	1
4-Bromofluorobenzene (Surr)	94		59 - 120					03/14/19 19:22	1
Toluene-d8 (Surr)	113		70 - 123					03/14/19 19:22	1
Dibromofluoromethane (Surr)	98		75 - 128					03/14/19 19:22	1

Client Sample Results

Client: ARCADIS U.S., Inc.
 Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

Client Sample ID: MW-83S_030519

Lab Sample ID: 240-109102-4

Date Collected: 03/05/19 14:07

Matrix: Water

Date Received: 03/08/19 08:00

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

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,4-Dioxane	2.0	U	2.0	0.86	ug/L			03/13/19 13:45	1
Surrogate	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	98		63 - 125					03/13/19 13:45	1

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

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/14/19 19:44	1
cis-1,2-Dichloroethene	1.0	U	1.0	0.16	ug/L			03/14/19 19:44	1
Tetrachloroethene	1.0	U	1.0	0.15	ug/L			03/14/19 19:44	1
trans-1,2-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/14/19 19:44	1
Trichloroethene	0.20	J B	1.0	0.10	ug/L			03/14/19 19:44	1
Vinyl chloride	0.27	J *	1.0	0.20	ug/L			03/14/19 19:44	1
Surrogate	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	105		70 - 121					03/14/19 19:44	1
4-Bromofluorobenzene (Surr)	88		59 - 120					03/14/19 19:44	1
Toluene-d8 (Surr)	104		70 - 123					03/14/19 19:44	1
Dibromofluoromethane (Surr)	95		75 - 128					03/14/19 19:44	1

Client Sample Results

Client: ARCADIS U.S., Inc.
 Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

Client Sample ID: MW-82SR_030519

Lab Sample ID: 240-109102-5

Date Collected: 03/05/19 17:16

Matrix: Water

Date Received: 03/08/19 08:00

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

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,4-Dioxane	2.0	U	2.0	0.86	ug/L			03/13/19 14:11	1
Surrogate	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	101		63 - 125					03/13/19 14:11	1

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

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/14/19 20:06	1
cis-1,2-Dichloroethene	1.0	U	1.0	0.16	ug/L			03/14/19 20:06	1
Tetrachloroethene	1.0	U	1.0	0.15	ug/L			03/14/19 20:06	1
trans-1,2-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/14/19 20:06	1
Trichloroethene	0.16	J B	1.0	0.10	ug/L			03/14/19 20:06	1
Vinyl chloride	0.27	J *	1.0	0.20	ug/L			03/14/19 20:06	1
Surrogate	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	110		70 - 121					03/14/19 20:06	1
4-Bromofluorobenzene (Surr)	90		59 - 120					03/14/19 20:06	1
Toluene-d8 (Surr)	109		70 - 123					03/14/19 20:06	1
Dibromofluoromethane (Surr)	96		75 - 128					03/14/19 20:06	1

Client Sample Results

Client: ARCADIS U.S., Inc.
 Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

Client Sample ID: TRIP BLANK

Lab Sample ID: 240-109102-6

Date Collected: 03/05/19 00:00

Matrix: Water

Date Received: 03/08/19 08:00

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

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/14/19 20:28	1
cis-1,2-Dichloroethene	1.0	U	1.0	0.16	ug/L			03/14/19 20:28	1
Tetrachloroethene	1.0	U	1.0	0.15	ug/L			03/14/19 20:28	1
trans-1,2-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/14/19 20:28	1
Trichloroethene	0.16	J B	1.0	0.10	ug/L			03/14/19 20:28	1
Vinyl chloride	1.0	U *	1.0	0.20	ug/L			03/14/19 20:28	1

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	113		70 - 121		03/14/19 20:28	1
4-Bromofluorobenzene (Surr)	91		59 - 120		03/14/19 20:28	1
Toluene-d8 (Surr)	110		70 - 123		03/14/19 20:28	1
Dibromofluoromethane (Surr)	97		75 - 128		03/14/19 20:28	1

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: ICV 240-339391/11 Calibration Date: 08/03/2018 19:55
 Instrument ID: A3UX2 Calib Start Date: 08/03/2018 16:37
 GC Column: ZB-624 ID: 0.53 (mm) Calib End Date: 08/03/2018 19:31
 Lab File ID: X23824.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,4-Dioxane	Lin1		1.046		0.00898	0.0100	-10.2	
1,2-Dichloroethane-d4 (Surr)	Ave	0.3184	0.3275		0.0103	0.0100	2.9	

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23824.D
 Lims ID: ICV
 Client ID:
 Sample Type: ICV
 Inject. Date: 03-Aug-2018 19:55:30 ALS Bottle#: 10 Worklist Smp#: 11
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: icv
 Misc. Info.: 240-0078614-011
 Operator ID: 002808 Instrument ID: A3UX2
 Sublist:
 Method: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 07-Aug-2018 13:50:40 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK012

First Level Reviewer: macenczaks Date: 06-Aug-2018 08:48:54

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.053	4.056	-0.003	100	1953113	10.0	10.3	
* 7 Fluorobenzene	96	4.623	4.623	0.000	100	5963265	10.0	10.0	
* 9 Dioxane-d8 (IS)	96	5.899	5.899	0.000	99	130933	200.0	200.0	M
8 1,4-Dioxane	88	5.977	5.969	0.008	93	6851	10.0	8.98	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

vmdiox_spike_00154	Amount Added: 3.00	Units: uL	
vmDist_H2o_00121	Amount Added: 0.00	Units:	Run Reagent
vm150is_00117	Amount Added: 1.00	Units: uL	Run Reagent

TestAmerica Canton

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23824.D

Injection Date: 03-Aug-2018 19:55:30

Instrument ID: A3UX2

Operator ID: 002808

Lims ID: ICV

Worklist Smp#: 11

Client ID:

Purge Vol: 15.000 mL

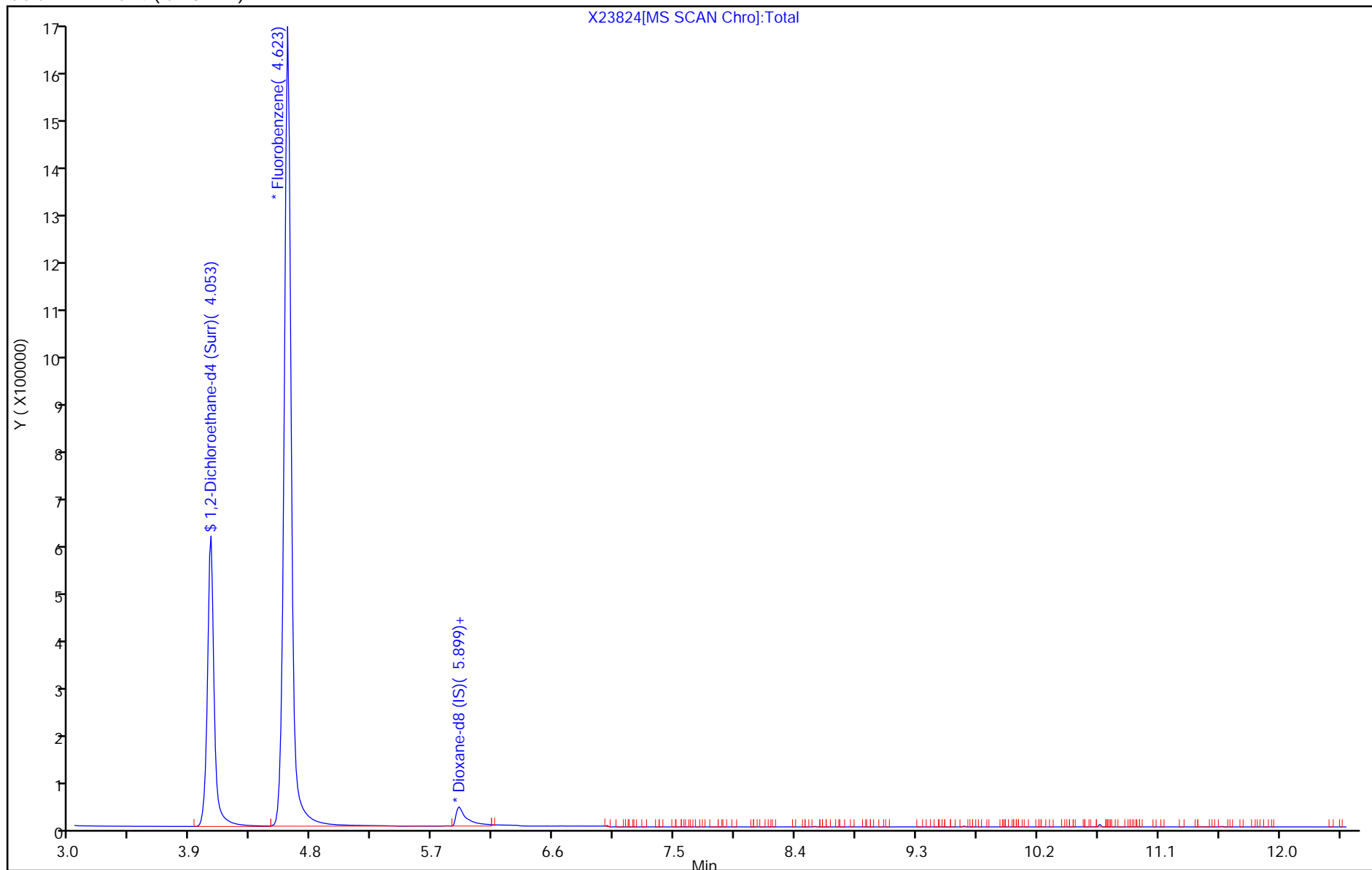
Dil. Factor: 1.0000

ALS Bottle#: 10

Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton

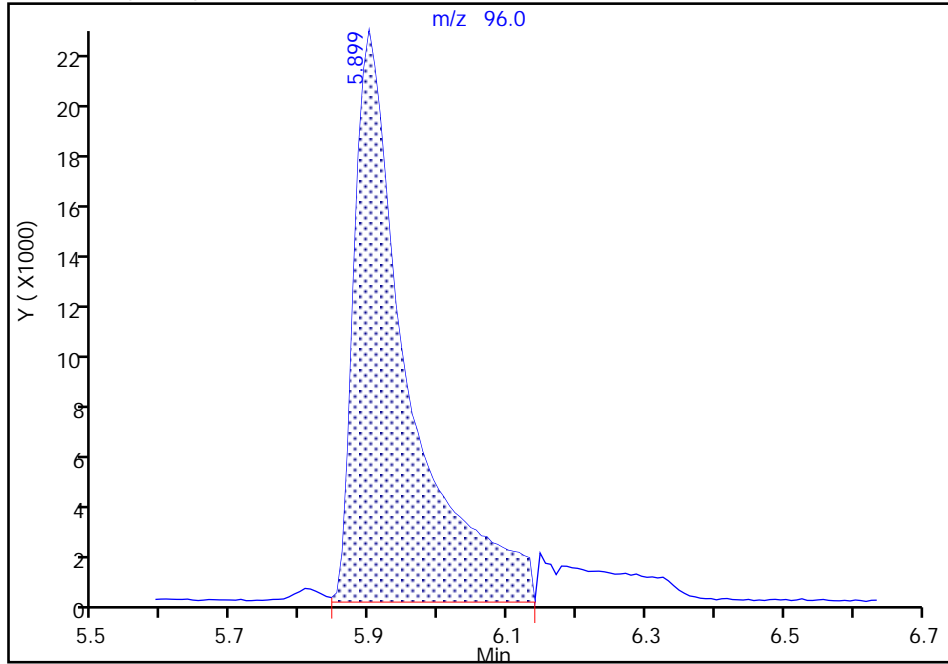
Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23824.D
Injection Date: 03-Aug-2018 19:55:30 Instrument ID: A3UX2
Lims ID: ICV
Client ID:
Operator ID: 002808 ALS Bottle#: 10 Worklist Smp#: 11
Purge Vol: 15.000 mL Dil. Factor: 1.0000
Method: 8260SIM_X2 Limit Group: MSV SIM ICAL
Column: DB-624 (0.18 mm) Detector: MS SCAN

* 9 Dioxane-d8 (IS), CAS: 17647-74-4

Signal: 1

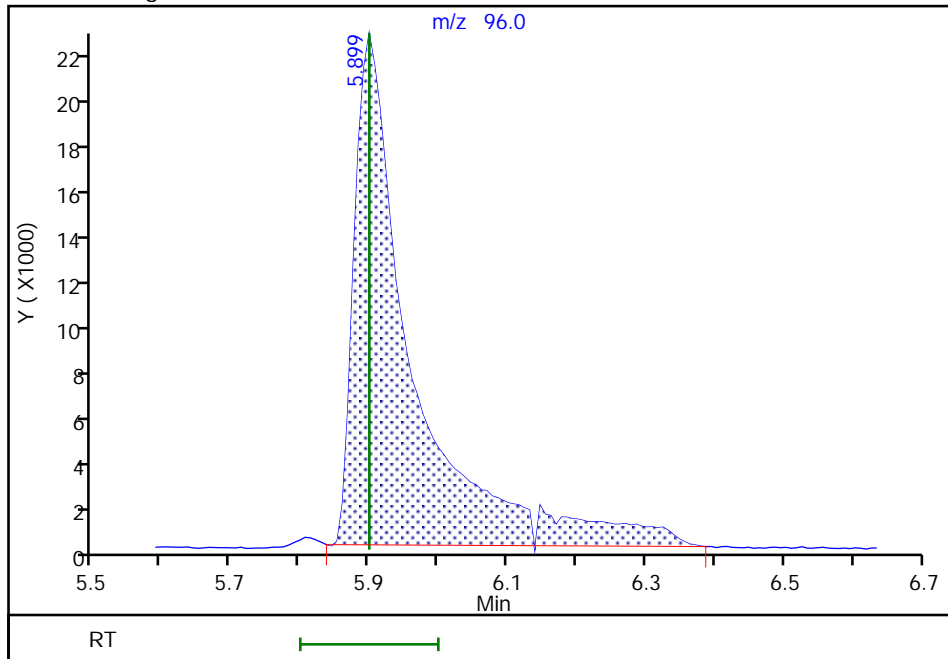
RT: 5.90
Area: 121822
Amount: 200.0000
Amount Units: ug/l

Processing Integration Results



RT: 5.90
Area: 130933
Amount: 200.0000
Amount Units: ug/l

Manual Integration Results



Reviewer: macenczaks, 06-Aug-2018 08:48:32

Audit Action: Manually Integrated

Audit Reason: Poor chromatography



FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: CCVIS 240-371373/3 Calibration Date: 03/13/2019 10:19
 Instrument ID: A3UX2 Calib Start Date: 08/03/2018 16:37
 GC Column: ZB-624 ID: 0.53 (mm) Calib End Date: 08/03/2018 19:31
 Lab File ID: X25661.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,4-Dioxane	Lin1		1.043		0.0174	0.0200	-12.9	35.0
1,2-Dichloroethane-d4 (Surr)	Ave	0.3184	0.2996		0.00941	0.0100	-5.9	35.0

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25661.D
 Lims ID: ccvis
 Client ID:
 Sample Type: CCVIS
 Inject. Date: 13-Mar-2019 10:19:30 ALS Bottle#: 2 Worklist Smp#: 3
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: ccvis
 Misc. Info.: 240-0085117-003
 Operator ID: 002808 Instrument ID: A3UX2
 Sublist: chrom-8260SIM_X2*sub2
 Method: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 14-Mar-2019 09:22:23 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0335

First Level Reviewer: macenczaks Date: 13-Mar-2019 10:48:16

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.040	4.040	0.000	100	1954572	10.0	9.41	M
* 7 Fluorobenzene	96	4.611	4.611	0.000	100	6524973	10.0	10.0	
* 9 Dioxane-d8 (IS)	96	5.891	5.891	0.000	92	128678	200.0	200.0	
8 1,4-Dioxane	88	5.969	5.969	0.000	98	13424	20.0	17.4	M

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

vmdioxanew_00165	Amount Added: 3.00	Units: uL	
vm150is_00135	Amount Added: 1.00	Units: uL	Run Reagent
vmDist_H2o_00139	Amount Added: 0.00	Units:	Run Reagent

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25661.D

Injection Date: 13-Mar-2019 10:19:30

Instrument ID: A3UX2

Operator ID: 002808

Lims ID: ccvis

Worklist Smp#: 3

Client ID:

Purge Vol: 15.000 mL

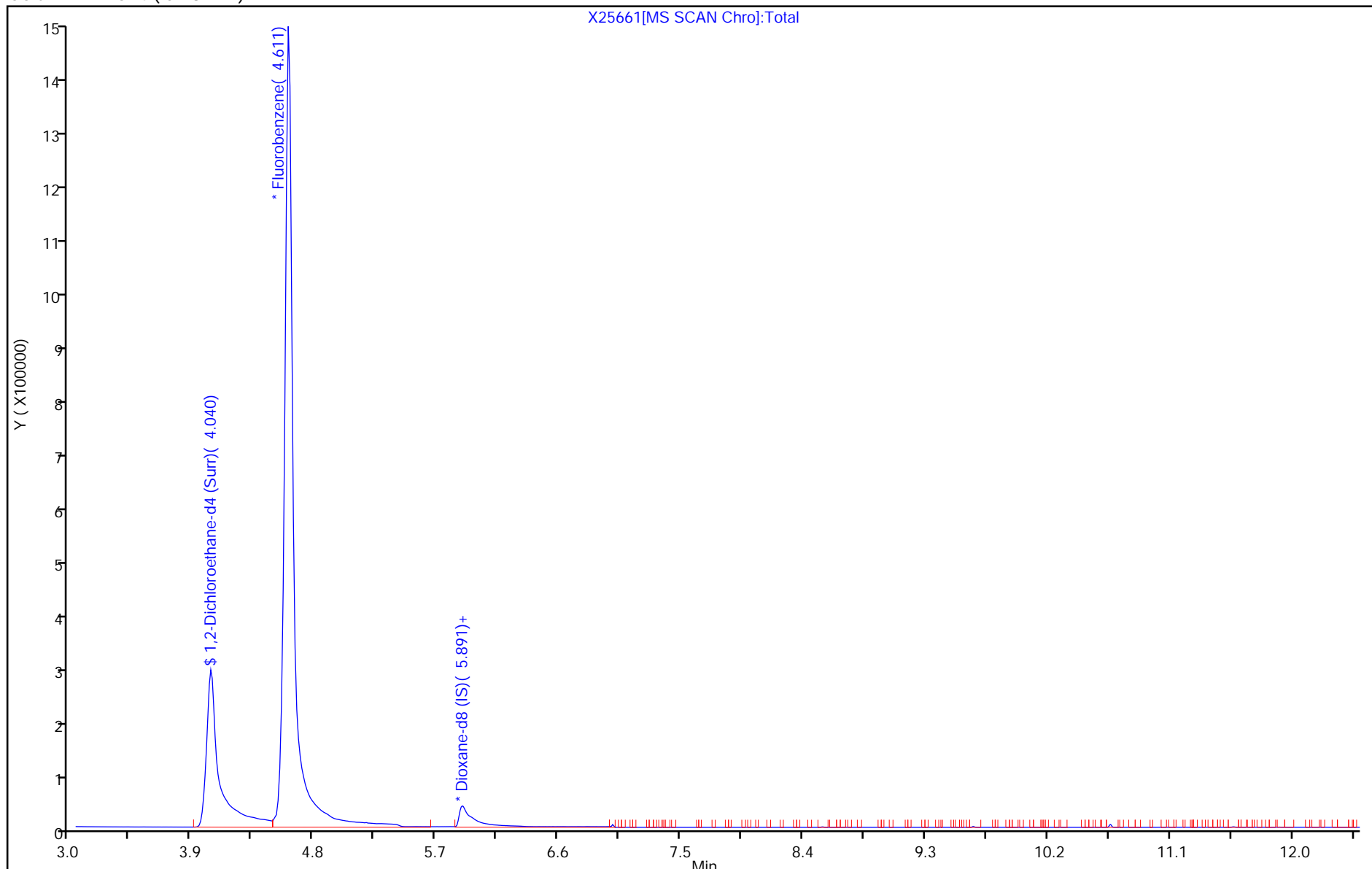
Dil. Factor: 1.0000

ALS Bottle#: 2

Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)



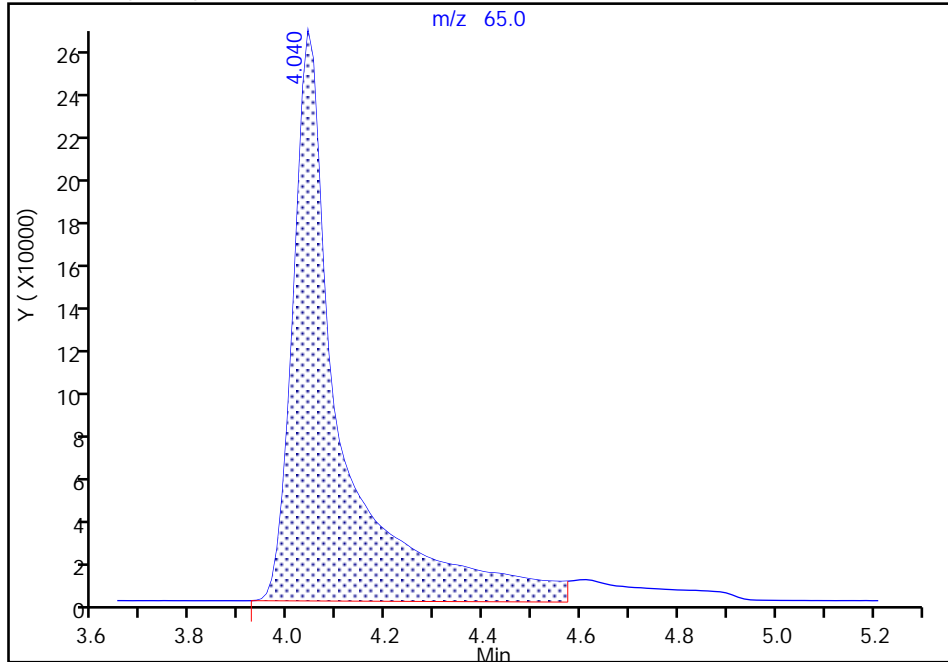
TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25661.D
Injection Date: 13-Mar-2019 10:19:30 Instrument ID: A3UX2
Lims ID: ccvis
Client ID:
Operator ID: 002808 ALS Bottle#: 2 Worklist Smp#: 3
Purge Vol: 15.000 mL Dil. Factor: 1.0000
Method: 8260SIM_X2 Limit Group: MSV SIM ICAL
Column: DB-624 (0.18 mm) Detector: MS SCAN

\$ 5 1,2-Dichloroethane-d4 (Surr), CAS: 17060-07-0
Signal: 1

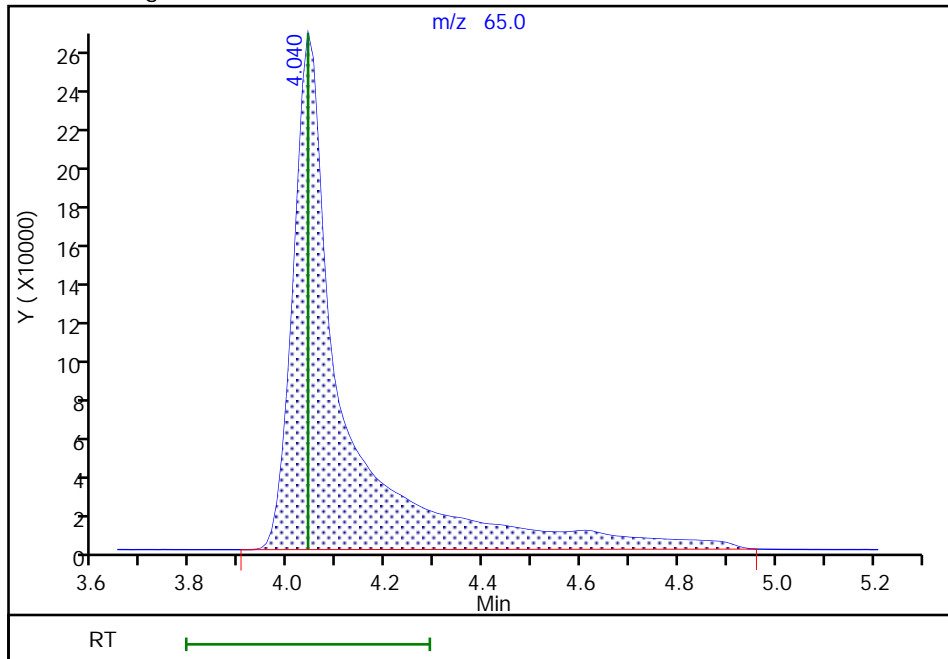
RT: 4.04
Area: 1844196
Amount: 8.875648
Amount Units: ug/l

Processing Integration Results



RT: 4.04
Area: 1954572
Amount: 9.406860
Amount Units: ug/l

Manual Integration Results



Reviewer: macenczaks, 13-Mar-2019 10:47:46
Audit Action: Manually Integrated

Audit Reason: Poor chromatography



TestAmerica Canton

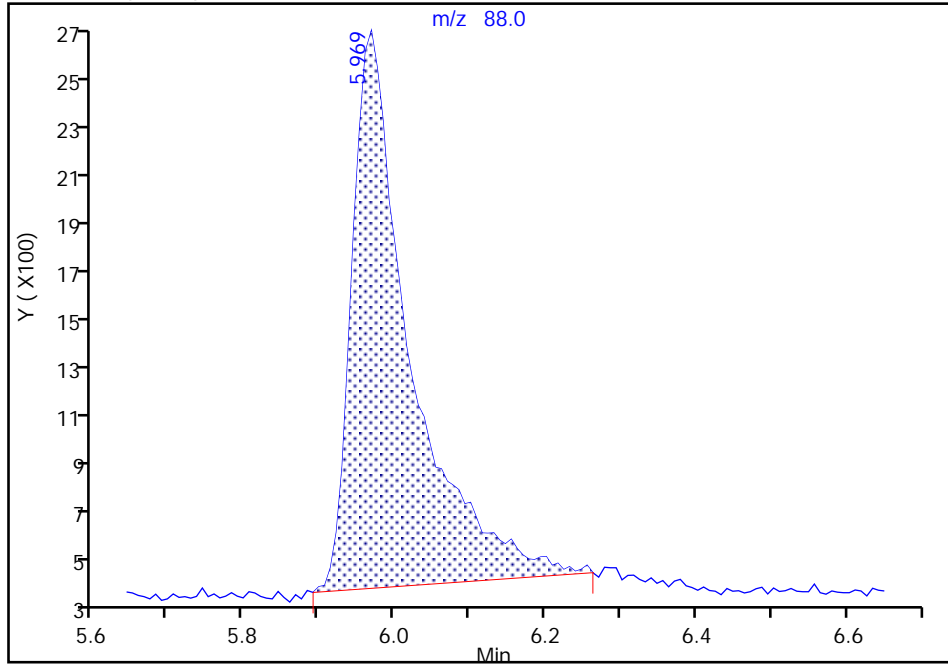
Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25661.D
Injection Date: 13-Mar-2019 10:19:30 Instrument ID: A3UX2
Lims ID: ccvis
Client ID:
Operator ID: 002808 ALS Bottle#: 2 Worklist Smp#: 3
Purge Vol: 15.000 mL Dil. Factor: 1.0000
Method: 8260SIM_X2 Limit Group: MSV SIM ICAL
Column: DB-624 (0.18 mm) Detector: MS SCAN

8 1,4-Dioxane, CAS: 123-91-1

Signal: 1

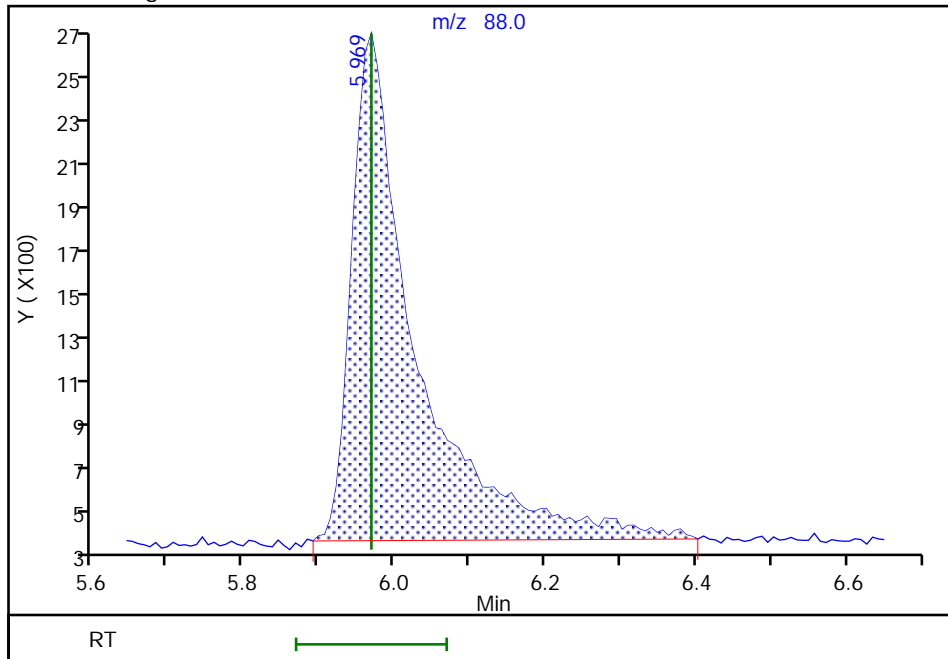
RT: 5.97
Area: 12203
Amount: 15.875539
Amount Units: ug/l

Processing Integration Results



RT: 5.97
Area: 13424
Amount: 17.414560
Amount Units: ug/l

Manual Integration Results



Reviewer: macenczaks, 13-Mar-2019 10:48:10

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 339391

SDG No.: _____

Instrument ID: A3UX2 GC Column: ZB-624 ID: 0.53 (mm) Heated Purge: (Y/N) Y

Calibration Start Date: 08/03/2018 16:37 Calibration End Date: 08/03/2018 19:31 Calibration ID: 46372

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	IC 240-339391/10	X23823.D
Level 2	IC 240-339391/9	X23822.D
Level 3	IC 240-339391/8	X23821.D
Level 4	IC 240-339391/7	X23820.D
Level 5	ICIS 240-339391/6	X23819.D
Level 6	IC 240-339391/5	X23818.D
Level 7	IC 240-339391/4	X23817.D
Level 8	IC 240-339391/3	X23816.D

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
	LVL 6	LVL 7	LVL 8														
1,4-Dioxane	0.8460 1.2319	0.8103 1.2785	0.9740 1.1991	1.0555	1.3209	Lin1	-0.609	1.2331						0.9980		0.9900	
1,2-Dichloroethane-d4 (Surr)	0.3176 0.3197	0.3262 0.3127	0.3228 0.2855	0.3194	0.3436	Ave		0.3184			5.1		35.0				

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

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 339391

SDG No.: _____

Instrument ID: A3UX2 GC Column: ZB-624 ID: 0.53 (mm) Heated Purge: (Y/N) Y

Calibration Start Date: 08/03/2018 16:37 Calibration End Date: 08/03/2018 19:31 Calibration ID: 46372

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	IC 240-339391/10	X23823.D
Level 2	IC 240-339391/9	X23822.D
Level 3	IC 240-339391/8	X23821.D
Level 4	IC 240-339391/7	X23820.D
Level 5	ICIS 240-339391/6	X23819.D
Level 6	IC 240-339391/5	X23818.D
Level 7	IC 240-339391/4	X23817.D
Level 8	IC 240-339391/3	X23816.D

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/L)				
			LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5
			LVL 6	LVL 7	LVL 8			LVL 6	LVL 7	LVL 8		
1,4-Dioxane	14DD 8	Lin1	565 42927	1077 90111	3285 178605	7165	17398	1.00 50.0	2.00 100	5.00 200	10.0	20.0
1,2-Dichloroethane-d4 (Surr)	FB	Ave	95458 5075034	191086 10489340	482900 20302296	957114	2094441	0.500 25.0	1.00 50.0	2.50 100	5.00	10.0

Curve Type Legend:

Ave = Average ISTD
Lin1 = Linear 1/conc ISTD

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23816.D
 Lims ID: IC
 Client ID:
 Sample Type: IC Calib Level: 8
 Inject. Date: 03-Aug-2018 16:37:30 ALS Bottle#: 2 Worklist Smp#: 3
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: ic
 Misc. Info.: 240-0078614-003
 Operator ID: 002808 Instrument ID: A3UX2
 Sublist: chrom-8260SIM_X2*sub2
 Method: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 07-Aug-2018 13:50:33 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK012

First Level Reviewer: macenczaks

Date: 03-Aug-2018 16:54:36

Compound	Sig	RT (min.)	Exp RT (min.)	Diff RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.055	4.056	-0.001	100	20302296	100.0	89.7	
* 7 Fluorobenzene	96	4.622	4.623	-0.001	100	7110946	10.0	10.0	
* 9 Dioxane-d8 (IS)	96	5.898	5.899	-0.001	98	148953	200.0	200.0	
8 1,4-Dioxane	88	5.968	5.969	-0.001	94	178605	200.0	195.0	

Reagents:

vmdioxanew_00141 Amount Added: 30.00 Units: uL
 vm50ss_00330 Amount Added: 30.00 Units: uL
 vm150is_00118 Amount Added: 1.00 Units: uL

TestAmerica Canton

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23816.D

Injection Date: 03-Aug-2018 16:37:30

Instrument ID: A3UX2

Operator ID: 002808

Lims ID: IC

Worklist Smp#: 3

Client ID:

Purge Vol: 15.000 mL

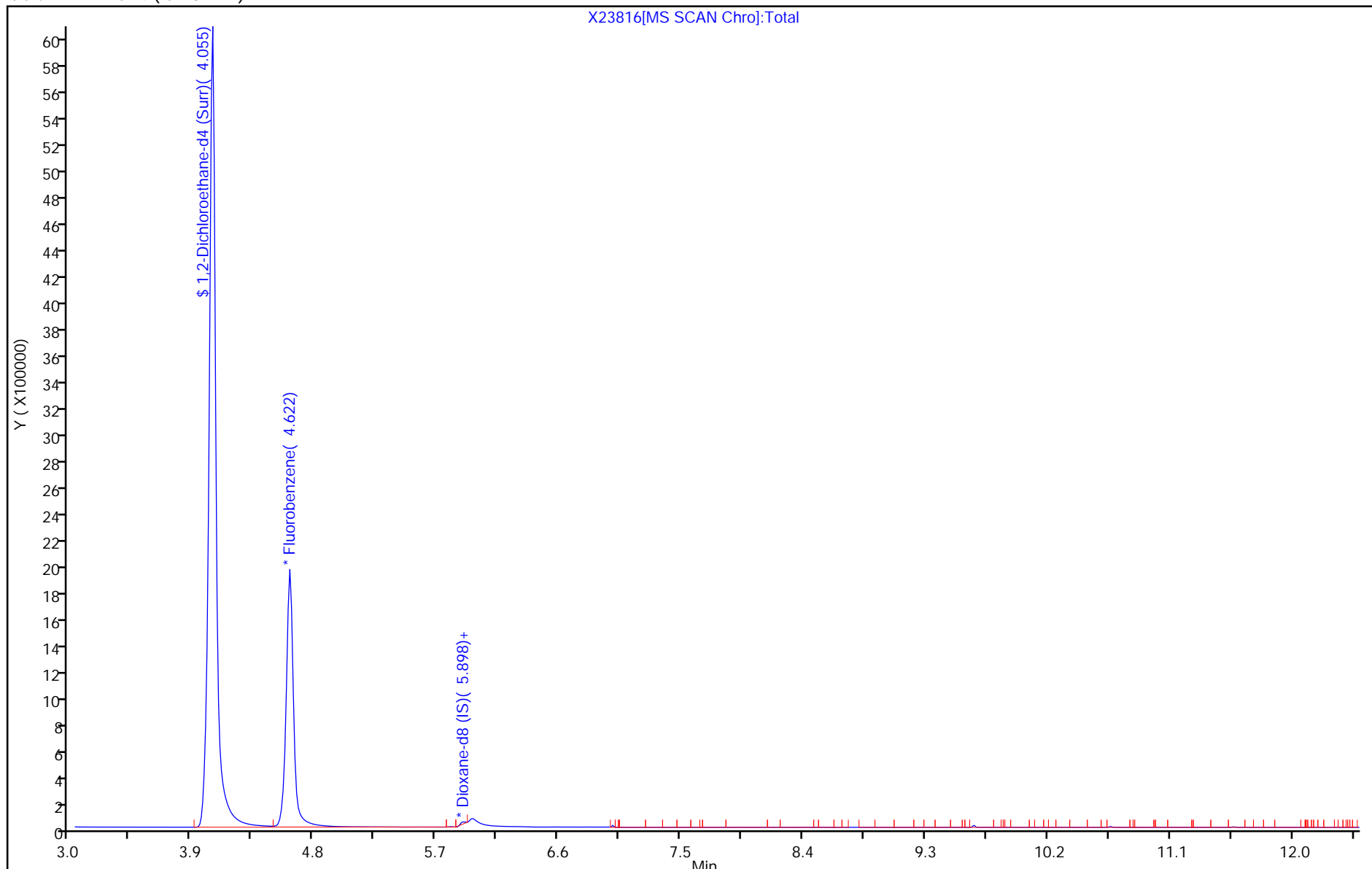
Dil. Factor: 1.0000

ALS Bottle#: 2

Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23817.D
 Lims ID: IC
 Client ID:
 Sample Type: IC Calib Level: 7
 Inject. Date: 03-Aug-2018 17:01:30 ALS Bottle#: 3 Worklist Smp#: 4
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: ic
 Misc. Info.: 240-0078614-004
 Operator ID: 002808 Instrument ID: A3UX2
 Sublist: chrom-8260SIM_X2*sub2
 Method: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 07-Aug-2018 13:50:34 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK012

First Level Reviewer: macenczaks

Date: 06-Aug-2018 08:43:03

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.055	4.056	-0.001	100	10489340	50.0	49.1	
* 7 Fluorobenzene	96	4.623	4.623	0.000	100	6708121	10.0	10.0	
* 9 Dioxane-d8 (IS)	96	5.899	5.899	0.000	99	140967	200.0	200.0	
8 1,4-Dioxane	88	5.968	5.969	-0.001	93	90111	100.0	104.2	

Reagents:

vmdioxanew_00141 Amount Added: 15.00 Units: uL
 vm50ss_00330 Amount Added: 15.00 Units: uL
 vm150is_00118 Amount Added: 1.00 Units: uL

TestAmerica Canton

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23817.D

Injection Date: 03-Aug-2018 17:01:30

Instrument ID: A3UX2

Operator ID: 002808

Lims ID: IC

Worklist Smp#: 4

Client ID:

Purge Vol: 15.000 mL

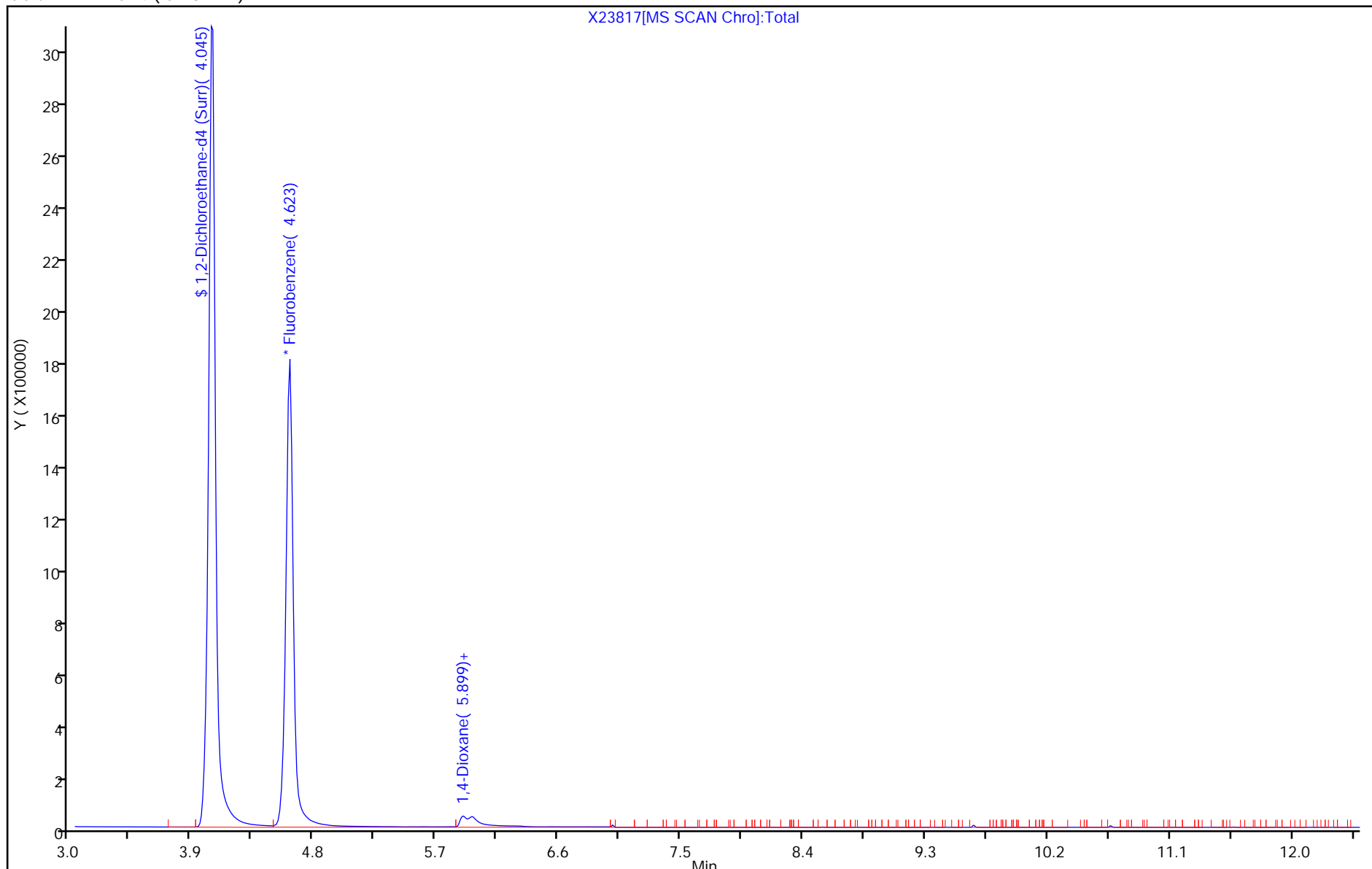
Dil. Factor: 1.0000

ALS Bottle#: 3

Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23818.D
 Lims ID: IC
 Client ID:
 Sample Type: IC Calib Level: 6
 Inject. Date: 03-Aug-2018 17:27:30 ALS Bottle#: 4 Worklist Smp#: 5
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: ic
 Misc. Info.: 240-0078614-005
 Operator ID: 002808 Instrument ID: A3UX2
 Sublist: chrom-8260SIM_X2*sub2
 Method: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 07-Aug-2018 13:50:35 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK012

First Level Reviewer: macenczaks

Date: 06-Aug-2018 08:43:11

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.057	4.056	0.001	100	5075034	25.0	25.1	
* 7 Fluorobenzene	96	4.622	4.623	-0.001	100	6349438	10.0	10.0	
* 9 Dioxane-d8 (IS)	96	5.898	5.899	-0.001	99	139382	200.0	200.0	
8 1,4-Dioxane	88	5.975	5.969	0.006	92	42927	50.0	50.4	

Reagents:

vmdioxanew_00141 Amount Added: 7.50 Units: uL
 vm50ss_00330 Amount Added: 7.50 Units: uL
 vm150is_00118 Amount Added: 1.00 Units: uL

TestAmerica Canton

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23818.D

Injection Date: 03-Aug-2018 17:27:30

Instrument ID: A3UX2

Operator ID: 002808

Lims ID: IC

Worklist Smp#: 5

Client ID:

Purge Vol: 15.000 mL

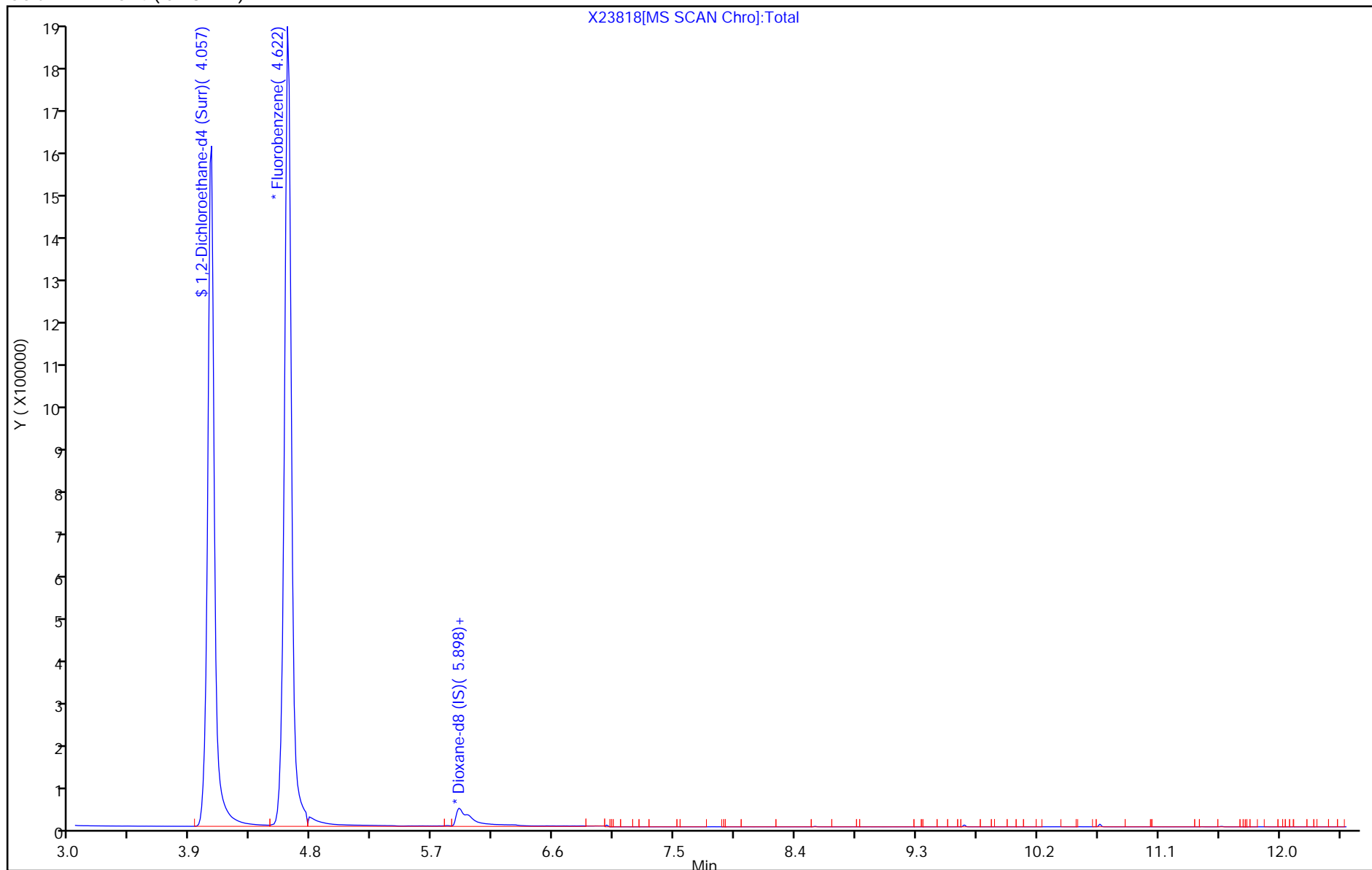
Dil. Factor: 1.0000

ALS Bottle#: 4

Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23819.D
 Lims ID: ICIS
 Client ID:
 Sample Type: ICIS Calib Level: 5
 Inject. Date: 03-Aug-2018 17:52:30 ALS Bottle#: 5 Worklist Smp#: 6
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: icis
 Misc. Info.: 240-0078614-006
 Operator ID: 002808 Instrument ID: A3UX2
 Sublist: chrom-8260SIM_X2*sub2
 Method: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 07-Aug-2018 13:50:35 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK012

First Level Reviewer: macenczaks Date: 06-Aug-2018 08:43:23

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.056	4.056	0.000	100	2094441	10.0	10.8	
* 7 Fluorobenzene	96	4.623	4.623	0.000	100	6095890	10.0	10.0	
* 9 Dioxane-d8 (IS)	96	5.899	5.899	0.000	98	131715	200.0	200.0	
8 1,4-Dioxane	88	5.969	5.969	0.000	93	17398	20.0	21.9	M

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

vmdioxanew_00141 Amount Added: 3.00 Units: uL
 vm50ss_00330 Amount Added: 3.00 Units: uL
 vm150is_00118 Amount Added: 1.00 Units: uL

TestAmerica Canton

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23819.D

Injection Date: 03-Aug-2018 17:52:30

Instrument ID: A3UX2

Operator ID: 002808

Lims ID: ICIS

Worklist Smp#: 6

Client ID:

Purge Vol: 15.000 mL

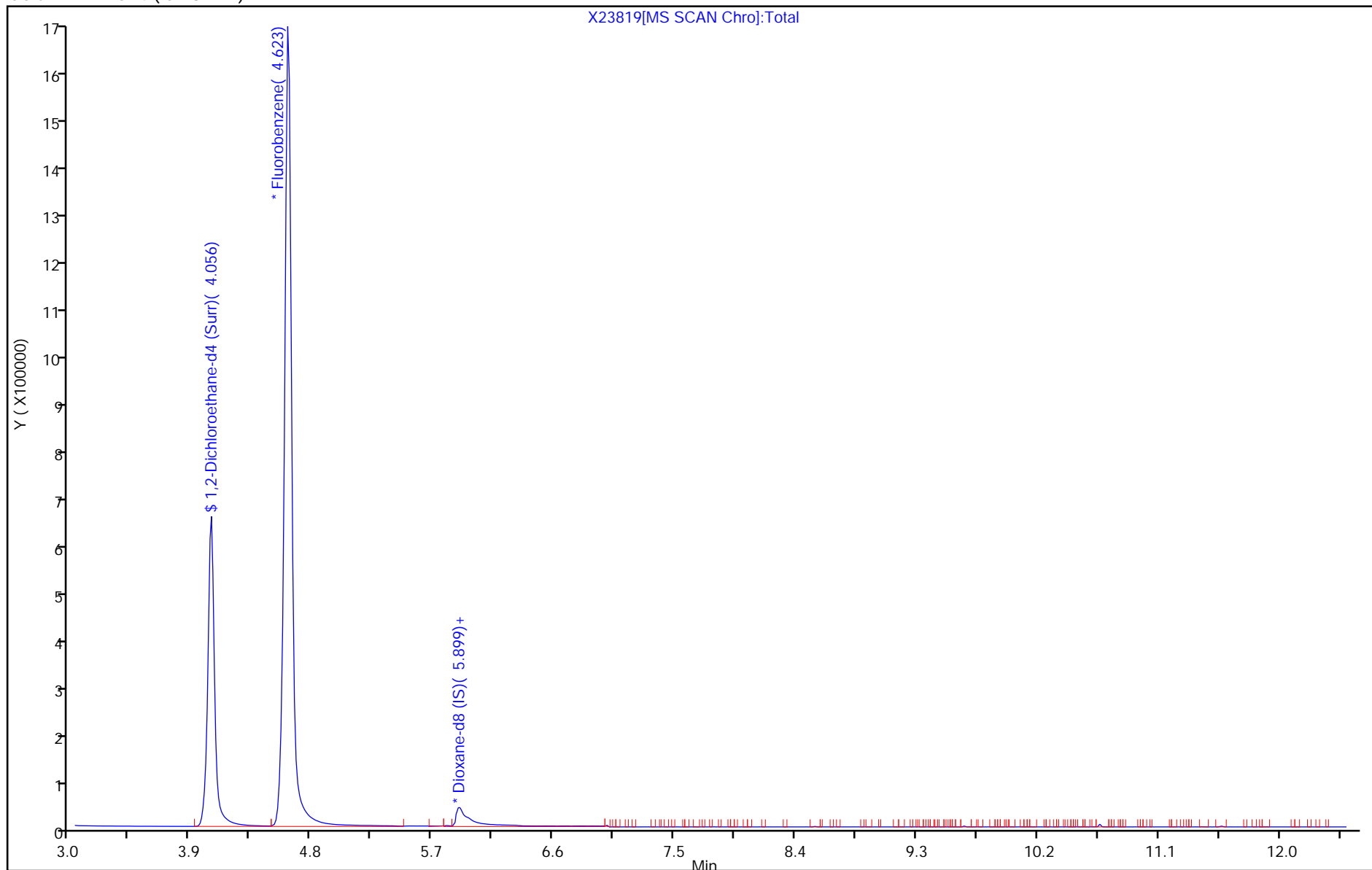
Dil. Factor: 1.0000

ALS Bottle#: 5

Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton

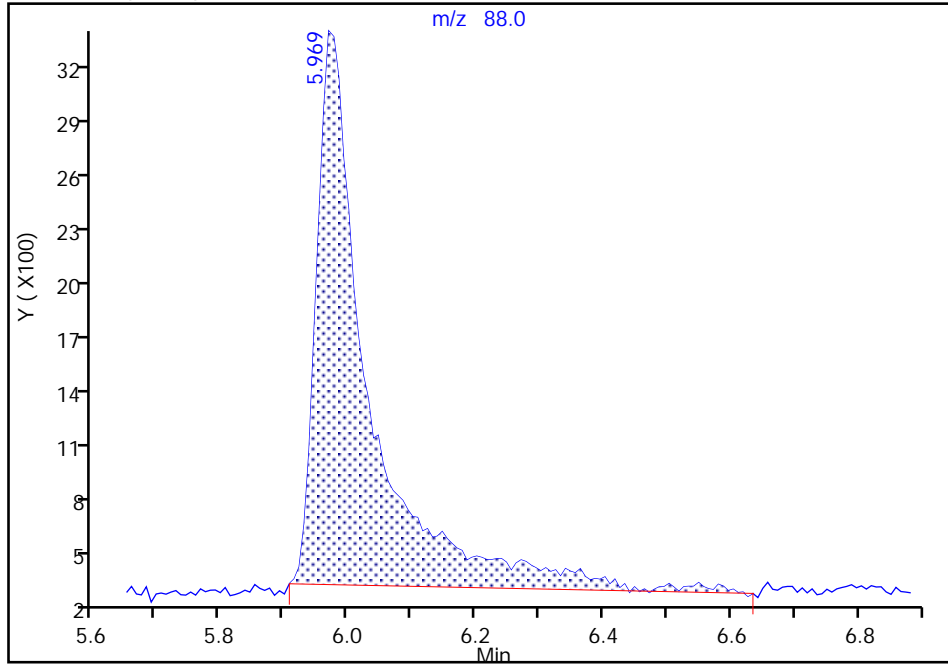
Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23819.D
Injection Date: 03-Aug-2018 17:52:30 Instrument ID: A3UX2
Lims ID: ICIS
Client ID:
Operator ID: 002808 ALS Bottle#: 5 Worklist Smp#: 6
Purge Vol: 15.000 mL Dil. Factor: 1.0000
Method: 8260SIM_X2 Limit Group: MSV SIM ICAL
Column: DB-624 (0.18 mm) Detector MS SCAN

8 1,4-Dioxane, CAS: 123-91-1

Signal: 1

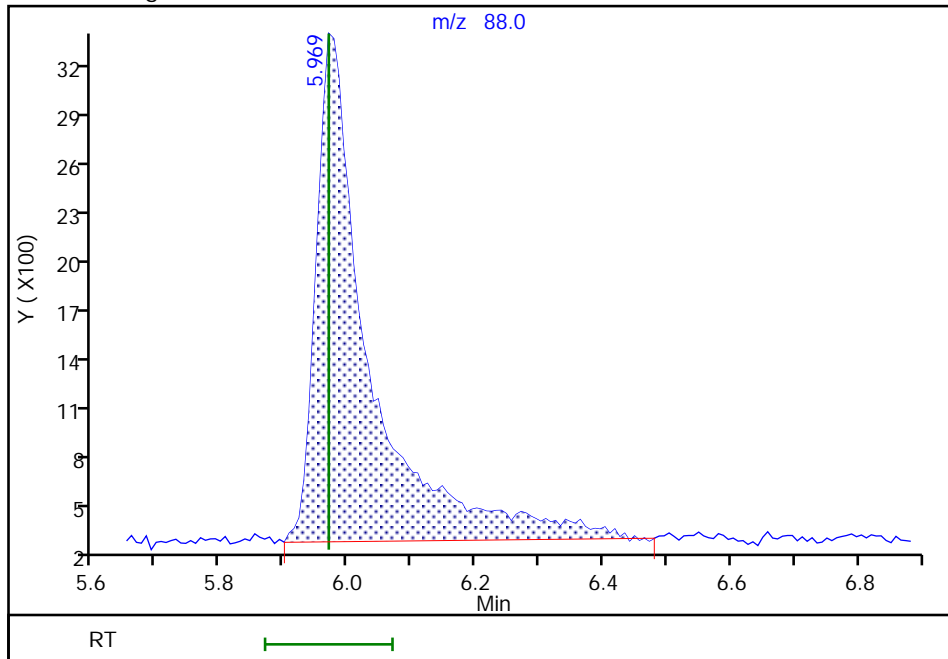
RT: 5.97
Area: 16835
Amount: 24.508842
Amount Units: ug/l

Processing Integration Results



RT: 5.97
Area: 17398
Amount: 21.917989
Amount Units: ug/l

Manual Integration Results



Reviewer: macenczaks, 06-Aug-2018 08:44:10

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23820.D
 Lims ID: IC
 Client ID:
 Sample Type: IC Calib Level: 4
 Inject. Date: 03-Aug-2018 18:16:30 ALS Bottle#: 6 Worklist Smp#: 7
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: ic
 Misc. Info.: 240-0078614-007
 Operator ID: 002808 Instrument ID: A3UX2
 Sublist: chrom-8260SIM_X2*sub2
 Method: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 07-Aug-2018 13:50:37 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK012

First Level Reviewer: macenczaks Date: 06-Aug-2018 08:43:55

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.046	4.056	-0.010	100	957114	5.00	5.01	
* 7 Fluorobenzene	96	4.623	4.623	0.000	100	5993400	10.0	10.0	
* 9 Dioxane-d8 (IS)	96	5.898	5.899	-0.001	99	135762	200.0	200.0	
8 1,4-Dioxane	88	5.968	5.969	-0.001	94	7165	10.0	9.05	M

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

vmdioxanew_00141 Amount Added: 1.50 Units: uL
 vm50ss_00330 Amount Added: 1.50 Units: uL
 vm150is_00118 Amount Added: 1.00 Units: uL

TestAmerica Canton

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23820.D

Injection Date: 03-Aug-2018 18:16:30

Instrument ID: A3UX2

Operator ID: 002808

Lims ID: IC

Worklist Smp#: 7

Client ID:

Purge Vol: 15.000 mL

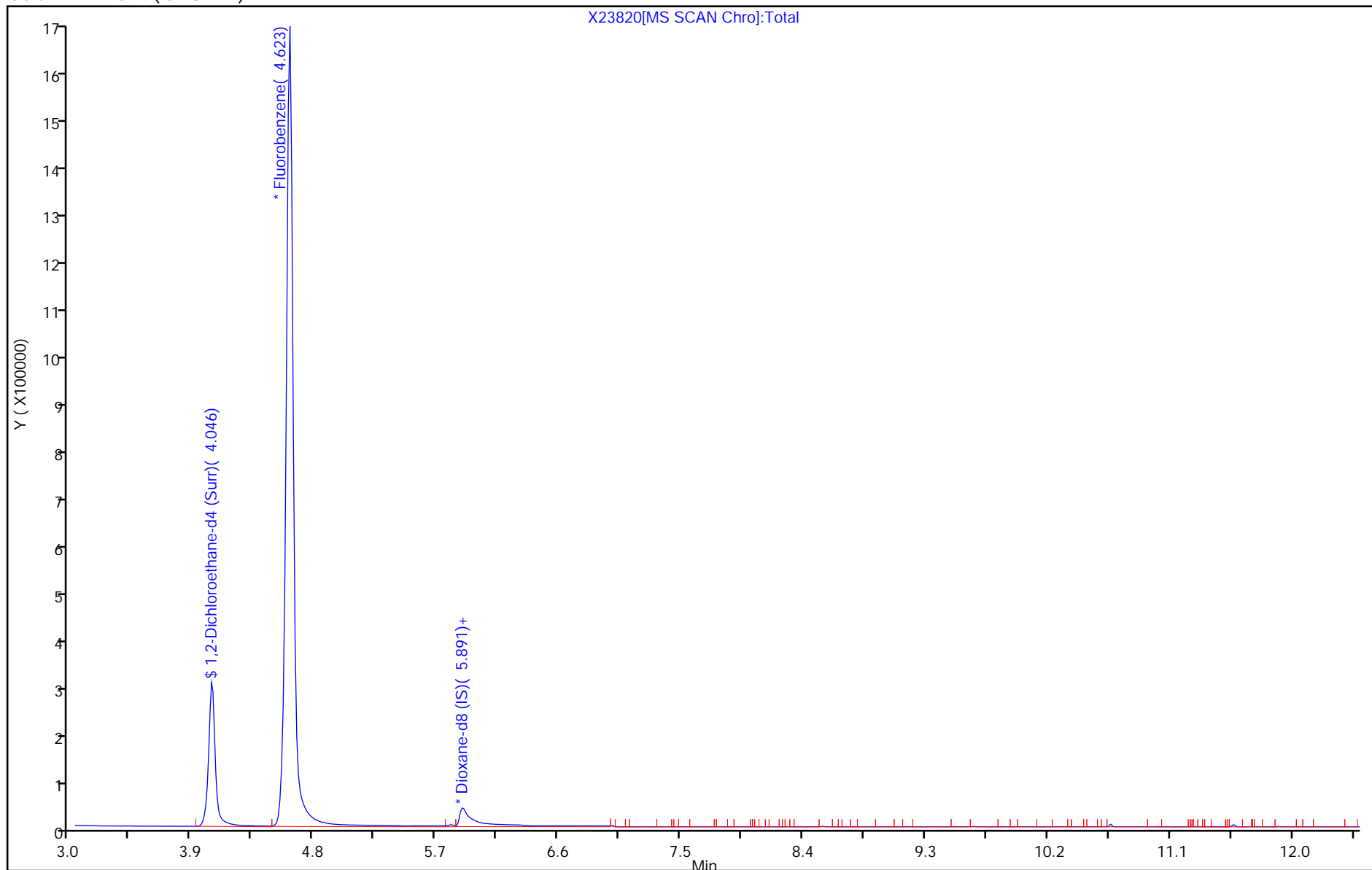
Dil. Factor: 1.0000

ALS Bottle#: 6

Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton

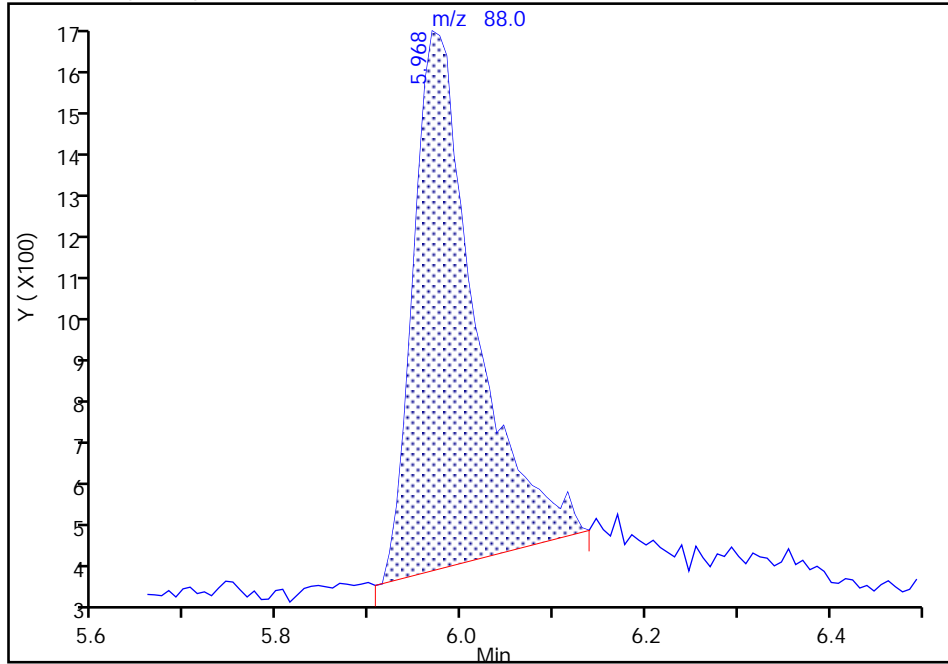
Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23820.D
Injection Date: 03-Aug-2018 18:16:30 Instrument ID: A3UX2
Lims ID: IC
Client ID:
Operator ID: 002808 ALS Bottle#: 6 Worklist Smp#: 7
Purge Vol: 15.000 mL Dil. Factor: 1.0000
Method: 8260SIM_X2 Limit Group: MSV SIM ICAL
Column: DB-624 (0.18 mm) Detector: MS SCAN

8 1,4-Dioxane, CAS: 123-91-1

Signal: 1

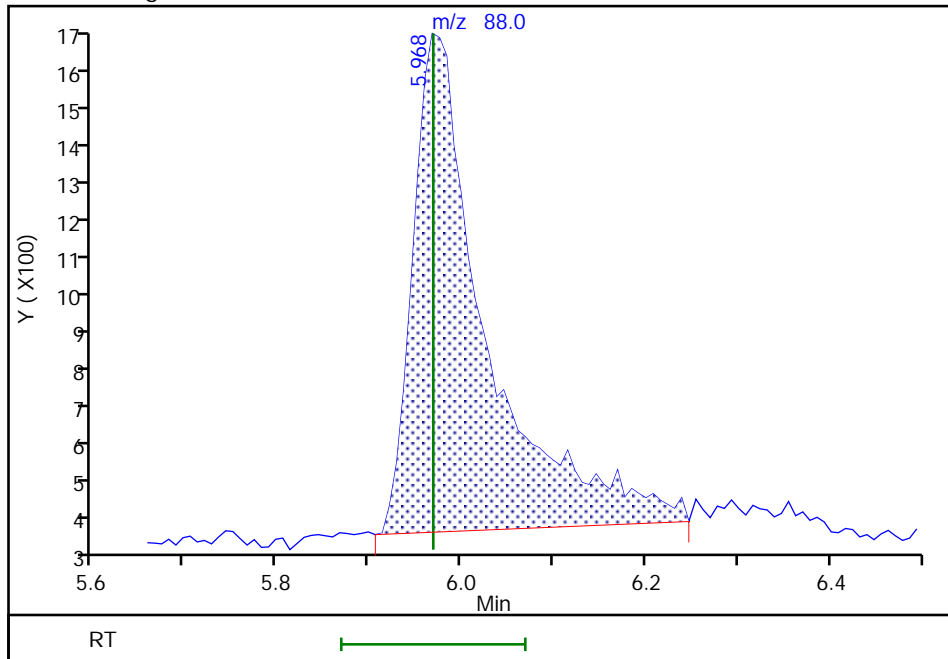
RT: 5.97
Area: 5915
Amount: 8.543057
Amount Units: ug/l

Processing Integration Results



RT: 5.97
Area: 7165
Amount: 9.054101
Amount Units: ug/l

Manual Integration Results



Reviewer: macenczaks, 06-Aug-2018 08:43:42

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23821.D
 Lims ID: IC
 Client ID:
 Sample Type: IC Calib Level: 3
 Inject. Date: 03-Aug-2018 18:41:30 ALS Bottle#: 7 Worklist Smp#: 8
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: ic
 Misc. Info.: 240-0078614-008
 Operator ID: 002808 Instrument ID: A3UX2
 Sublist: chrom-8260SIM_X2*sub2
 Method: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 07-Aug-2018 13:50:38 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK012

First Level Reviewer: macenczaks

Date: 06-Aug-2018 08:45:04

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.057	4.056	0.001	100	482900	2.50	2.53	
* 7 Fluorobenzene	96	4.622	4.623	-0.001	100	5983806	10.0	10.0	
* 9 Dioxane-d8 (IS)	96	5.898	5.899	-0.001	98	134907	200.0	200.0	
8 1,4-Dioxane	88	5.983	5.969	0.014	87	3285	5.00	4.44	M

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

vmdioxanew_00141 Amount Added: 0.75 Units: uL
 vm50ss_00330 Amount Added: 0.75 Units: uL
 vm150is_00118 Amount Added: 1.00 Units: uL

TestAmerica Canton

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23821.D

Injection Date: 03-Aug-2018 18:41:30

Instrument ID: A3UX2

Operator ID: 002808

Lims ID: IC

Worklist Smp#: 8

Client ID:

Purge Vol: 15.000 mL

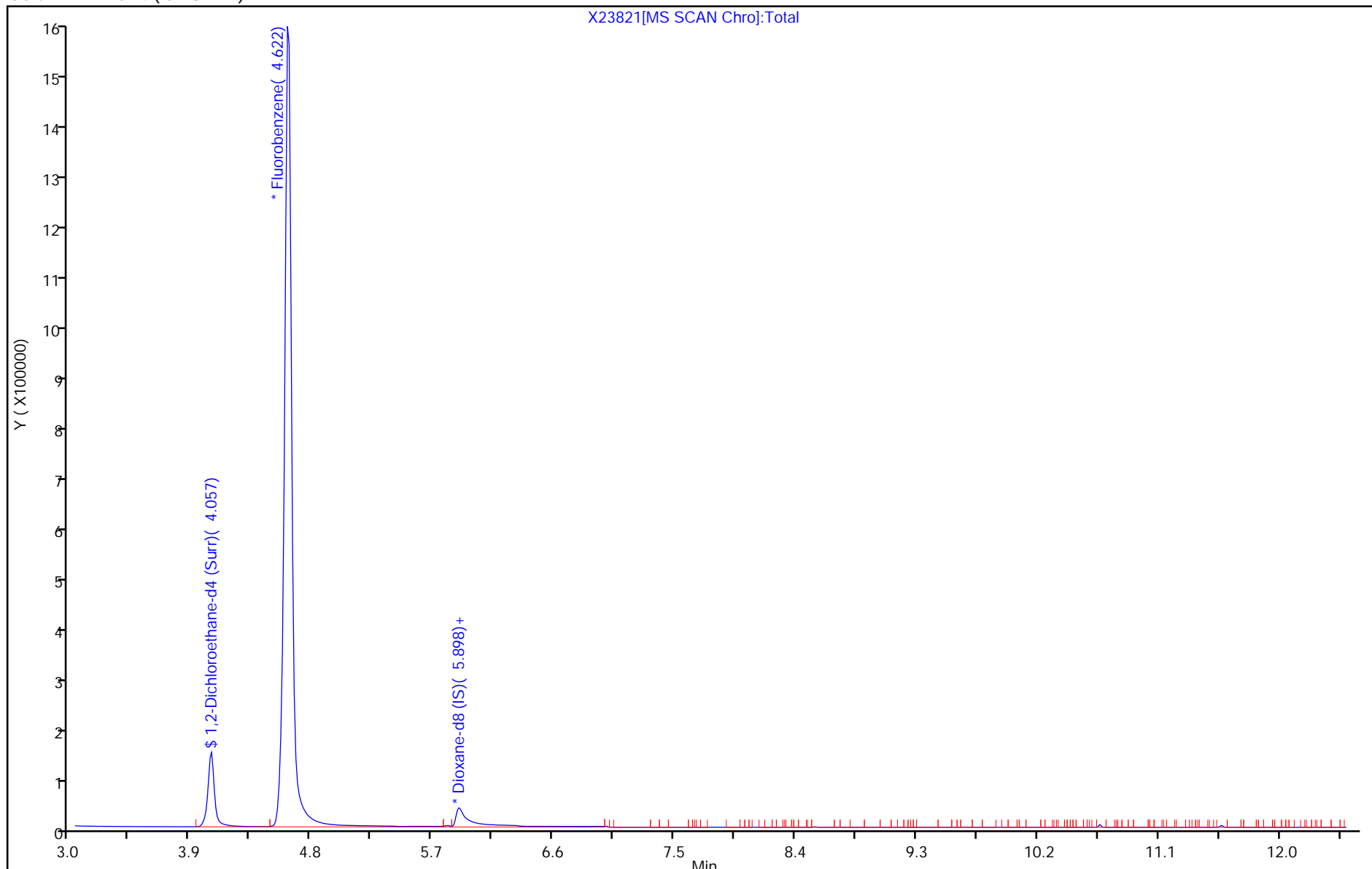
Dil. Factor: 1.0000

ALS Bottle#: 7

Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton

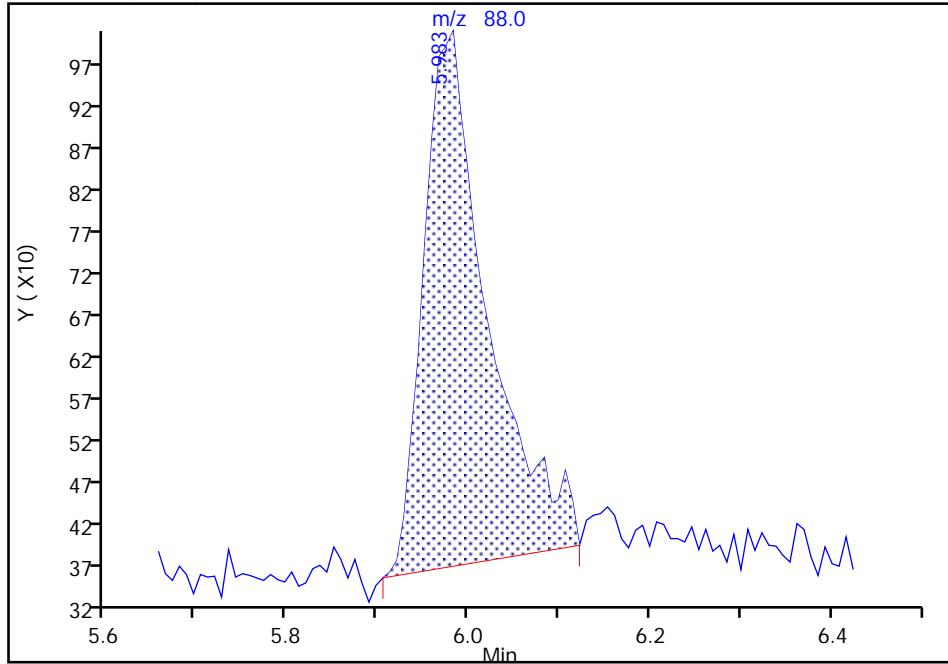
Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23821.D
Injection Date: 03-Aug-2018 18:41:30 Instrument ID: A3UX2
Lims ID: IC
Client ID:
Operator ID: 002808 ALS Bottle#: 7 Worklist Smp#: 8
Purge Vol: 15.000 mL Dil. Factor: 1.0000
Method: 8260SIM_X2 Limit Group: MSV SIM ICAL
Column: DB-624 (0.18 mm) Detector: MS SCAN

8 1,4-Dioxane, CAS: 123-91-1

Signal: 1

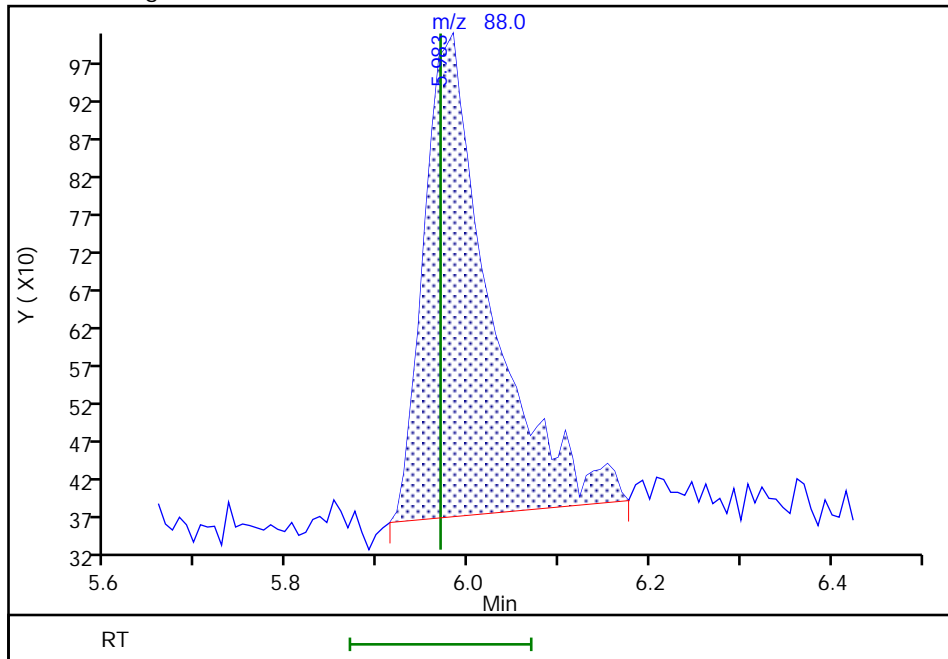
RT: 5.98
Area: 3156
Amount: 4.463015
Amount Units: ug/l

Processing Integration Results



RT: 5.98
Area: 3285
Amount: 4.443580
Amount Units: ug/l

Manual Integration Results



Reviewer: macenczaks, 06-Aug-2018 08:44:55

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23822.D
 Lims ID: IC
 Client ID:
 Sample Type: IC Calib Level: 2
 Inject. Date: 03-Aug-2018 19:06:30 ALS Bottle#: 8 Worklist Smp#: 9
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: ic
 Misc. Info.: 240-0078614-009
 Operator ID: 002808 Instrument ID: A3UX2
 Sublist: chrom-8260SIM_X2*sub2
 Method: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 07-Aug-2018 13:50:39 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK012

First Level Reviewer: macenczaks Date: 06-Aug-2018 08:45:41

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.056	4.056	0.000	100	191086	1.00	1.02	
* 7 Fluorobenzene	96	4.623	4.623	0.000	100	5858078	10.0	10.0	
* 9 Dioxane-d8 (IS)	96	5.898	5.899	-0.001	99	132918	200.0	200.0	
8 1,4-Dioxane	88	5.976	5.969	0.007	87	1077	2.00	1.81	M

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

vmdioxanew_00141 Amount Added: 0.30 Units: uL
 vm50ss_00330 Amount Added: 0.30 Units: uL
 vm150is_00118 Amount Added: 1.00 Units: uL

TestAmerica Canton

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23822.D

Injection Date: 03-Aug-2018 19:06:30

Instrument ID: A3UX2

Operator ID: 002808

Lims ID: IC

Worklist Smp#: 9

Client ID:

Purge Vol: 15.000 mL

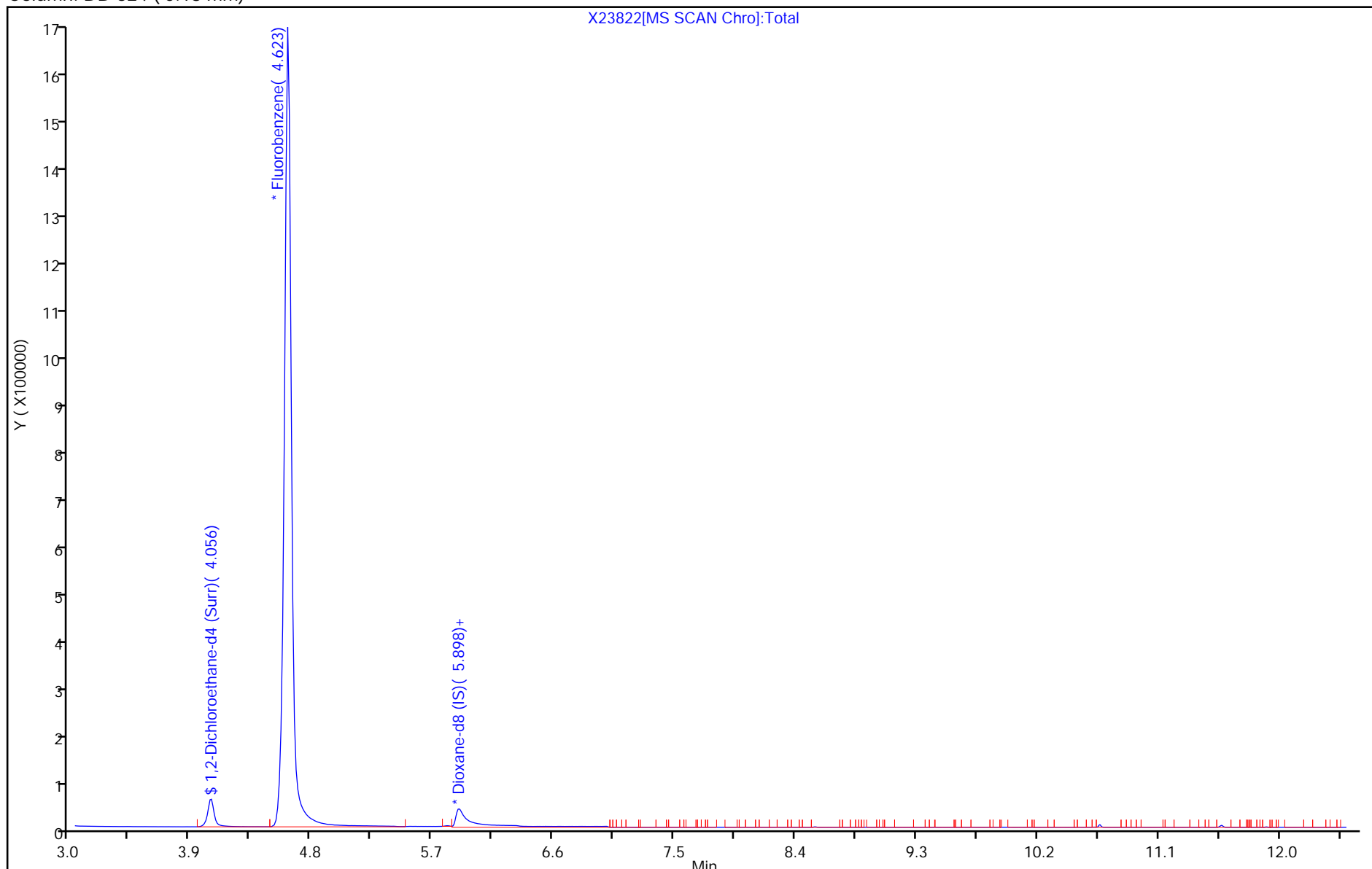
Dil. Factor: 1.0000

ALS Bottle#: 8

Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton

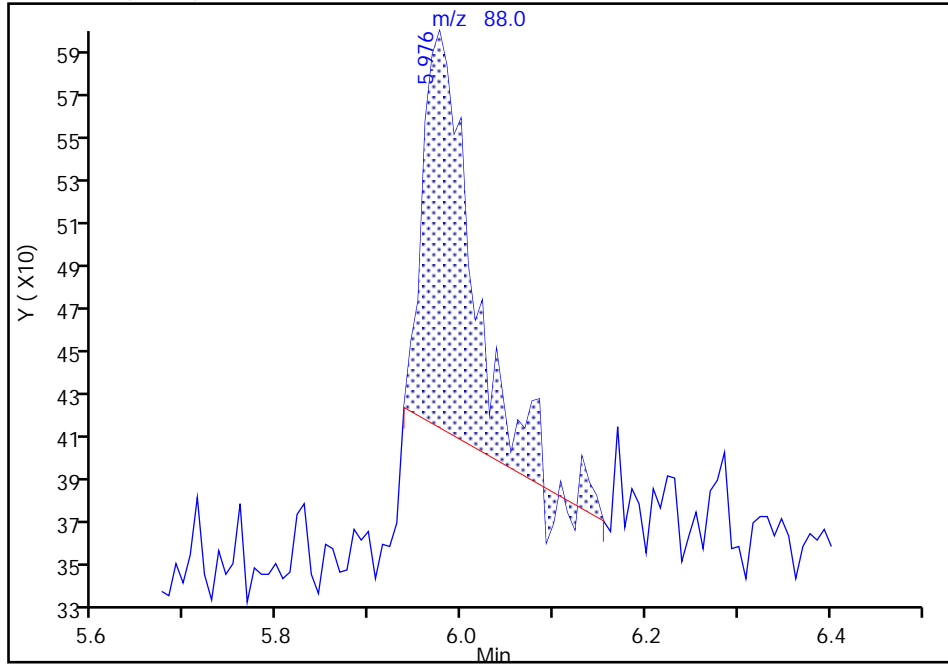
Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23822.D
Injection Date: 03-Aug-2018 19:06:30 Instrument ID: A3UX2
Lims ID: IC
Client ID:
Operator ID: 002808 ALS Bottle#: 8 Worklist Smp#: 9
Purge Vol: 15.000 mL Dil. Factor: 1.0000
Method: 8260SIM_X2 Limit Group: MSV SIM ICAL
Column: DB-624 (0.18 mm) Detector: MS SCAN

8 1,4-Dioxane, CAS: 123-91-1

Signal: 1

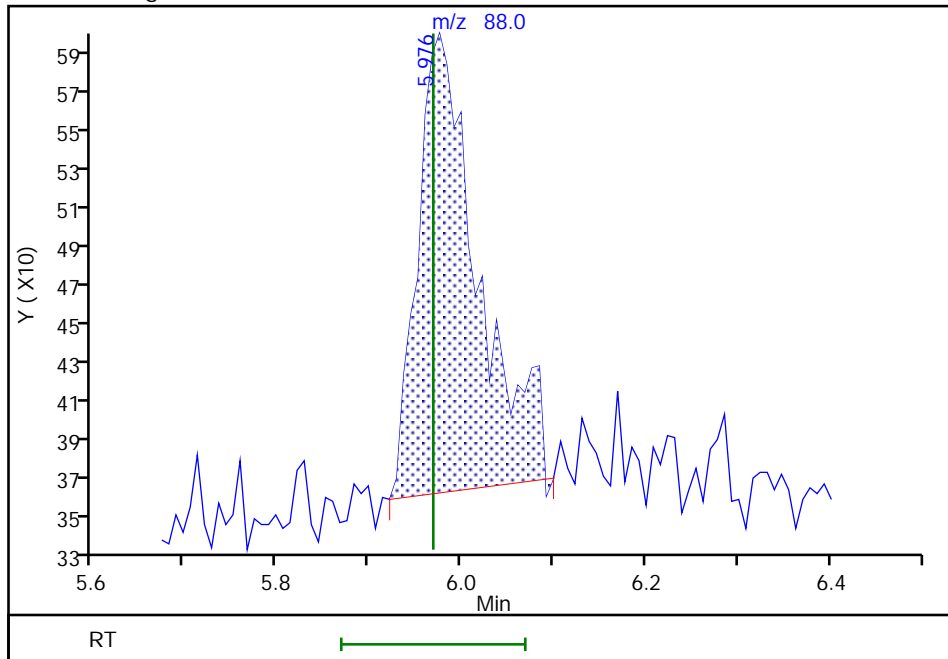
RT: 5.98
Area: 690
Amount: 0.985859
Amount Units: ug/l

Processing Integration Results



RT: 5.98
Area: 1077
Amount: 1.808364
Amount Units: ug/l

Manual Integration Results



Reviewer: macenczaks, 06-Aug-2018 08:45:32

Audit Action: Manually Integrated

Audit Reason: Poor chromatography

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Lims ID: IC
 Client ID:
 Sample Type: IC Calib Level: 1
 Inject. Date: 03-Aug-2018 19:31:30 ALS Bottle#: 9 Worklist Smp#: 10
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: ic
 Misc. Info.: 240-0078614-010
 Operator ID: 002808 Instrument ID: A3UX2
 Sublist: chrom-8260SIM_X2*sub2
 Method: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 07-Aug-2018 13:50:40 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: XAWRK012

First Level Reviewer: macenczaks

Date: 06-Aug-2018 08:47:24

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.055	4.056	-0.001	99	95458	0.5000	0.4987	
* 7 Fluorobenzene	96	4.622	4.623	-0.001	100	6011274	10.0	10.0	
* 9 Dioxane-d8 (IS)	96	5.899	5.899	0.000	99	133567	200.0	200.0	
8 1,4-Dioxane	88	5.976	5.969	0.007	96	565	1.00	1.18	

Reagents:

vmdioxanew_00141 Amount Added: 0.15 Units: uL
 vm50ss_00330 Amount Added: 0.15 Units: uL
 vm150is_00118 Amount Added: 1.00 Units: uL

TestAmerica Canton

Data File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D

Injection Date: 03-Aug-2018 19:31:30

Instrument ID: A3UX2

Operator ID: 002808

Lims ID: IC

Worklist Smp#: 10

Client ID:

Purge Vol: 15.000 mL

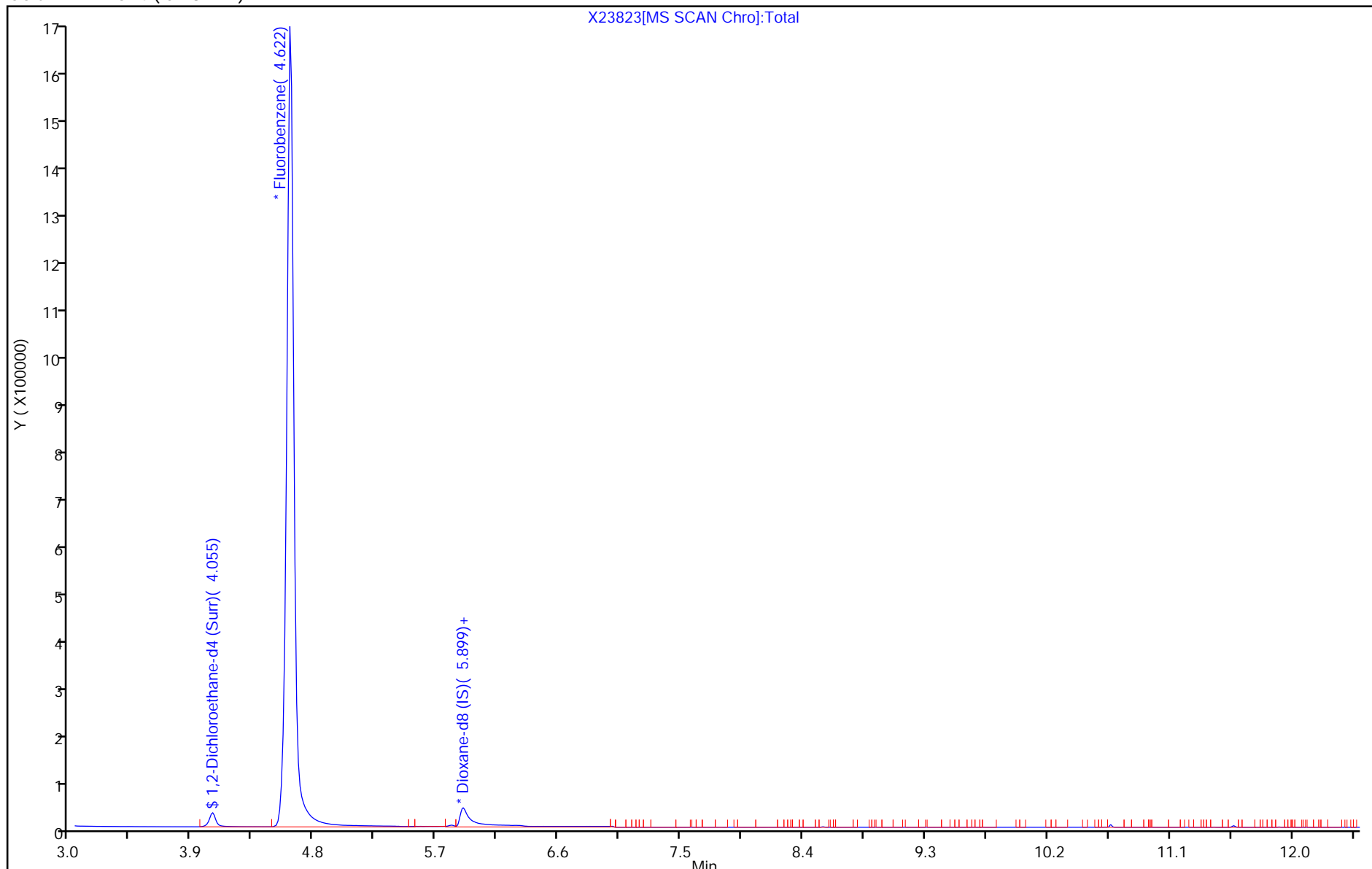
Dil. Factor: 1.0000

ALS Bottle#: 9

Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Client Sample ID: MW-72_030519 Lab Sample ID: 240-109102-1
 Matrix: Water Lab File ID: X25666.D
 Analysis Method: 8260B SIM Date Collected: 03/05/2019 14:20
 Sample wt/vol: 15 (mL) Date Analyzed: 03/13/2019 12:28
 Soil Aliquot Vol: _____ Dilution Factor: 1
 Soil Extract Vol.: _____ GC Column: ZB-624 ID: 0.53 (mm)
 % Moisture: _____ Level: (low/med) Low
 Analysis Batch No.: 371373 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
123-91-1	1,4-Dioxane	0.91	J	2.0	0.86

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	98		63-125

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25666.D
 Lims ID: 240-109102-A-1
 Client ID: MW-72_030519
 Sample Type: Client
 Inject. Date: 13-Mar-2019 12:28:30 ALS Bottle#: 7 Worklist Smp#: 8
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: 240-109102-a-1
 Misc. Info.: 240-0085117-008
 Operator ID: 002808 Instrument ID: A3UX2
 Method: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 14-Mar-2019 09:22:23 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0335

First Level Reviewer: macenczaks

Date: 13-Mar-2019 15:48:15

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/l	Flags
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.043	4.040	0.003	100	2030756	9.80	
* 7 Fluorobenzene	96	4.611	4.611	0.000	100	6504193	10.0	
* 9 Dioxane-d8 (IS)	96	5.890	5.891	-0.001	90	130518	200.0	
8 1,4-Dioxane	88	5.967	5.969	-0.002	50	338	0.9142	M

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

vm150is_00135	Amount Added: 1.00	Units: uL	Run Reagent
vmDist_H2o_00139	Amount Added: 0.00	Units:	Run Reagent

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25666.D

Injection Date: 13-Mar-2019 12:28:30

Instrument ID: A3UX2

Operator ID: 002808

Lims ID: 240-109102-A-1

Lab Sample ID: 240-109102-1

Worklist Smp#: 8

Client ID: MW-72_030519

Purge Vol: 15.000 mL

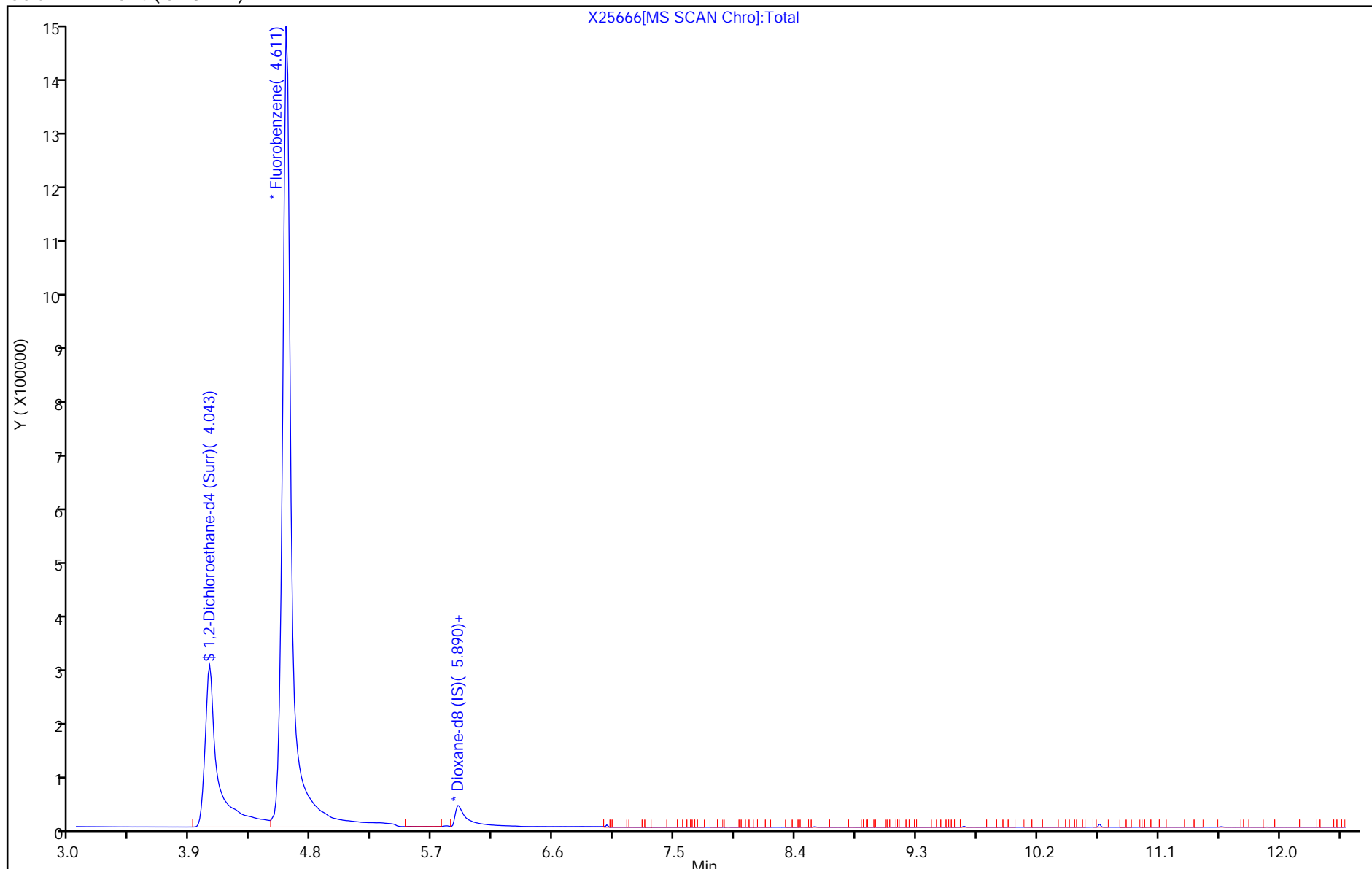
Dil. Factor: 1.0000

ALS Bottle#: 7

Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Recovery Report

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25666.D
 Lims ID: 240-109102-A-1
 Client ID: MW-72_030519
 Sample Type: Client
 Inject. Date: 13-Mar-2019 12:28:30 ALS Bottle#: 7 Worklist Smp#: 8
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: 240-109102-a-1
 Misc. Info.: 240-0085117-008
 Operator ID: 002808 Instrument ID: A3UX2
 Method: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 14-Mar-2019 09:22:23 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0335

First Level Reviewer: macenczaks Date: 13-Mar-2019 15:48:15

Compound	Amount Added	Amount Recovered	% Rec.
\$ 5 1,2-Dichloroethane-d4 (Surr)	10.0	9.80	98.05

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25666.D

Injection Date: 13-Mar-2019 12:28:30

Instrument ID: A3UX2

Lims ID: 240-109102-A-1

Lab Sample ID: 240-109102-1

Client ID: MW-72_030519

Operator ID: 002808

ALS Bottle#: 7 Worklist Smp#: 8

Purge Vol: 15.000 mL

Dil. Factor: 1.0000

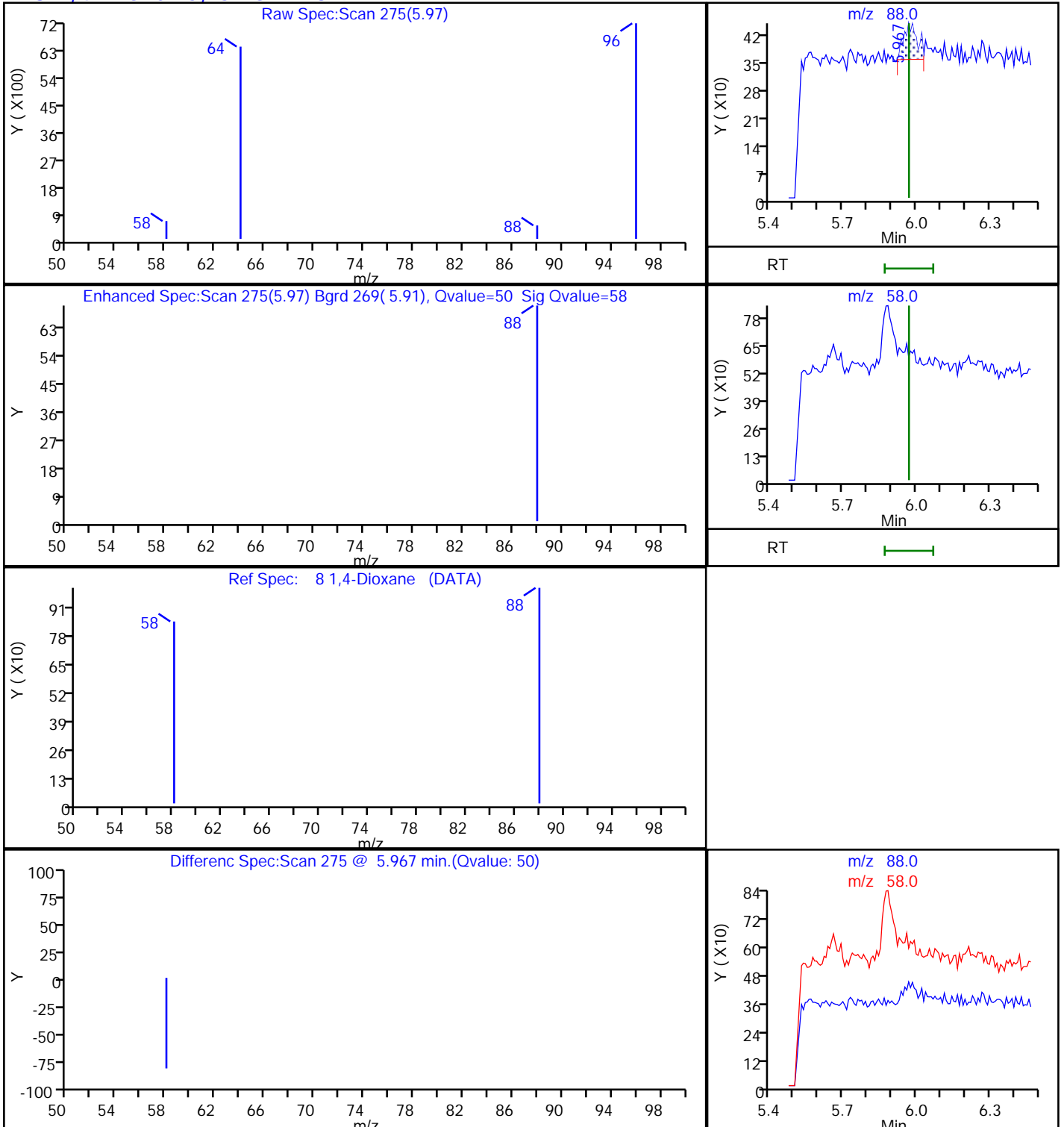
Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)

Detector: MS SCAN

8 1,4-Dioxane, CAS: 123-91-1



TestAmerica Canton

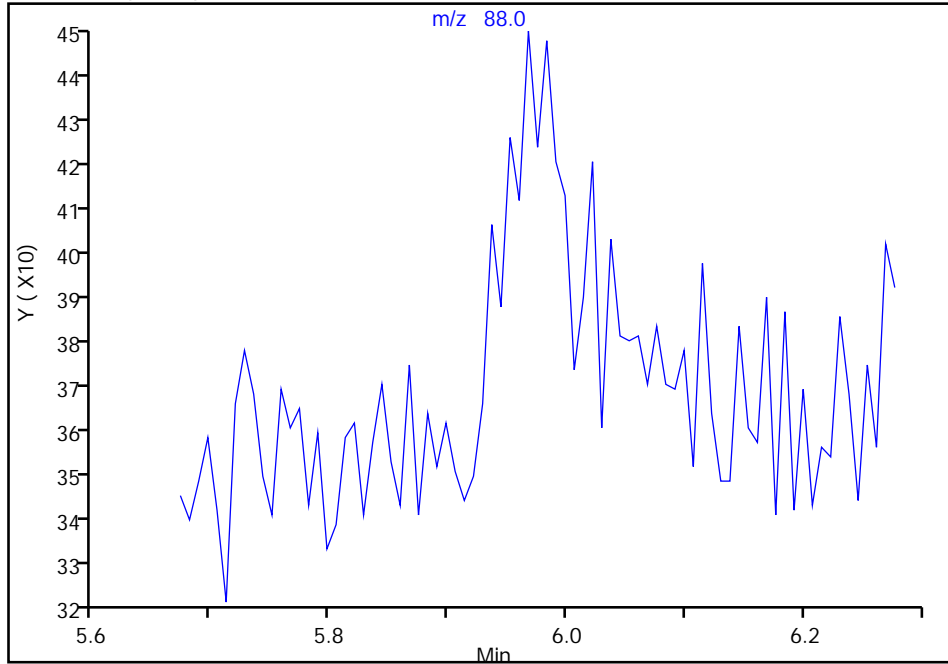
Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25666.D
Injection Date: 13-Mar-2019 12:28:30 Instrument ID: A3UX2
Lims ID: 240-109102-A-1 Lab Sample ID: 240-109102-1
Client ID: MW-72_030519
Operator ID: 002808 ALS Bottle#: 7 Worklist Smp#: 8
Purge Vol: 15.000 mL Dil. Factor: 1.0000
Method: 8260SIM_X2 Limit Group: MSV SIM ICAL
Column: DB-624 (0.18 mm) Detector: MS SCAN

8 1,4-Dioxane, CAS: 123-91-1

Signal: 1

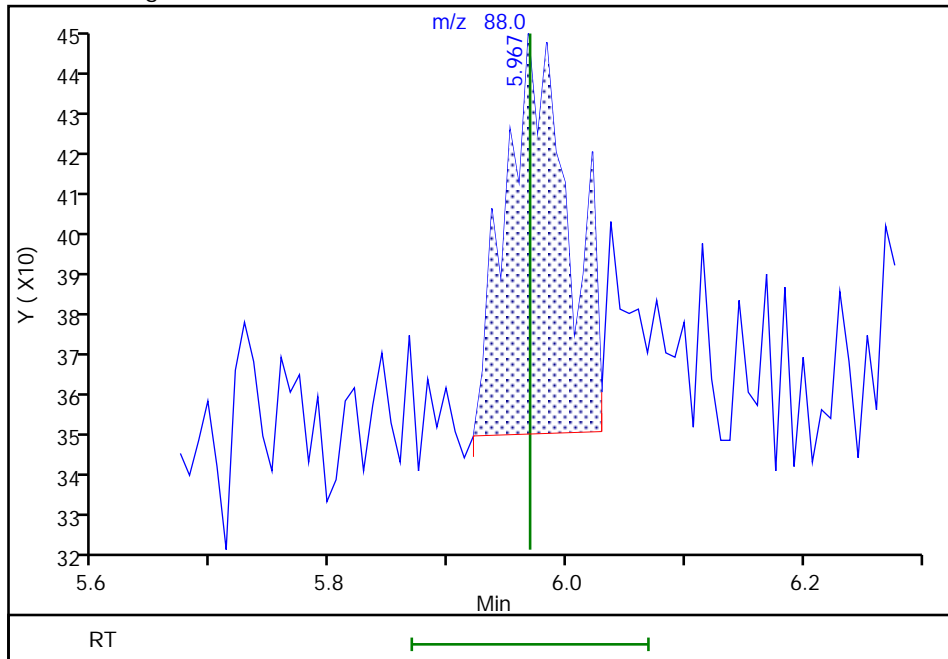
Not Detected
Expected RT: 5.97

Processing Integration Results



RT: 5.97
Area: 338
Amount: 0.914182
Amount Units: ug/l

Manual Integration Results



Reviewer: macenczaks, 13-Mar-2019 15:48:11

Audit Action: Split an Integrated Peak

Audit Reason: Poor chromatography



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Client Sample ID: MW-72S_030519 Lab Sample ID: 240-109102-2
 Matrix: Water Lab File ID: X25667.D
 Analysis Method: 8260B SIM Date Collected: 03/05/2019 15:50
 Sample wt/vol: 15 (mL) Date Analyzed: 03/13/2019 12:53
 Soil Aliquot Vol: _____ Dilution Factor: 1
 Soil Extract Vol.: _____ GC Column: ZB-624 ID: 0.53 (mm)
 % Moisture: _____ Level: (low/med) Low
 Analysis Batch No.: 371373 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
123-91-1	1,4-Dioxane	2.0	U	2.0	0.86

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	99		63-125

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25667.D
 Lims ID: 240-109102-A-2
 Client ID: MW-72S_030519
 Sample Type: Client
 Inject. Date: 13-Mar-2019 12:53:30 ALS Bottle#: 8 Worklist Smp#: 9
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: 240-109102-a-2
 Misc. Info.: 240-0085117-009
 Operator ID: 002808 Instrument ID: A3UX2
 Method: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 14-Mar-2019 09:22:23 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0335

First Level Reviewer: macenczaks

Date: 13-Mar-2019 15:48:56

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/l	Flags
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.042	4.040	0.002	100	1901417	9.88	M
* 7 Fluorobenzene	96	4.610	4.611	-0.001	100	6044946	10.0	
* 9 Dioxane-d8 (IS)	96	5.899	5.891	0.008	92	125263	200.0	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

vm150is_00135	Amount Added: 1.00	Units: uL	Run Reagent
vmDist_H2o_00139	Amount Added: 0.00	Units:	Run Reagent

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25667.D

Injection Date: 13-Mar-2019 12:53:30

Instrument ID: A3UX2

Operator ID: 002808

Lims ID: 240-109102-A-2

Lab Sample ID: 240-109102-2

Worklist Smp#: 9

Client ID: MW-72S_030519

Purge Vol: 15.000 mL

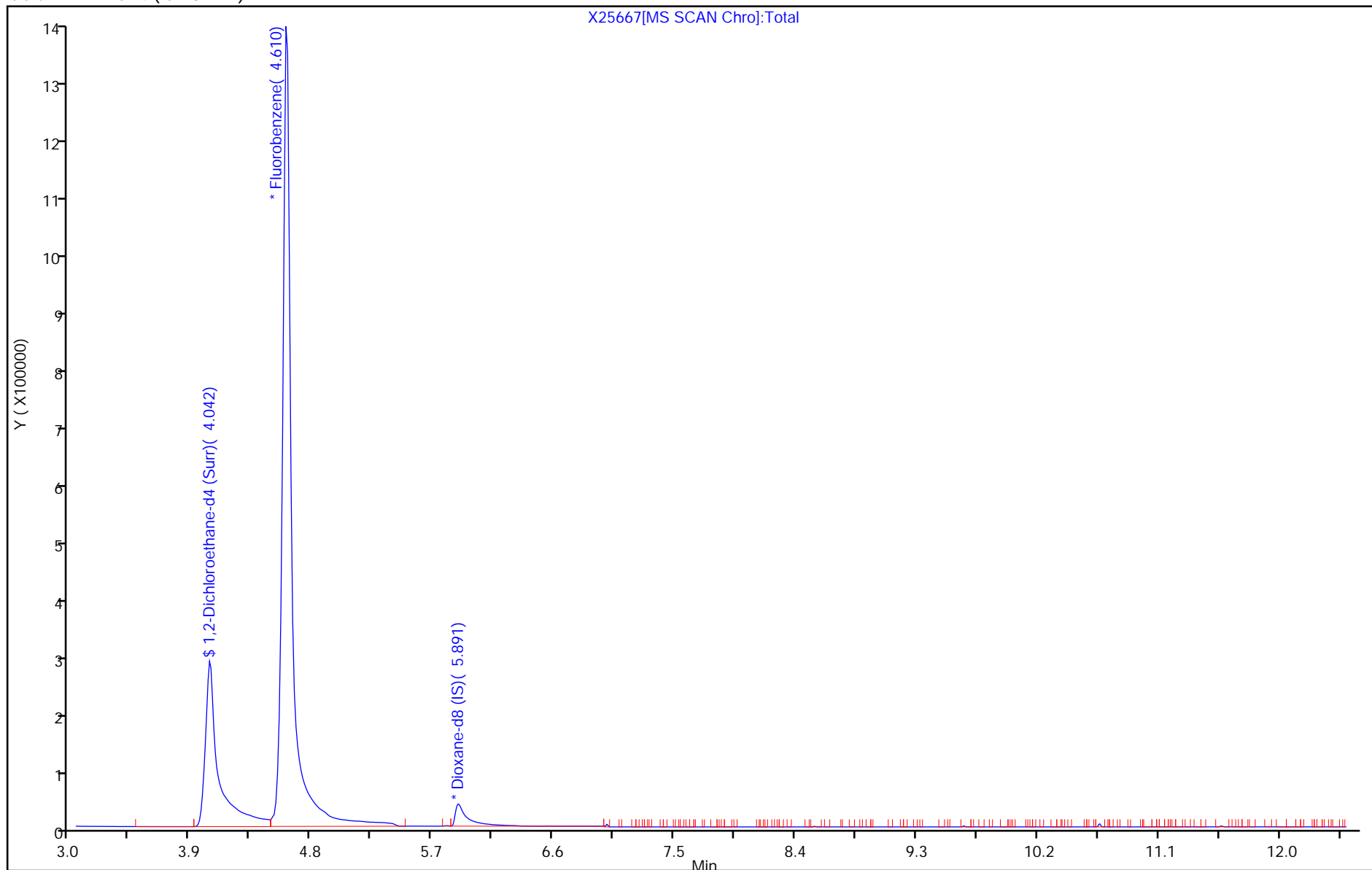
Dil. Factor: 1.0000

ALS Bottle#: 8

Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Recovery Report

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25667.D
 Lims ID: 240-109102-A-2
 Client ID: MW-72S_030519
 Sample Type: Client
 Inject. Date: 13-Mar-2019 12:53:30 ALS Bottle#: 8 Worklist Smp#: 9
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: 240-109102-a-2
 Misc. Info.: 240-0085117-009
 Operator ID: 002808 Instrument ID: A3UX2
 Method: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 14-Mar-2019 09:22:23 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0335

First Level Reviewer: macenczaks Date: 13-Mar-2019 15:48:56

Compound	Amount Added	Amount Recovered	% Rec.
\$ 5 1,2-Dichloroethane-d4 (Surr)	10.0	9.88	98.78

TestAmerica Canton

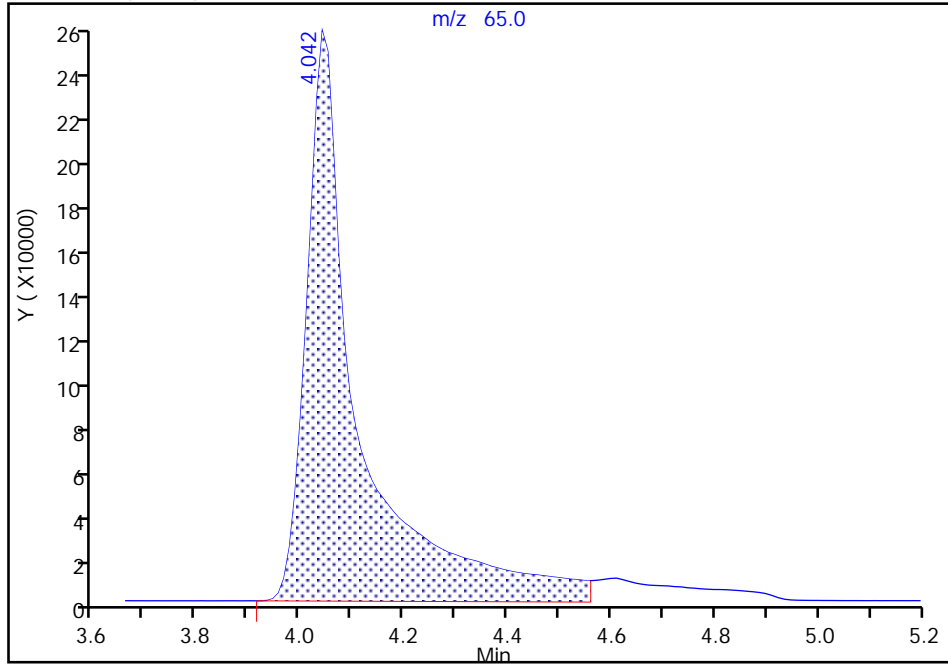
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Injection Date: 13-Mar-2019 12:53:30 Instrument ID: A3UX2
Lims ID: 240-109102-A-2 Lab Sample ID: 240-109102-2
Client ID: MW-72S_030519
Operator ID: 002808 ALS Bottle#: 8 Worklist Smp#: 9
Purge Vol: 15.000 mL Dil. Factor: 1.0000
Method: 8260SIM_X2 Limit Group: MSV SIM ICAL
Column: DB-624 (0.18 mm) Detector MS SCAN

\$ 5 1,2-Dichloroethane-d4 (Surr), CAS: 17060-07-0

Signal: 1

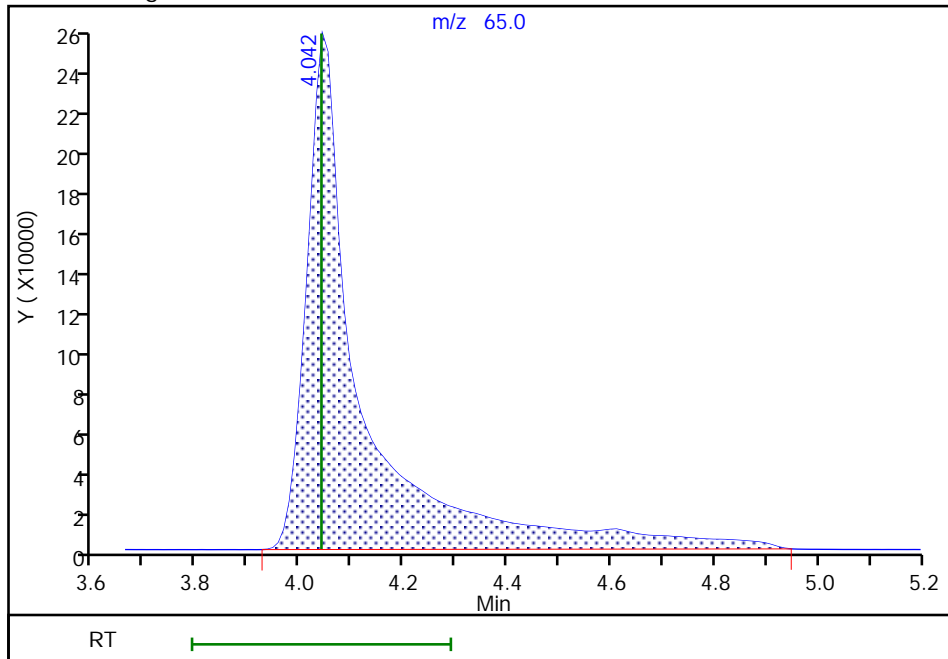
RT: 4.04
Area: 1788805
Amount: 9.292708
Amount Units: ug/l

Processing Integration Results



RT: 4.04
Area: 1901417
Amount: 9.877719
Amount Units: ug/l

Manual Integration Results



Reviewer: macenczaks, 13-Mar-2019 15:48:49

Audit Action: Manually Integrated

Audit Reason: Poor chromatography



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Client Sample ID: MW-83_030519 Lab Sample ID: 240-109102-3
 Matrix: Water Lab File ID: X25668.D
 Analysis Method: 8260B SIM Date Collected: 03/05/2019 12:13
 Sample wt/vol: 15 (mL) Date Analyzed: 03/13/2019 13:19
 Soil Aliquot Vol: _____ Dilution Factor: 1
 Soil Extract Vol.: _____ GC Column: ZB-624 ID: 0.53 (mm)
 % Moisture: _____ Level: (low/med) Low
 Analysis Batch No.: 371373 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
123-91-1	1,4-Dioxane	2.0	U	2.0	0.86

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	101		63-125

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25668.D
 Lims ID: 240-109102-A-3
 Client ID: MW-83_030519
 Sample Type: Client
 Inject. Date: 13-Mar-2019 13:19:30 ALS Bottle#: 9 Worklist Smp#: 10
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: 240-109102-a-3
 Misc. Info.: 240-0085117-010
 Operator ID: 002808 Instrument ID: A3UX2
 Method: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 14-Mar-2019 09:22:23 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0335

First Level Reviewer: macenczaks

Date: 13-Mar-2019 15:49:17

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/l	Flags
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.039	4.040	-0.001	100	1907926	10.1	
* 7 Fluorobenzene	96	4.610	4.611	-0.001	100	5951939	10.0	
* 9 Dioxane-d8 (IS)	96	5.898	5.891	0.007	92	124614	200.0	

Reagents:

vm150is_00135 Amount Added: 1.00 Units: uL Run Reagent
 vmDist_H2o_00139 Amount Added: 0.00 Units: Run Reagent

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25668.D

Injection Date: 13-Mar-2019 13:19:30

Instrument ID: A3UX2

Operator ID: 002808

Lims ID: 240-109102-A-3

Lab Sample ID: 240-109102-3

Worklist Smp#: 10

Client ID: MW-83_030519

Purge Vol: 15.000 mL

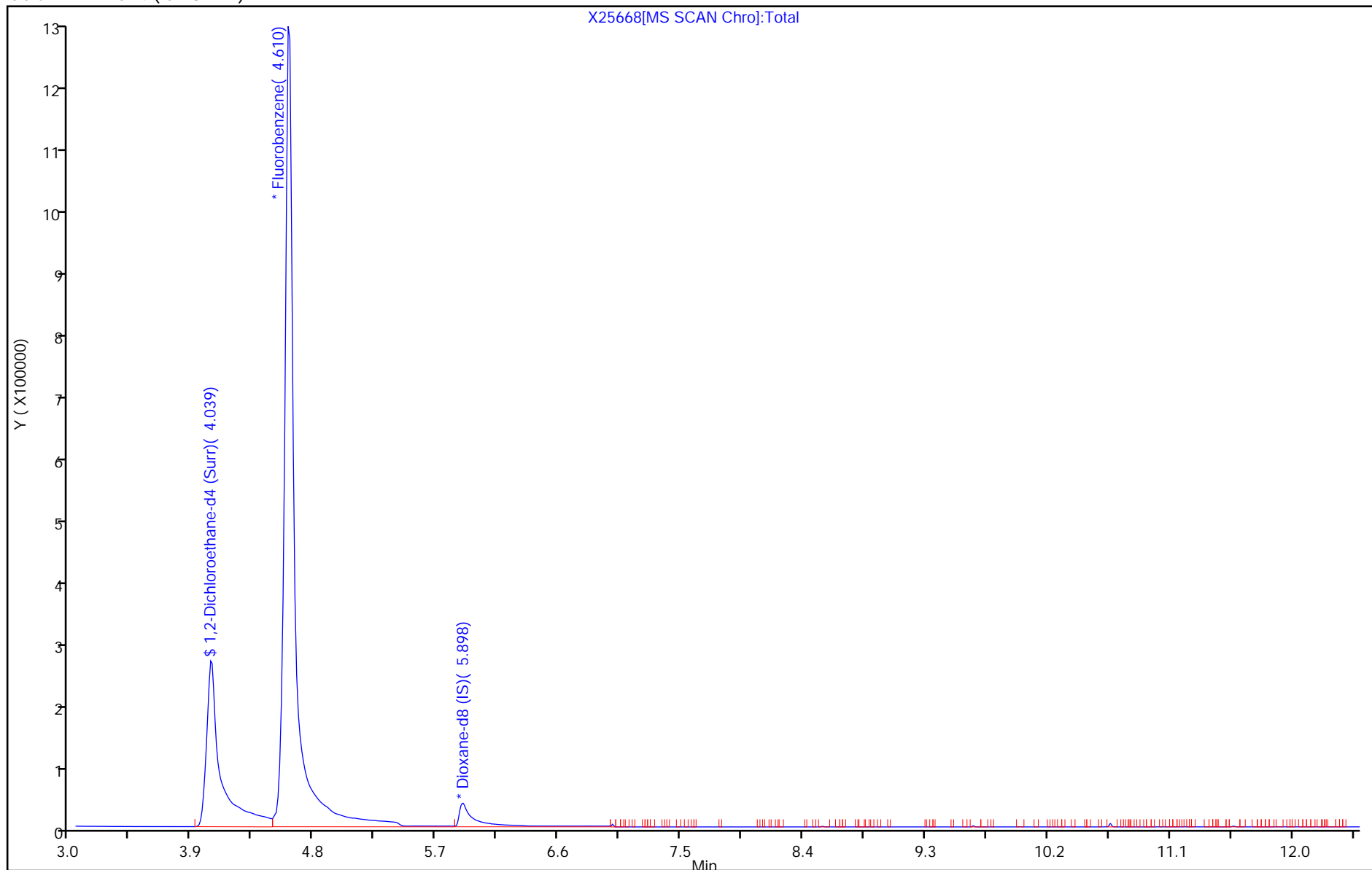
Dil. Factor: 1.0000

ALS Bottle#: 9

Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Recovery Report

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25668.D
 Lims ID: 240-109102-A-3
 Client ID: MW-83_030519
 Sample Type: Client
 Inject. Date: 13-Mar-2019 13:19:30 ALS Bottle#: 9 Worklist Smp#: 10
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: 240-109102-a-3
 Misc. Info.: 240-0085117-010
 Operator ID: 002808 Instrument ID: A3UX2
 Method: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 14-Mar-2019 09:22:23 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0335

First Level Reviewer: macenczaks Date: 13-Mar-2019 15:49:17

Compound	Amount Added	Amount Recovered	% Rec.
\$ 5 1,2-Dichloroethane-d4 (Surr)	10.0	10.1	100.66

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Client Sample ID: MW-83S_030519 Lab Sample ID: 240-109102-4
 Matrix: Water Lab File ID: X25669.D
 Analysis Method: 8260B SIM Date Collected: 03/05/2019 14:07
 Sample wt/vol: 15 (mL) Date Analyzed: 03/13/2019 13:45
 Soil Aliquot Vol: _____ Dilution Factor: 1
 Soil Extract Vol.: _____ GC Column: ZB-624 ID: 0.53 (mm)
 % Moisture: _____ Level: (low/med) Low
 Analysis Batch No.: 371373 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
123-91-1	1,4-Dioxane	2.0	U	2.0	0.86

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	98		63-125

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25669.D
 Lims ID: 240-109102-A-4
 Client ID: MW-83S_030519
 Sample Type: Client
 Inject. Date: 13-Mar-2019 13:45:30 ALS Bottle#: 10 Worklist Smp#: 11
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: 240-109102-a-4
 Misc. Info.: 240-0085117-011
 Operator ID: 002808 Instrument ID: A3UX2
 Method: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 14-Mar-2019 09:22:23 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0335

First Level Reviewer: macenczaks

Date: 13-Mar-2019 15:49:53

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/l	Flags
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.040	4.040	0.000	100	1938308	9.80	M
* 7 Fluorobenzene	96	4.611	4.611	0.000	100	6208534	10.0	
* 9 Dioxane-d8 (IS)	96	5.892	5.891	0.001	90	127666	200.0	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

vm150is_00135	Amount Added: 1.00	Units: uL	Run Reagent
vmDist_H2o_00139	Amount Added: 0.00	Units:	Run Reagent

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25669.D

Injection Date: 13-Mar-2019 13:45:30

Instrument ID: A3UX2

Operator ID: 002808

Lims ID: 240-109102-A-4

Lab Sample ID: 240-109102-4

Worklist Smp#: 11

Client ID: MW-83S_030519

Purge Vol: 15.000 mL

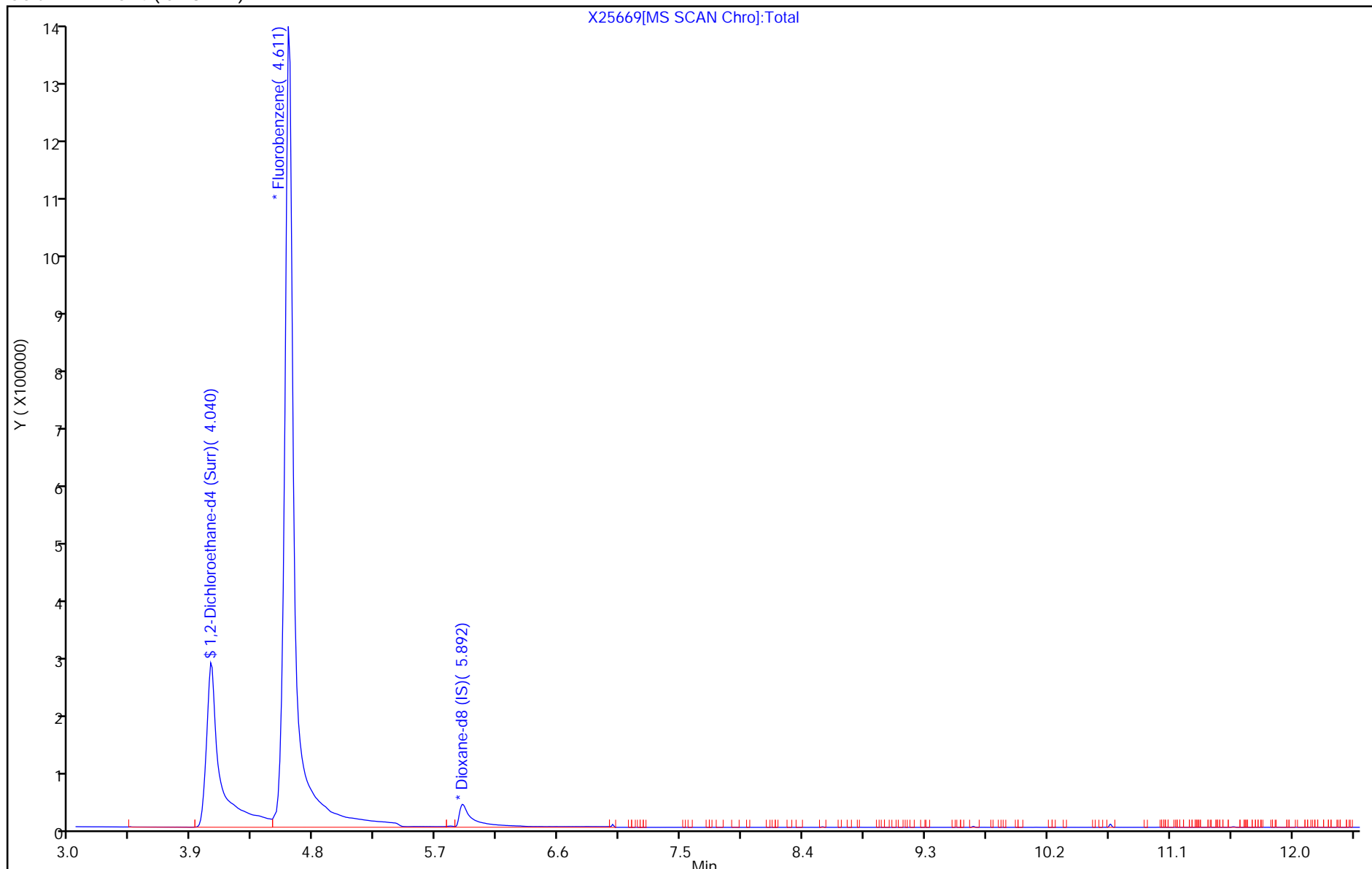
Dil. Factor: 1.0000

ALS Bottle#: 10

Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Recovery Report

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25669.D
 Lims ID: 240-109102-A-4
 Client ID: MW-83S_030519
 Sample Type: Client
 Inject. Date: 13-Mar-2019 13:45:30 ALS Bottle#: 10 Worklist Smp#: 11
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: 240-109102-a-4
 Misc. Info.: 240-0085117-011
 Operator ID: 002808 Instrument ID: A3UX2
 Method: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 14-Mar-2019 09:22:23 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0335

First Level Reviewer: macenczaks

Date: 13-Mar-2019 15:49:53

Compound	Amount Added	Amount Recovered	% Rec.
\$ 5 1,2-Dichloroethane-d4 (Surr)	10.0	9.80	98.04

TestAmerica Canton

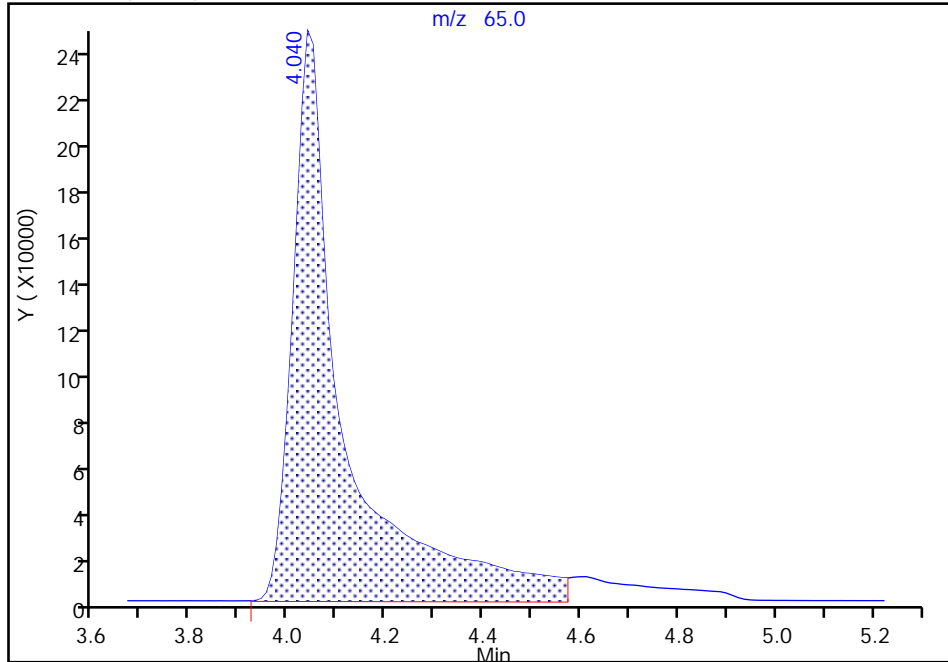
Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25669.D
Injection Date: 13-Mar-2019 13:45:30 Instrument ID: A3UX2
Lims ID: 240-109102-A-4 Lab Sample ID: 240-109102-4
Client ID: MW-83S_030519
Operator ID: 002808 ALS Bottle#: 10 Worklist Smp#: 11
Purge Vol: 15.000 mL Dil. Factor: 1.0000
Method: 8260SIM_X2 Limit Group: MSV SIM ICAL
Column: DB-624 (0.18 mm) Detector: MS SCAN

\$ 5 1,2-Dichloroethane-d4 (Surr), CAS: 17060-07-0

Signal: 1

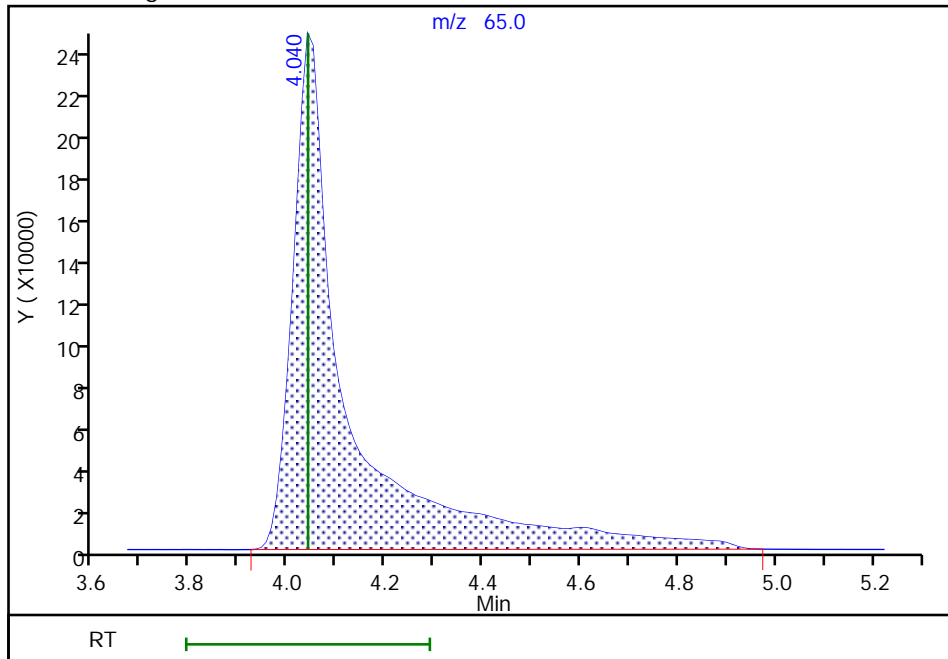
RT: 4.04
Area: 1827908
Amount: 9.245640
Amount Units: ug/l

Processing Integration Results



RT: 4.04
Area: 1938308
Amount: 9.804048
Amount Units: ug/l

Manual Integration Results



Reviewer: macenczaks, 13-Mar-2019 15:49:45

Audit Action: Manually Integrated

Audit Reason: Poor chromatography



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Client Sample ID: MW-82SR_030519 Lab Sample ID: 240-109102-5
 Matrix: Water Lab File ID: X25670.D
 Analysis Method: 8260B SIM Date Collected: 03/05/2019 17:16
 Sample wt/vol: 15 (mL) Date Analyzed: 03/13/2019 14:11
 Soil Aliquot Vol: _____ Dilution Factor: 1
 Soil Extract Vol.: _____ GC Column: ZB-624 ID: 0.53 (mm)
 % Moisture: _____ Level: (low/med) Low
 Analysis Batch No.: 371373 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
123-91-1	1,4-Dioxane	2.0	U	2.0	0.86

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	101		63-125

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25670.D
 Lims ID: 240-109102-A-5
 Client ID: MW-82SR_030519
 Sample Type: Client
 Inject. Date: 13-Mar-2019 14:11:30 ALS Bottle#: 11 Worklist Smp#: 12
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: 240-109102-a-5
 Misc. Info.: 240-0085117-012
 Operator ID: 002808 Instrument ID: A3UX2
 Method: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 14-Mar-2019 09:22:23 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0335

First Level Reviewer: macenczaks Date: 13-Mar-2019 15:50:14

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/l	Flags
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.042	4.040	0.002	100	1977970	10.1	
* 7 Fluorobenzene	96	4.611	4.611	0.000	100	6149477	10.0	
* 9 Dioxane-d8 (IS)	96	5.898	5.891	0.007	92	129396	200.0	

Reagents:

vm150is_00135 Amount Added: 1.00 Units: uL Run Reagent
 vmDist_H2o_00139 Amount Added: 0.00 Units: Run Reagent

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25670.D

Injection Date: 13-Mar-2019 14:11:30

Instrument ID: A3UX2

Operator ID: 002808

Lims ID: 240-109102-A-5

Lab Sample ID: 240-109102-5

Worklist Smp#: 12

Client ID: MW-82SR_030519

Purge Vol: 15.000 mL

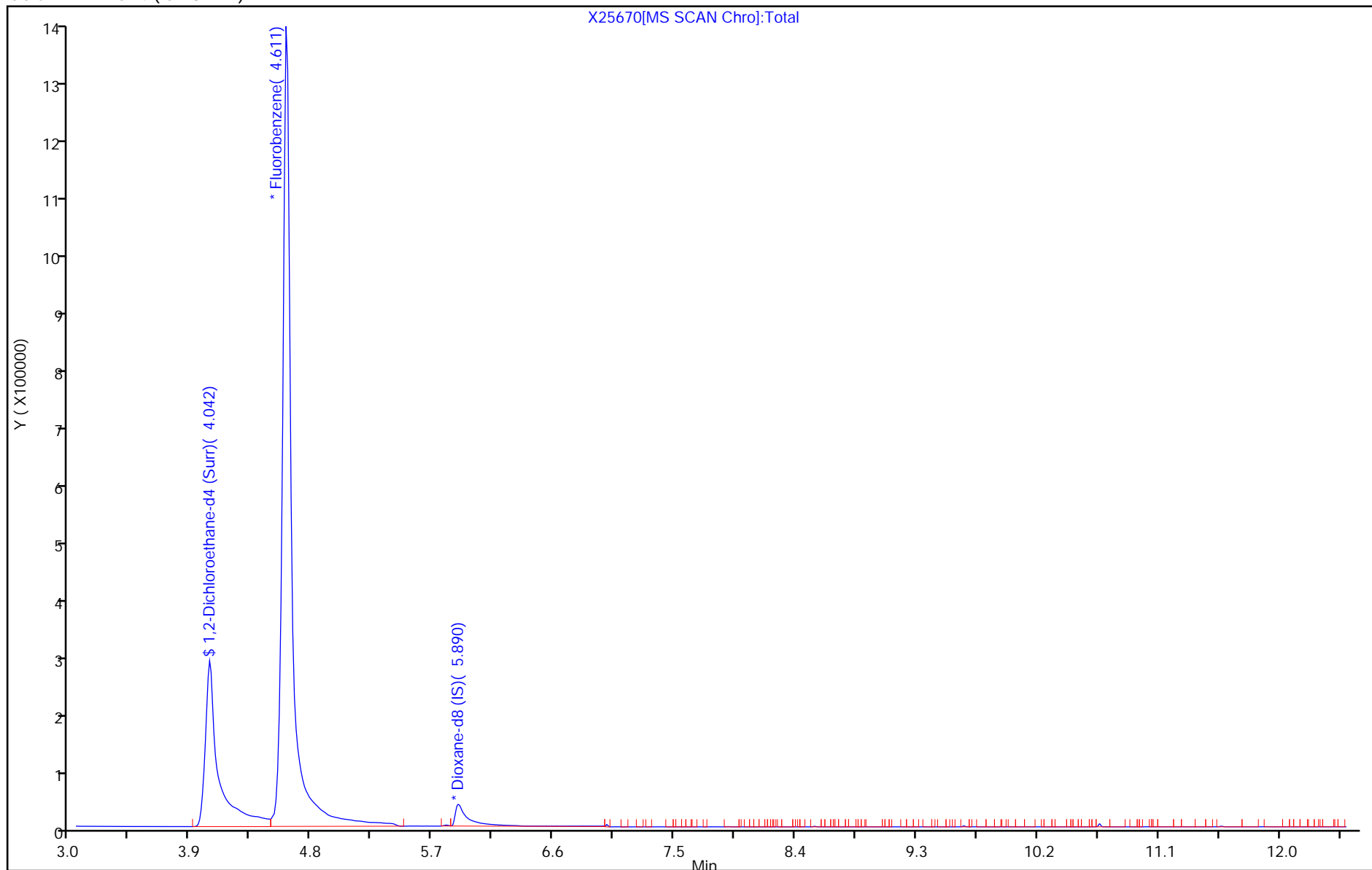
Dil. Factor: 1.0000

ALS Bottle#: 11

Method: 8260SIM_X2

Limit Group: MSV SIM ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Recovery Report

Data File: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\X25670.D
 Lims ID: 240-109102-A-5
 Client ID: MW-82SR_030519
 Sample Type: Client
 Inject. Date: 13-Mar-2019 14:11:30 ALS Bottle#: 11 Worklist Smp#: 12
 Purge Vol: 15.000 mL Dil. Factor: 1.0000
 Sample Info: 240-109102-a-5
 Misc. Info.: 240-0085117-012
 Operator ID: 002808 Instrument ID: A3UX2
 Method: \\chromna\Canton\ChromData\A3UX2\20190313-85117.b\8260SIM_X2.m
 Limit Group: MSV SIM ICAL
 Last Update: 14-Mar-2019 09:22:23 Calib Date: 03-Aug-2018 19:31:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Canton\ChromData\A3UX2\20180803-78614.b\X23823.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0335

First Level Reviewer: macenczaks

Date: 13-Mar-2019 15:50:14

Compound	Amount Added	Amount Recovered	% Rec.
\$ 5 1,2-Dichloroethane-d4 (Surr)	10.0	10.1	101.01

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: ICV 240-367104/14 Calibration Date: 02/07/2019 15:22
 Instrument ID: A3UX11 Calib Start Date: 02/07/2019 13:09
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 02/07/2019 15:00
 Lab File ID: UXJ8027.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Dichlorodifluoromethane	Ave	0.2965	0.2155		0.00727	0.0100	-27.3	50.0
Chloromethane	Ave	0.2437	0.1916	0.1000	0.00786	0.0100	-21.4	50.0
Butadiene	Ave	0.2273	0.1923		0.00846	0.0100	-15.4	50.0
Vinyl chloride	Ave	0.2648	0.2502		0.00945	0.0100	-5.5	20.0
Bromomethane	Ave	0.1749	0.1599		0.00914	0.0100	-8.6	50.0
Chloroethane	Ave	0.1395	0.1398		0.0100	0.0100	0.3	50.0
Dichlorofluoromethane	Ave	0.3514	0.3586		0.0102	0.0100	2.0	50.0
Trichlorofluoromethane	Ave	0.3720	0.3714		0.00998	0.0100	-0.2	50.0
Ethyl ether	Ave	0.1858	0.1927		0.0104	0.0100	3.7	50.0
Acrolein	Ave	0.0478	0.0476		0.0498	0.0500	-0.3	50.0
1,1-Dichloroethene	Ave	0.2297	0.2301		0.0100	0.0100	0.2	20.0
1,1,2-Trichloro-1,2,2-trifluoroethane	Ave	0.1930	0.1932		0.0100	0.0100	0.1	50.0
Acetone	Ave	0.1072	0.1040		0.0194	0.0200	-3.0	50.0
Iodomethane	Ave	0.4080	0.4225		0.0104	0.0100	3.5	50.0
Carbon disulfide	Ave	0.7535	0.7034		0.00934	0.0100	-6.6	50.0
3-Chloro-1-propene	Ave	0.1694	0.1694		0.0100	0.0100	-0.0	50.0
Methyl acetate	Ave	0.2292	0.2239		0.0195	0.0200	-2.3	50.0
Methylene Chloride	Lin1		0.2589		0.00975	0.0100	-2.5	50.0
2-Methyl-2-propanol	Ave	0.0387	0.0366		0.0946	0.100	-5.4	50.0
Acrylonitrile	Ave	0.1197	0.1160		0.0969	0.100	-3.1	50.0
trans-1,2-Dichloroethene	Ave	0.2733	0.2926		0.0107	0.0100	7.1	50.0
Methyl tert-butyl ether	Ave	0.6104	0.6251		0.0102	0.0100	2.4	50.0
Hexane	Lin1		0.0663		0.0101	0.0100	0.8	20.0
1,1-Dichloroethane	Ave	0.4478	0.4612	0.1000	0.0103	0.0100	3.0	50.0
Vinyl acetate	Ave	0.4212	0.4555		0.0108	0.0100	8.1	50.0
2,2-Dichloropropane	Ave	0.0660	0.0608		0.00921	0.0100	-7.9	50.0
cis-1,2-Dichloroethene	Ave	0.2845	0.2960		0.0104	0.0100	4.0	50.0
2-Butanone (MEK)	Ave	0.1351	0.1368		0.0203	0.0200	1.3	50.0
Chlorobromomethane	Ave	0.1457	0.1446		0.00993	0.0100	-0.7	50.0
Tetrahydrofuran	Ave	0.0906	0.0830		0.0183	0.0200	-8.4	50.0
Chloroform	Ave	0.4338	0.4439		0.0102	0.0100	2.3	20.0
1,1,1-Trichloroethane	Ave	0.3774	0.3805		0.0101	0.0100	0.8	50.0
Cyclohexane	Ave	0.3509	0.3668		0.0105	0.0100	4.5	50.0
1,1-Dichloropropene	Ave	0.3456	0.3570		0.0103	0.0100	3.3	50.0
Carbon tetrachloride	Ave	0.3572	0.3548		0.00993	0.0100	-0.7	50.0
Isobutyl alcohol	Ave	0.0197	0.0199		0.254	0.250	1.4	50.0
Benzene	Ave	1.001	1.000		0.00998	0.0100	-0.2	50.0
1,2-Dichloroethane	Ave	0.3480	0.3422		0.00983	0.0100	-1.7	50.0
n-Heptane	Lin1		0.0515		0.00896	0.0100	-10.4	50.0
Trichloroethene	Ave	0.2978	0.2982		0.0100	0.0100	0.2	50.0

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: ICV 240-367104/14 Calibration Date: 02/07/2019 15:22
 Instrument ID: A3UX11 Calib Start Date: 02/07/2019 13:09
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 02/07/2019 15:00
 Lab File ID: UXJ8027.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Methylcyclohexane	Ave	0.3885	0.3731		0.00960	0.0100	-4.0	50.0
1,2-Dichloropropane	Ave	0.2100	0.2111		0.0101	0.0100	0.5	20.0
Dibromomethane	Ave	0.1475	0.1557		0.0106	0.0100	5.6	50.0
1,4-Dioxane	Lin1		0.0027		0.163	0.200	-18.4	50.0
Dichlorobromomethane	Ave	0.3202	0.3096		0.00967	0.0100	-3.3	50.0
2-Chloroethyl vinyl ether	Ave	0.1287	0.1154		0.00896	0.0100	-10.4	50.0
cis-1,3-Dichloropropene	Ave	0.3323	0.3447		0.0104	0.0100	3.8	50.0
4-Methyl-2-pentanone (MIBK)	Ave	0.2408	0.2395		0.0199	0.0200	-0.5	50.0
Toluene	Ave	1.444	1.486		0.0103	0.0100	2.9	20.0
trans-1,3-Dichloropropene	Ave	0.4702	0.4369		0.00929	0.0100	-7.1	50.0
Ethyl methacrylate	Ave	0.3915	0.3637		0.00929	0.0100	-7.1	50.0
1,1,2-Trichloroethane	Ave	0.3033	0.2899		0.00956	0.0100	-4.4	50.0
Tetrachloroethene	Ave	0.3229	0.3283		0.0102	0.0100	1.7	50.0
1,3-Dichloropropane	Ave	0.4858	0.4860		0.0100	0.0100	0.0	50.0
2-Hexanone	Ave	0.2590	0.2563		0.0198	0.0200	-1.0	50.0
Chlorodibromomethane	Ave	0.3521	0.3693		0.0105	0.0100	4.9	50.0
Ethylene Dibromide	Ave	0.3126	0.2892		0.00925	0.0100	-7.5	50.0
Chlorobenzene	Ave	0.9680	0.9501	0.3000	0.00982	0.0100	-1.8	50.0
1,1,1,2-Tetrachloroethane	Ave	0.3713	0.3714		0.0100	0.0100	0.0	50.0
Ethylbenzene	Ave	0.4859	0.4717		0.00971	0.0100	-2.9	20.0
m-Xylene & p-Xylene	Ave	0.5666	0.5960		0.0105	0.0100	5.2	50.0
o-Xylene	Ave	0.5483	0.5526		0.0101	0.0100	0.8	50.0
Styrene	Ave	0.9241	0.9199		0.00995	0.0100	-0.5	50.0
Bromoform	Ave	0.2804	0.2771	0.1000	0.00988	0.0100	-1.2	50.0
Isopropylbenzene	Ave	1.410	1.425		0.0101	0.0100	1.0	50.0
1,1,2,2-Tetrachloroethane	Ave	0.8110	0.7825	0.3000	0.00965	0.0100	-3.5	50.0
Bromobenzene	Ave	0.7495	0.7090		0.00946	0.0100	-5.4	50.0
1,2,3-Trichloropropane	Ave	0.2816	0.2766		0.00982	0.0100	-1.8	50.0
trans-1,4-Dichloro-2-butene	Ave	0.2588	0.2519		0.00973	0.0100	-2.7	50.0
N-Propylbenzene	Ave	0.7593	0.7548		0.00994	0.0100	-0.6	50.0
2-Chlorotoluene	Ave	0.6805	0.7074		0.0104	0.0100	4.0	50.0
1,3,5-Trimethylbenzene	Ave	2.068	2.065		0.00999	0.0100	-0.1	50.0
4-Chlorotoluene	Ave	0.6970	0.7266		0.0104	0.0100	4.2	50.0
tert-Butylbenzene	Ave	1.769	1.706		0.00965	0.0100	-3.5	50.0
1,2,4-Trimethylbenzene	Ave	2.086	2.075		0.00995	0.0100	-0.5	50.0
sec-Butylbenzene	Ave	2.519	2.561		0.0102	0.0100	1.7	50.0
1,3-Dichlorobenzene	Ave	1.363	1.309		0.00961	0.0100	-3.9	50.0
4-Isopropyltoluene	Ave	2.229	2.248		0.0101	0.0100	0.9	50.0
1,4-Dichlorobenzene	Ave	1.462	1.373		0.00939	0.0100	-6.1	50.0
1,2-Dichlorobenzene	Ave	1.275	1.209		0.00948	0.0100	-5.2	50.0
n-Butylbenzene	Ave	1.818	1.699		0.00935	0.0100	-6.5	50.0

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: ICV 240-367104/14 Calibration Date: 02/07/2019 15:22
 Instrument ID: A3UX11 Calib Start Date: 02/07/2019 13:09
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 02/07/2019 15:00
 Lab File ID: UXJ8027.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,2-Dibromo-3-Chloropropane	Lin1		0.2374		0.00979	0.0100	-2.1	50.0
1,2,4-Trichlorobenzene	Ave	0.8021	0.7506		0.00936	0.0100	-6.4	50.0
Hexachlorobutadiene	Ave	0.3705	0.2825		0.00762	0.0100	-23.8	50.0
Naphthalene	Ave	2.134	1.995		0.00935	0.0100	-6.5	50.0
1,2,3-Trichlorobenzene	Ave	0.7522	0.7514		0.00999	0.0100	-0.1	50.0
Dibromofluoromethane (Surr)	Ave	0.2333	0.2347		0.0201	0.0200	0.6	50.0
1,2-Dichloroethane-d4 (Surr)	Ave	0.2824	0.2726		0.0193	0.0200	-3.5	50.0
Toluene-d8 (Surr)	Ave	1.187	1.203		0.0203	0.0200	1.4	50.0
4-Bromofluorobenzene (Surr)	Ave	0.3796	0.3597		0.0190	0.0200	-5.2	50.0

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8027.D
 Lims ID: ICV
 Client ID:
 Sample Type: ICV
 Inject. Date: 07-Feb-2019 15:22:30 ALS Bottle#: 7 Worklist Smp#: 14
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0084305-014
 Misc. Info.: J90207A-IC,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Sublist:
 Method: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 08-Feb-2019 09:15:38 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0324

First Level Reviewer: evansle

Date: 08-Feb-2019 08:15:49

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	4.659	4.659	0.000	98	1136874	20.0	20.0	
* 2 Chlorobenzene-d5	117	7.309	7.309	0.000	83	742699	20.0	20.0	
* 3 1,4-Dichlorobenzene-d4	152	9.522	9.522	0.000	94	406319	20.0	20.0	
\$ 4 Dibromofluoromethane (Surr	113	4.091	4.103	-0.012	98	266836	20.0	20.1	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.375	4.375	0.000	97	309934	20.0	19.3	
\$ 6 Toluene-d8 (Surr)	98	6.008	6.008	0.000	92	893512	20.0	20.3	
\$ 7 4-Bromofluorobenzene (Surr	95	8.410	8.410	0.000	86	267153	20.0	19.0	
9 Dichlorodifluoromethane	85	1.156	1.168	-0.012	99	122492	10.0	7.27	
10 Chloromethane	50	1.322	1.310	0.012	98	108930	10.0	7.86	
13 Butadiene	54	1.369	1.369	0.000	95	109324	10.0	8.46	
12 Vinyl chloride	62	1.393	1.393	0.000	98	142219	10.0	9.45	
15 Bromomethane	94	1.594	1.606	-0.012	92	90908	10.0	9.14	
16 Chloroethane	64	1.653	1.665	-0.012	98	79487	10.0	10.0	
17 Dichlorofluoromethane	67	1.807	1.807	0.000	98	203824	10.0	10.2	
18 Trichlorofluoromethane	101	1.854	1.831	0.023	95	211117	10.0	9.98	
19 Ethyl ether	59	2.055	2.067	-0.012	89	109517	10.0	10.4	
20 Acrolein	56	2.162	2.162	0.000	98	135366	50.0	49.8	
21 1,1-Dichloroethene	96	2.245	2.245	0.000	98	130822	10.0	10.0	
22 1,1,2-Trichloro-1,2,2-trif	151	2.280	2.280	0.000	91	109807	10.0	10.0	
23 Acetone	43	2.292	2.292	0.000	100	118234	20.0	19.4	
25 Iodomethane	142	2.363	2.363	0.000	95	240134	10.0	10.4	
26 Carbon disulfide	76	2.410	2.411	0.000	99	399812	10.0	9.34	
28 3-Chloro-1-propene	76	2.541	2.541	0.000	88	96291	10.0	10.0	
29 Methyl acetate	43	2.564	2.564	0.000	97	254495	20.0	19.5	
30 Methylene Chloride	84	2.635	2.647	-0.012	87	147143	10.0	9.75	
31 2-Methyl-2-propanol	59	2.754	2.754	0.000	96	208181	100.0	94.6	
32 Acrylonitrile	53	2.848	2.848	0.000	98	659215	100.0	96.9	
34 trans-1,2-Dichloroethene	96	2.872	2.872	0.000	98	166333	10.0	10.7	
33 Methyl tert-butyl ether	73	2.884	2.884	0.000	95	355339	10.0	10.2	
35 Hexane	86	3.097	3.109	-0.012	92	37666	10.0	10.1	

Compound	Sig	RT (min.)	Exp RT (min.)	Diff RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
36 1,1-Dichloroethane	63	3.215	3.215	0.000	97	262156	10.0	10.3	
37 Vinyl acetate	43	3.262	3.274	-0.012	97	258898	10.0	10.8	
41 cis-1,2-Dichloroethene	96	3.700	3.700	0.000	80	168258	10.0	10.4	
43 2,2-Dichloropropane	97	3.700	3.700	0.000	63	34552	10.0	9.21	
42 2-Butanone (MEK)	43	3.712	3.712	0.000	99	155571	20.0	20.3	
47 Chlorobromomethane	128	3.890	3.890	0.000	88	82193	10.0	9.93	
48 Tetrahydrofuran	42	3.937	3.937	0.000	80	94338	20.0	18.3	
49 Chloroform	83	3.961	3.961	0.000	94	252351	10.0	10.2	
50 1,1,1-Trichloroethane	97	4.114	4.114	0.000	98	216310	10.0	10.1	
51 Cyclohexane	56	4.162	4.162	0.000	87	208525	10.0	10.5	
53 Carbon tetrachloride	117	4.256	4.256	0.000	93	201657	10.0	9.93	
52 1,1-Dichloropropene	75	4.245	4.256	-0.011	95	202920	10.0	10.3	
54 Isobutyl alcohol	41	4.363	4.363	0.000	87	185080	250.0	253.6	
55 Benzene	78	4.422	4.422	0.000	95	568169	10.0	9.98	
56 1,2-Dichloroethane	62	4.434	4.434	0.000	97	194510	10.0	9.83	
58 n-Heptane	100	4.659	4.659	0.000	44	29291	10.0	8.96	
60 Trichloroethene	130	4.966	4.966	0.000	93	169532	10.0	10.0	
62 Methylcyclohexane	83	5.132	5.132	0.000	89	212078	10.0	9.60	
63 1,2-Dichloropropane	63	5.156	5.156	0.000	95	119996	10.0	10.1	
65 Dibromomethane	93	5.250	5.250	0.000	90	88497	10.0	10.6	
66 1,4-Dioxane	88	5.274	5.274	0.000	94	31141	200.0	163.1	
67 Dichlorobromomethane	83	5.392	5.392	0.000	100	175974	10.0	9.67	
69 2-Chloroethyl vinyl ether	63	5.653	5.653	0.000	91	65573	10.0	8.96	
70 cis-1,3-Dichloropropene	75	5.771	5.771	0.000	97	195960	10.0	10.4	
71 4-Methyl-2-pentanone (MIBK)	43	5.913	5.913	0.000	95	272309	20.0	19.9	
72 Toluene	91	6.067	6.067	0.000	97	551967	10.0	10.3	
73 trans-1,3-Dichloropropene	75	6.268	6.268	0.000	92	162243	10.0	9.29	
74 Ethyl methacrylate	69	6.363	6.363	0.000	87	135071	10.0	9.29	
75 1,1,2-Trichloroethane	97	6.422	6.422	0.000	90	107634	10.0	9.56	
76 Tetrachloroethene	164	6.552	6.552	0.000	97	121909	10.0	10.2	
77 1,3-Dichloropropane	76	6.576	6.576	0.000	89	180490	10.0	10.0	
78 2-Hexanone	43	6.658	6.659	-0.001	95	190355	20.0	19.8	
80 Chlorodibromomethane	129	6.777	6.777	0.000	88	137141	10.0	10.5	
82 Ethylene Dibromide	107	6.871	6.872	-0.001	99	107383	10.0	9.25	
84 Chlorobenzene	112	7.333	7.333	0.000	98	352832	10.0	9.82	
85 1,1,1,2-Tetrachloroethane	131	7.404	7.404	0.000	93	137909	10.0	10.0	
86 Ethylbenzene	106	7.439	7.439	0.000	98	175181	10.0	9.71	
87 m-Xylene & p-Xylene	106	7.546	7.546	0.000	98	221333	10.0	10.5	
88 o-Xylene	106	7.913	7.913	0.000	96	205210	10.0	10.1	
89 Styrene	104	7.925	7.925	0.000	92	341620	10.0	9.95	
90 Bromoform	173	8.090	8.090	0.000	98	102912	10.0	9.88	
91 Isopropylbenzene	105	8.268	8.268	0.000	95	529237	10.0	10.1	
95 Bromobenzene	156	8.552	8.540	0.012	88	144029	10.0	9.46	
94 1,1,2,2-Tetrachloroethane	83	8.552	8.552	0.000	97	158969	10.0	9.65	
96 1,2,3-Trichloropropane	110	8.587	8.587	0.000	85	56194	10.0	9.82	
97 trans-1,4-Dichloro-2-buten	53	8.611	8.611	0.000	82	51170	10.0	9.73	
98 N-Propylbenzene	120	8.658	8.658	0.000	99	153347	10.0	9.94	
99 2-Chlorotoluene	126	8.741	8.741	0.000	98	143712	10.0	10.4	
100 1,3,5-Trimethylbenzene	105	8.836	8.836	0.000	95	419613	10.0	9.99	
101 4-Chlorotoluene	126	8.836	8.848	-0.012	97	147608	10.0	10.4	
102 tert-Butylbenzene	119	9.143	9.155	-0.012	91	346660	10.0	9.65	
104 1,2,4-Trimethylbenzene	105	9.191	9.191	0.000	97	421589	10.0	9.95	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
105 sec-Butylbenzene	105	9.356	9.356	0.000	93	520257	10.0	10.2	
106 1,3-Dichlorobenzene	146	9.451	9.451	0.000	99	266004	10.0	9.61	
107 4-Isopropyltoluene	119	9.510	9.510	0.000	97	456786	10.0	10.1	
108 1,4-Dichlorobenzene	146	9.546	9.546	0.000	97	279004	10.0	9.39	
111 n-Butylbenzene	91	9.901	9.901	0.000	97	345182	10.0	9.35	
112 1,2-Dichlorobenzene	146	9.901	9.901	0.000	96	245592	10.0	9.48	
114 1,2-Dibromo-3-Chloropropan	157	10.670	10.670	0.000	93	48232	10.0	9.79	
116 1,2,4-Trichlorobenzene	180	11.486	11.486	0.000	91	152496	10.0	9.36	
117 Hexachlorobutadiene	225	11.676	11.676	0.000	97	57388	10.0	7.62	
118 Naphthalene	128	11.723	11.723	0.000	96	405313	10.0	9.35	
119 1,2,3-Trichlorobenzene	180	11.971	11.960	0.011	97	152661	10.0	9.99	
S 133 Trihalomethanes, Total	1				0		40.0	40.3	
S 159 Total BTEX	1				0		50.0	50.6	
S 132 Xylenes, Total	106				0		20.0	20.6	

Reagents:

VMFASPW_00287	Amount Added: 8.00	Units: uL	
VMFASAW_00266	Amount Added: 8.00	Units: uL	
VMFASGW_00295	Amount Added: 8.00	Units: uL	
vm50is_stk_A_00002	Amount Added: 2.00	Units: uL	Run Reagent
vm50ss_stk_00079	Amount Added: 2.00	Units: uL	Run Reagent
vmDist_H2o_00138	Amount Added: 0.00	Units:	Run Reagent

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8027.D

Injection Date: 07-Feb-2019 15:22:30

Instrument ID: A3UX11

Operator ID: 43582

Lims ID: ICV

Worklist Smp#: 14

Client ID:

Purge Vol: 5.000 mL

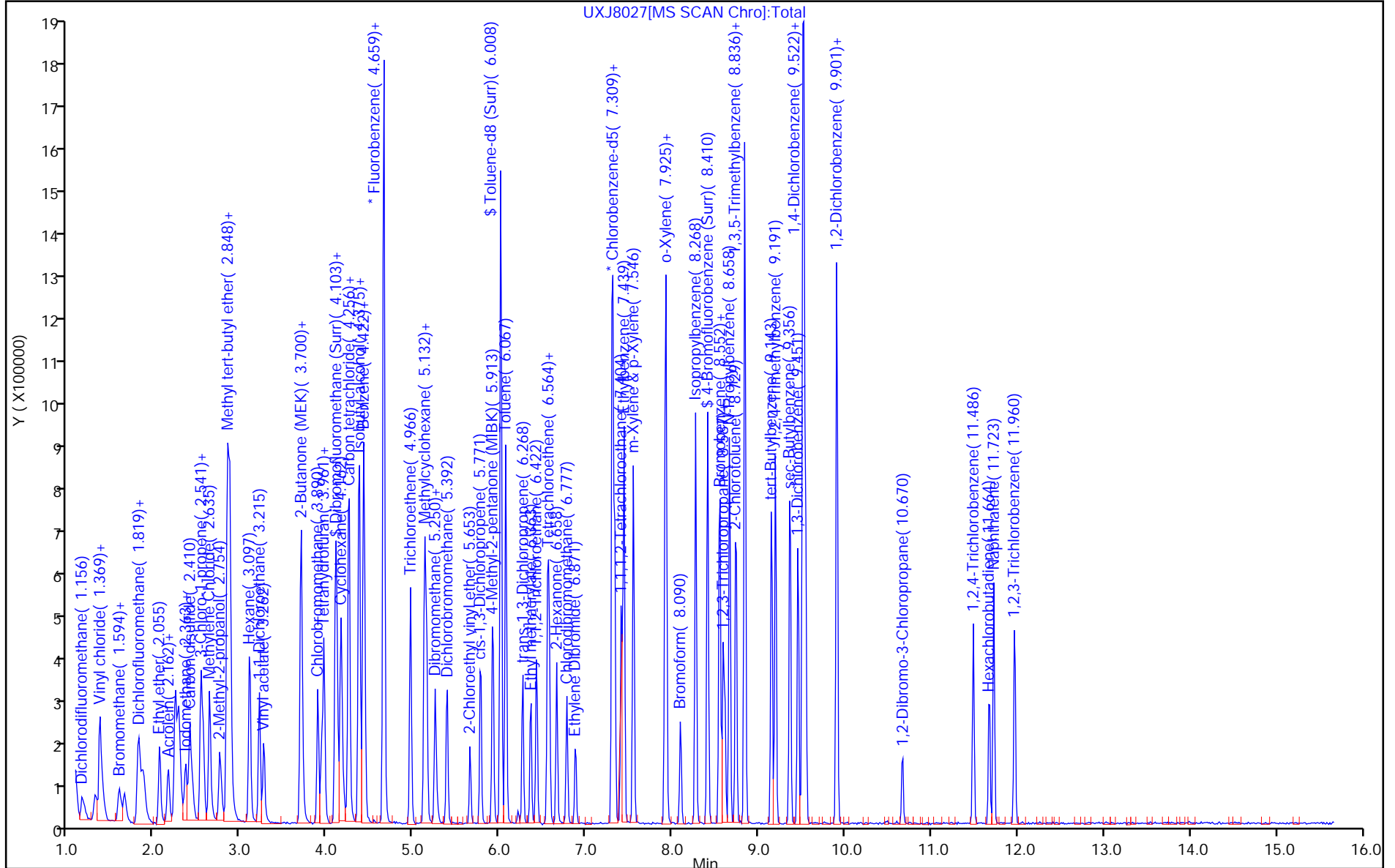
Dil. Factor: 1.0000

ALS Bottle#: 7

Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: ICV 240-367104/15 Calibration Date: 02/07/2019 17:58
 Instrument ID: A3UX11 Calib Start Date: 02/07/2019 13:09
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 02/07/2019 15:00
 Lab File ID: UXJ8034.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Dibromofluoromethane (Surr)	Ave	0.2333	0.2472		0.0212	0.0200	5.9	50.0
1,2-Dichloroethane-d4 (Surr)	Ave	0.2824	0.2744		0.0194	0.0200	-2.8	50.0
Toluene-d8 (Surr)	Ave	1.187	1.173		0.0198	0.0200	-1.2	50.0
4-Bromofluorobenzene (Surr)	Ave	0.3796	0.3662		0.0193	0.0200	-3.5	50.0

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8034.D
 Lims ID: ICV A9
 Client ID:
 Sample Type: ICV
 Inject. Date: 07-Feb-2019 17:58:30 ALS Bottle#: 14 Worklist Smp#: 15
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0084305-015
 Misc. Info.: J90207A-IC,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Sublist:
 Method: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 08-Feb-2019 09:15:38 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0324

First Level Reviewer: evansle

Date: 08-Feb-2019 08:53:04

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	4.659	4.659	0.000	100	1095747	20.0	20.0	
* 2 Chlorobenzene-d5	117	7.309	7.309	0.000	77	685491	20.0	20.0	
* 3 1,4-Dichlorobenzene-d4	152	9.522	9.522	0.000	91	350426	20.0	20.0	
\$ 4 Dibromofluoromethane (Surr	113	4.091	4.103	-0.012	95	270885	20.0	21.2	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.375	4.375	0.000	94	300650	20.0	19.4	
\$ 6 Toluene-d8 (Surr)	98	6.008	6.008	0.000	92	803880	20.0	19.8	
\$ 7 4-Bromofluorobenzene (Surr	95	8.410	8.410	0.000	87	251011	20.0	19.3	
27 Acetonitrile	41	2.517	2.517	0.000	99	186717	100.0	105.4	
39 2-Chloro-1,3-butadiene	53	3.298	3.286	0.012	85	215755	10.0	10.4	
38 Isopropyl ether	87	3.286	3.286	0.000	90	108209	10.0	10.5	
40 Tert-butyl ethyl ether	59	3.594	3.582	0.012	94	315929	10.0	10.5	
44 Propionitrile	54	3.759	3.759	0.000	23	241667	100.0	101.0	
45 Ethyl acetate	43	3.771	3.771	0.000	96	247987	20.0	18.8	
46 Methacrylonitrile	41	3.889	3.890	-0.001	91	907357	100.0	99.8	
57 Tert-amyl methyl ether	73	4.528	4.529	-0.001	92	321983	10.0	10.6	
59 n-Butanol	56	4.919	4.919	0.000	88	130328	250.0	237.6	
61 Ethyl acrylate	55	5.061	5.061	0.000	97	135427	10.0	9.07	
64 Methyl methacrylate	41	5.262	5.262	0.000	85	188872	20.0	19.6	
68 2-Nitropropane	41	5.582	5.582	0.000	94	84065	20.0	18.9	
79 n-Butyl acetate	43	6.777	6.789	-0.012	97	141764	10.0	9.48	
83 1-Chlorohexane	91	7.321	7.321	0.000	94	150288	10.0	10.3	
92 Cyclohexanone	55	8.339	8.339	0.000	82	43960	100.0	91.9	
103 Pentachloroethane	167	9.155	9.155	0.000	92	126476	20.0	14.8	
109 1,2,3-Trimethylbenzene	105	9.605	9.605	0.000	94	388080	10.0	10.3	
110 Benzyl chloride	126	9.676	9.676	0.000	99	59598	10.0	9.03	
115 1,3,5-Trichlorobenzene	180	10.883	10.883	0.000	93	168026	10.0	10.1	
120 2-Methylnaphthalene	142	13.048	13.048	0.000	67	198099	20.0	12.0	

Reagents:

VMFASA9W_00218	Amount Added: 8.00	Units: uL	
vm50is_stk_A_00002	Amount Added: 2.00	Units: uL	Run Reagent
vm50ss_stk_00079	Amount Added: 2.00	Units: uL	Run Reagent
vmDist_H2o_00138	Amount Added: 0.00	Units:	Run Reagent

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Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8034.D

Injection Date: 07-Feb-2019 17:58:30

Instrument ID: A3UX11

Operator ID: 43582

Lims ID: ICV A9

Worklist Smp#: 15

Client ID:

Purge Vol: 5.000 mL

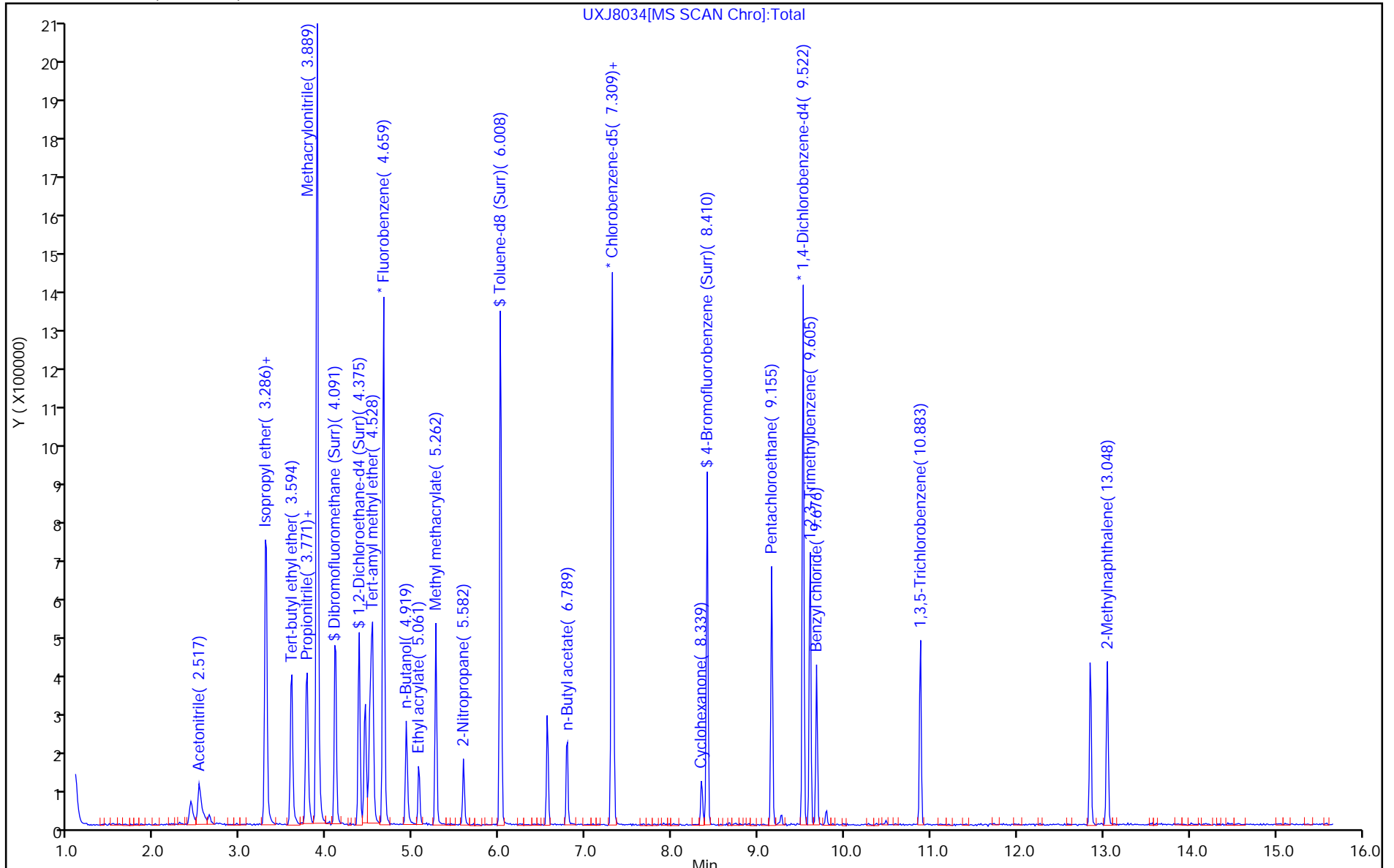
Dil. Factor: 1.0000

ALS Bottle#: 14

Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: ICV 240-367104/15 Calibration Date: 02/07/2019 17:58
 Instrument ID: A3UX11 Calib Start Date: 02/07/2019 15:44
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 02/07/2019 17:35
 Lab File ID: UXJ8034.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Acetonitrile	Ave	0.0323	0.0341		0.105	0.100	5.4	50.0
Isopropyl ether	Ave	0.1872	0.1975		0.0105	0.0100	5.5	50.0
2-Chloro-1,3-butadiene	Ave	0.3783	0.3938		0.0104	0.0100	4.1	50.0
Tert-butyl ethyl ether	Ave	0.5509	0.5767		0.0105	0.0100	4.7	50.0
Propionitrile	Ave	0.0437	0.0441		0.101	0.100	1.0	50.0
Ethyl acetate	Ave	0.2406	0.2263		0.0188	0.0200	-5.9	50.0
Methacrylonitrile	Ave	0.1660	0.1656		0.0998	0.100	-0.2	50.0
Tert-amyl methyl ether	Ave	0.5547	0.5877		0.0106	0.0100	5.9	50.0
n-Butanol	Ave	0.0160	0.0152		0.238	0.250	-5.0	50.0
Ethyl acrylate	Ave	0.2725	0.2472		0.00907	0.0100	-9.3	50.0
Methyl methacrylate	Ave	0.1760	0.1724		0.0196	0.0200	-2.1	50.0
2-Nitropropane	Ave	0.0812	0.0767		0.0189	0.0200	-5.6	50.0
n-Butyl acetate	Ave	0.2731	0.2588		0.00948	0.0100	-5.2	50.0
1-Chlorohexane	Ave	0.4268	0.4385		0.0103	0.0100	2.7	50.0
Cyclohexanone	Lin1		0.0251		0.0919	0.100	-8.1	50.0
Pentachloroethane	Ave	0.2485	0.1845		0.0148	0.0200	-25.8	50.0
1,2,3-Trimethylbenzene	Ave	2.159	2.215		0.0103	0.0100	2.6	50.0
Benzyl chloride	Ave	0.3768	0.3402		0.00903	0.0100	-9.7	50.0
1,3,5-Trichlorobenzene	Ave	0.9457	0.9590		0.0101	0.0100	1.4	50.0
2-Methylnaphthalene	Lin		0.5653		0.0120	0.0200	-40.0	50.0

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8034.D
 Lims ID: ICV A9
 Client ID:
 Sample Type: ICV
 Inject. Date: 07-Feb-2019 17:58:30 ALS Bottle#: 14 Worklist Smp#: 15
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0084305-015
 Misc. Info.: J90207A-IC,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Sublist:
 Method: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 08-Feb-2019 09:15:38 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0324

First Level Reviewer: evansle

Date: 08-Feb-2019 08:53:04

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	4.659	4.659	0.000	100	1095747	20.0	20.0	
* 2 Chlorobenzene-d5	117	7.309	7.309	0.000	77	685491	20.0	20.0	
* 3 1,4-Dichlorobenzene-d4	152	9.522	9.522	0.000	91	350426	20.0	20.0	
\$ 4 Dibromofluoromethane (Surr	113	4.091	4.103	-0.012	95	270885	20.0	21.2	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.375	4.375	0.000	94	300650	20.0	19.4	
\$ 6 Toluene-d8 (Surr)	98	6.008	6.008	0.000	92	803880	20.0	19.8	
\$ 7 4-Bromofluorobenzene (Surr	95	8.410	8.410	0.000	87	251011	20.0	19.3	
27 Acetonitrile	41	2.517	2.517	0.000	99	186717	100.0	105.4	
39 2-Chloro-1,3-butadiene	53	3.298	3.286	0.012	85	215755	10.0	10.4	
38 Isopropyl ether	87	3.286	3.286	0.000	90	108209	10.0	10.5	
40 Tert-butyl ethyl ether	59	3.594	3.582	0.012	94	315929	10.0	10.5	
44 Propionitrile	54	3.759	3.759	0.000	23	241667	100.0	101.0	
45 Ethyl acetate	43	3.771	3.771	0.000	96	247987	20.0	18.8	
46 Methacrylonitrile	41	3.889	3.890	-0.001	91	907357	100.0	99.8	
57 Tert-amyl methyl ether	73	4.528	4.529	-0.001	92	321983	10.0	10.6	
59 n-Butanol	56	4.919	4.919	0.000	88	130328	250.0	237.6	
61 Ethyl acrylate	55	5.061	5.061	0.000	97	135427	10.0	9.07	
64 Methyl methacrylate	41	5.262	5.262	0.000	85	188872	20.0	19.6	
68 2-Nitropropane	41	5.582	5.582	0.000	94	84065	20.0	18.9	
79 n-Butyl acetate	43	6.777	6.789	-0.012	97	141764	10.0	9.48	
83 1-Chlorohexane	91	7.321	7.321	0.000	94	150288	10.0	10.3	
92 Cyclohexanone	55	8.339	8.339	0.000	82	43960	100.0	91.9	
103 Pentachloroethane	167	9.155	9.155	0.000	92	126476	20.0	14.8	
109 1,2,3-Trimethylbenzene	105	9.605	9.605	0.000	94	388080	10.0	10.3	
110 Benzyl chloride	126	9.676	9.676	0.000	99	59598	10.0	9.03	
115 1,3,5-Trichlorobenzene	180	10.883	10.883	0.000	93	168026	10.0	10.1	
120 2-Methylnaphthalene	142	13.048	13.048	0.000	67	198099	20.0	12.0	

Reagents:

VMFASA9W_00218	Amount Added: 8.00	Units: uL	
vm50is_stk_A_00002	Amount Added: 2.00	Units: uL	Run Reagent
vm50ss_stk_00079	Amount Added: 2.00	Units: uL	Run Reagent
vmDist_H2o_00138	Amount Added: 0.00	Units:	Run Reagent

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Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8034.D

Injection Date: 07-Feb-2019 17:58:30

Instrument ID: A3UX11

Operator ID: 43582

Lims ID: ICV A9

Worklist Smp#: 15

Client ID:

Purge Vol: 5.000 mL

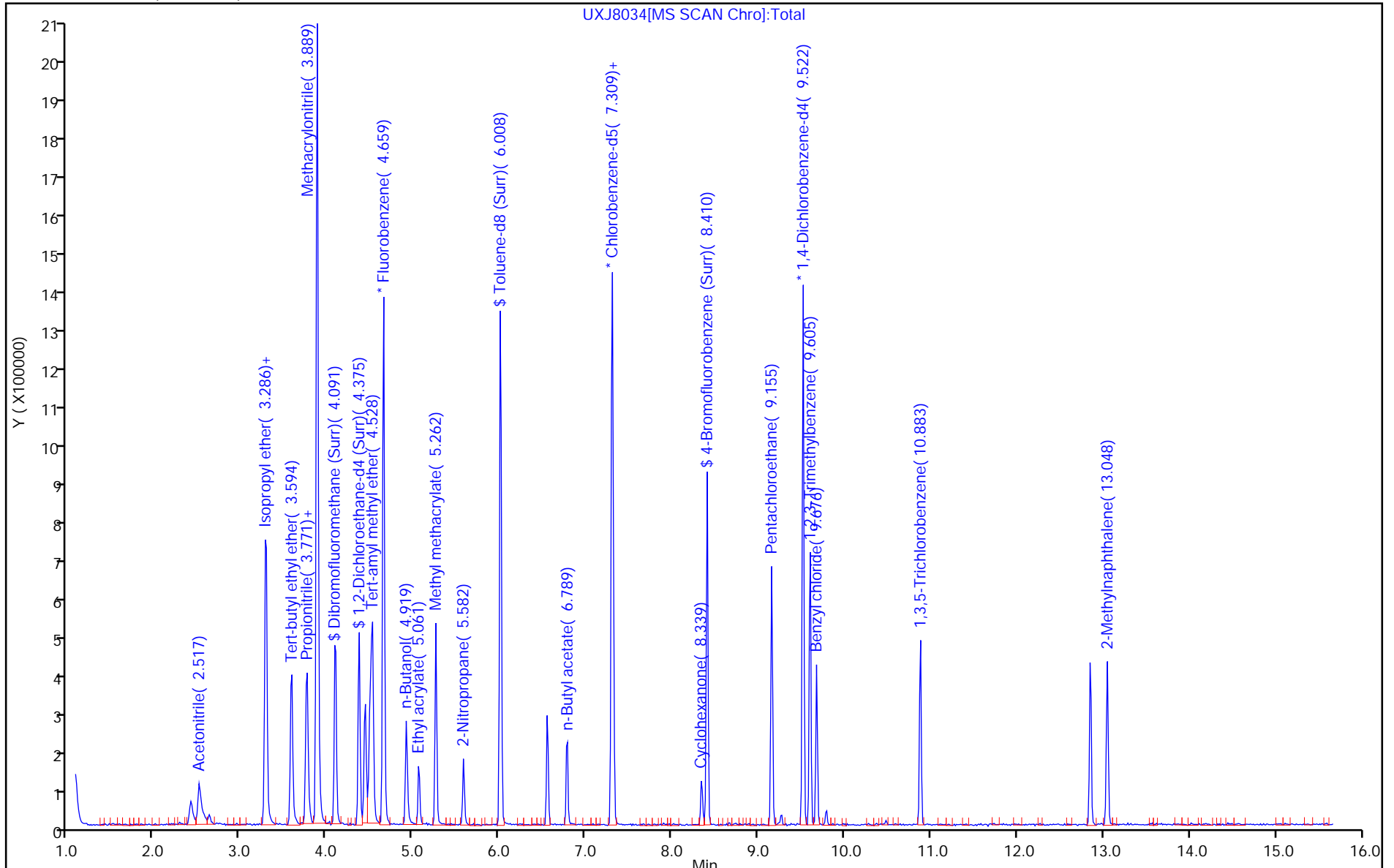
Dil. Factor: 1.0000

ALS Bottle#: 14

Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: CCVIS 240-371553/2 Calibration Date: 03/14/2019 10:07
 Instrument ID: A3UX11 Calib Start Date: 02/07/2019 13:09
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 02/07/2019 15:00
 Lab File ID: UXJ8779.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Dichlorodifluoromethane	Ave	0.2965	0.2591		0.00874	0.0100	-12.6	50.0
Chloromethane	Ave	0.2437	0.3738	0.1000	0.0153	0.0100	53.4*	50.0
Butadiene	Ave	0.2273	0.3069		0.0135	0.0100	35.0	50.0
Vinyl chloride	Ave	0.2648	0.3525		0.0133	0.0100	33.1*	20.0
Bromomethane	Ave	0.1749	0.1843		0.0105	0.0100	5.4	50.0
Chloroethane	Ave	0.1395	0.2054		0.0147	0.0100	47.2	50.0
Dichlorofluoromethane	Ave	0.3514	0.4322		0.0123	0.0100	23.0	50.0
Trichlorofluoromethane	Ave	0.3720	0.3608		0.00970	0.0100	-3.0	50.0
Ethyl ether	Ave	0.1858	0.2590		0.0139	0.0100	39.4	50.0
Acrolein	Ave	0.0478	0.0613		0.0641	0.0500	28.2	50.0
1,1-Dichloroethene	Ave	0.2297	0.2394		0.0104	0.0100	4.2	20.0
1,1,2-Trichloro-1,2,2-trifluoroethane	Ave	0.1930	0.1495		0.00775	0.0100	-22.5	50.0
Acetone	Ave	0.1072	0.1288		0.0240	0.0200	20.1	50.0
Iodomethane	Ave	0.4080	0.3385		0.00830	0.0100	-17.0	50.0
Carbon disulfide	Ave	0.7535	0.8742		0.0116	0.0100	16.0	50.0
3-Chloro-1-propene	Ave	0.1694	0.1859		0.0110	0.0100	9.7	50.0
Methyl acetate	Ave	0.2292	0.3511		0.0306	0.0200	53.2*	50.0
Methylene Chloride	Lin1		0.2858		0.0109	0.0100	8.9	50.0
2-Methyl-2-propanol	Ave	0.0387	0.0331		0.0854	0.100	-14.6	50.0
Acrylonitrile	Ave	0.1197	0.1650		0.138	0.100	37.9	50.0
Methyl tert-butyl ether	Ave	0.6104	0.4213		0.00690	0.0100	-31.0	50.0
trans-1,2-Dichloroethene	Ave	0.2733	0.2611		0.00955	0.0100	-4.5	50.0
Hexane	Lin1		0.0724		0.0109	0.0100	9.4	20.0
1,1-Dichloroethane	Ave	0.4478	0.5416	0.1000	0.0121	0.0100	20.9	50.0
Vinyl acetate	Ave	0.4212	0.6228		0.0148	0.0100	47.9	50.0
cis-1,2-Dichloroethene	Ave	0.2845	0.3085		0.0108	0.0100	8.4	50.0
2,2-Dichloropropane	Ave	0.0660	0.0590		0.00894	0.0100	-10.6	50.0
2-Butanone (MEK)	Ave	0.1351	0.2198		0.0326	0.0200	62.8*	50.0
Chlorobromomethane	Ave	0.1457	0.1144		0.00786	0.0100	-21.4	50.0
Tetrahydrofuran	Ave	0.0906	0.1562		0.0345	0.0200	72.3*	50.0
Chloroform	Ave	0.4338	0.4451		0.0103	0.0100	2.6	20.0
1,1,1-Trichloroethane	Ave	0.3774	0.3274		0.00867	0.0100	-13.3	50.0
Cyclohexane	Ave	0.3509	0.4914		0.0140	0.0100	40.0	50.0
1,1-Dichloropropene	Ave	0.3456	0.3959		0.0115	0.0100	14.5	50.0
Carbon tetrachloride	Ave	0.3572	0.2985		0.00836	0.0100	-16.4	50.0
Isobutyl alcohol	Ave	0.0197	0.0308		0.392	0.250	56.7*	50.0
Benzene	Ave	1.001	1.112		0.0111	0.0100	11.0	50.0
1,2-Dichloroethane	Ave	0.3480	0.3644		0.0105	0.0100	4.7	50.0
n-Heptane	Lin1		0.0560		0.00972	0.0100	-2.8	50.0
Trichloroethene	Ave	0.2978	0.2841		0.00954	0.0100	-4.6	50.0

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: CCVIS 240-371553/2 Calibration Date: 03/14/2019 10:07
 Instrument ID: A3UX11 Calib Start Date: 02/07/2019 13:09
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 02/07/2019 15:00
 Lab File ID: UXJ8779.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Methylcyclohexane	Ave	0.3885	0.4074		0.0105	0.0100	4.9	50.0
1,2-Dichloropropane	Ave	0.2100	0.2870		0.0137	0.0100	36.6*	20.0
Dibromomethane	Ave	0.1475	0.1572		0.0107	0.0100	6.6	50.0
1,4-Dioxane	Lin1		0.0029		0.170	0.200	-15.0	50.0
Dichlorobromomethane	Ave	0.3202	0.3286		0.0103	0.0100	2.6	50.0
2-Chloroethyl vinyl ether	Ave	0.1287	0.1534		0.0238	0.0200	19.1	50.0
cis-1,3-Dichloropropene	Ave	0.3323	0.3766		0.0113	0.0100	13.3	50.0
4-Methyl-2-pentanone (MIBK)	Ave	0.2408	0.3724		0.0309	0.0200	54.7*	50.0
Toluene	Ave	1.444	1.732		0.0120	0.0100	20.0	20.0
trans-1,3-Dichloropropene	Ave	0.4702	0.5750		0.0122	0.0100	22.3	50.0
Ethyl methacrylate	Ave	0.3915	0.5297		0.0135	0.0100	35.3	50.0
1,1,2-Trichloroethane	Ave	0.3033	0.3445		0.0114	0.0100	13.6	50.0
Tetrachloroethene	Ave	0.3229	0.2535		0.00785	0.0100	-21.5	50.0
1,3-Dichloropropane	Ave	0.4858	0.6483		0.0133	0.0100	33.4	50.0
2-Hexanone	Ave	0.2590	0.4697		0.0363	0.0200	81.4*	50.0
Chlorodibromomethane	Ave	0.3521	0.3527		0.0100	0.0100	0.2	50.0
Ethylene Dibromide	Ave	0.3126	0.3335		0.0107	0.0100	6.7	50.0
Chlorobenzene	Ave	0.9680	0.9705	0.3000	0.0100	0.0100	0.3	50.0
1,1,1,2-Tetrachloroethane	Ave	0.3713	0.3150		0.00848	0.0100	-15.2	50.0
Ethylbenzene	Ave	0.4859	0.4867		0.0100	0.0100	0.2	20.0
m-Xylene & p-Xylene	Ave	0.5666	0.5923		0.0105	0.0100	4.5	50.0
o-Xylene	Ave	0.5483	0.5481		0.0100	0.0100	-0.0	50.0
Styrene	Ave	0.9241	0.9793		0.0106	0.0100	6.0	50.0
Bromoform	Ave	0.2804	0.2118	0.1000	0.00755	0.0100	-24.5	50.0
Isopropylbenzene	Ave	1.410	1.383		0.00980	0.0100	-2.0	50.0
Bromobenzene	Ave	0.7495	0.6704		0.00894	0.0100	-10.6	50.0
1,1,2,2-Tetrachloroethane	Ave	0.8110	1.150	0.3000	0.0142	0.0100	41.8	50.0
1,2,3-Trichloropropane	Ave	0.2816	0.3103		0.0110	0.0100	10.2	50.0
trans-1,4-Dichloro-2-butene	Ave	0.2588	0.3832		0.0148	0.0100	48.1	50.0
N-Propylbenzene	Ave	0.7593	0.8252		0.0109	0.0100	8.7	50.0
2-Chlorotoluene	Ave	0.6805	0.7016		0.0103	0.0100	3.1	50.0
1,3,5-Trimethylbenzene	Ave	2.068	2.264		0.0110	0.0100	9.5	50.0
4-Chlorotoluene	Ave	0.6970	0.7230		0.0104	0.0100	3.7	50.0
tert-Butylbenzene	Ave	1.769	1.806		0.0102	0.0100	2.1	50.0
1,2,4-Trimethylbenzene	Ave	2.086	2.350		0.0113	0.0100	12.6	50.0
sec-Butylbenzene	Ave	2.519	2.769		0.0110	0.0100	9.9	50.0
1,3-Dichlorobenzene	Ave	1.363	1.225		0.00898	0.0100	-10.2	50.0
4-Isopropyltoluene	Ave	2.229	2.229		0.0100	0.0100	-0.0	50.0
1,4-Dichlorobenzene	Ave	1.462	1.293		0.00884	0.0100	-11.6	50.0
1,2-Dichlorobenzene	Ave	1.275	1.116		0.00875	0.0100	-12.5	50.0
n-Butylbenzene	Ave	1.818	2.001		0.0110	0.0100	10.0	50.0

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: CCVIS 240-371553/2 Calibration Date: 03/14/2019 10:07
 Instrument ID: A3UX11 Calib Start Date: 02/07/2019 13:09
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 02/07/2019 15:00
 Lab File ID: UXJ8779.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,2-Dibromo-3-Chloropropane	Lin1		0.1831		0.00749	0.0100	-25.1	50.0
1,2,4-Trichlorobenzene	Ave	0.8021	0.5255		0.00655	0.0100	-34.5	50.0
Hexachlorobutadiene	Ave	0.3705	0.2081		0.00562	0.0100	-43.8	50.0
Naphthalene	Ave	2.134	1.549		0.00726	0.0100	-27.4	50.0
1,2,3-Trichlorobenzene	Ave	0.7522	0.5231		0.00695	0.0100	-30.5	50.0
Dibromofluoromethane (Surr)	Ave	0.2333	0.2173		0.0186	0.0200	-6.9	50.0
1,2-Dichloroethane-d4 (Surr)	Ave	0.2824	0.2925		0.0207	0.0200	3.6	50.0
Toluene-d8 (Surr)	Ave	1.187	1.481		0.0250	0.0200	24.8	50.0
4-Bromofluorobenzene (Surr)	Ave	0.3796	0.4285		0.0226	0.0200	12.9	50.0

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8779.D
 Lims ID: CCVIS L4 8260
 Client ID:
 Sample Type: CCVIS
 Inject. Date: 14-Mar-2019 10:07:30 ALS Bottle#: 1 Worklist Smp#: 2
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0085154-002
 Misc. Info.: J90314A,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Sublist: chrom-8260_11*sub86
 Method: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 15-Mar-2019 07:48:16 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0312

First Level Reviewer: evansle

Date: 14-Mar-2019 11:27:00

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	4.659	4.659	0.000	97	1933598	20.0	20.0	
* 2 Chlorobenzene-d5	117	7.297	7.297	0.000	91	1147246	20.0	20.0	
* 3 1,4-Dichlorobenzene-d4	152	9.522	9.522	0.000	97	539090	20.0	20.0	
\$ 4 Dibromofluoromethane (Surr	113	4.091	4.091	0.000	97	420131	20.0	18.6	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.375	4.375	0.000	99	565539	20.0	20.7	
\$ 6 Toluene-d8 (Surr)	98	6.007	6.007	0.000	94	1699510	20.0	25.0	
\$ 7 4-Bromofluorobenzene (Surr	95	8.410	8.410	0.000	99	491578	20.0	22.6	
9 Dichlorodifluoromethane	85	1.156	1.156	0.000	98	250498	10.0	8.74	
10 Chloromethane	50	1.298	1.298	0.000	99	361372	10.0	15.3	
13 Butadiene	54	1.357	1.357	0.000	96	296753	10.0	13.5	
12 Vinyl chloride	62	1.369	1.369	0.000	97	340748	10.0	13.3	
15 Bromomethane	94	1.582	1.582	0.000	91	178169	10.0	10.5	
16 Chloroethane	64	1.653	1.653	0.000	100	198544	10.0	14.7	
17 Dichlorofluoromethane	67	1.795	1.795	0.000	99	417871	10.0	12.3	
18 Trichlorofluoromethane	101	1.830	1.830	0.000	99	348803	10.0	9.70	
19 Ethyl ether	59	2.055	2.055	0.000	95	250378	10.0	13.9	
20 Acrolein	56	2.162	2.162	0.000	98	296107	50.0	64.1	
21 1,1-Dichloroethene	96	2.233	2.233	0.000	94	231409	10.0	10.4	
22 1,1,2-Trichloro-1,2,2-trif	151	2.268	2.268	0.000	96	144523	10.0	7.75	
23 Acetone	43	2.280	2.280	0.000	100	248964	20.0	24.0	
25 Iodomethane	142	2.363	2.363	0.000	98	327286	10.0	8.30	
26 Carbon disulfide	76	2.410	2.410	0.000	99	845164	10.0	11.6	
28 3-Chloro-1-propene	76	2.540	2.540	0.000	93	179727	10.0	11.0	
29 Methyl acetate	43	2.564	2.564	0.000	98	678886	20.0	30.6	
30 Methylene Chloride	84	2.635	2.635	0.000	99	276267	10.0	10.9	
31 2-Methyl-2-propanol	59	2.753	2.753	0.000	93	319722	100.0	85.4	
32 Acrylonitrile	53	2.836	2.836	0.000	99	1595601	100.0	137.9	
33 Methyl tert-butyl ether	73	2.872	2.872	0.000	96	407331	10.0	6.90	
34 trans-1,2-Dichloroethene	96	2.872	2.872	0.000	97	252399	10.0	9.55	
35 Hexane	86	3.108	3.108	0.000	93	69952	10.0	10.9	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
36 1,1-Dichloroethane	63	3.215	3.215	0.000	96	523592	10.0	12.1	
37 Vinyl acetate	43	3.262	3.262	0.000	97	602126	10.0	14.8	
41 cis-1,2-Dichloroethene	96	3.688	3.688	0.000	88	298282	10.0	10.8	
43 2,2-Dichloropropane	97	3.700	3.700	0.000	57	57055	10.0	8.94	
42 2-Butanone (MEK)	43	3.712	3.712	0.000	99	425066	20.0	32.6	
47 Chlorobromomethane	128	3.889	3.889	0.000	92	110639	10.0	7.86	
48 Tetrahydrofuran	42	3.937	3.937	0.000	92	301948	20.0	34.5	
49 Chloroform	83	3.960	3.960	0.000	95	430288	10.0	10.3	
50 1,1,1-Trichloroethane	97	4.114	4.114	0.000	96	316514	10.0	8.67	
51 Cyclohexane	56	4.162	4.162	0.000	92	475077	10.0	14.0	
52 1,1-Dichloropropene	75	4.244	4.244	0.000	94	382710	10.0	11.5	
53 Carbon tetrachloride	117	4.256	4.256	0.000	93	288622	10.0	8.36	
54 Isobutyl alcohol	41	4.363	4.363	0.000	92	441698	250.0	391.8	
55 Benzene	78	4.422	4.422	0.000	97	1074700	10.0	11.1	
56 1,2-Dichloroethane	62	4.434	4.434	0.000	97	352316	10.0	10.5	
58 n-Heptane	100	4.659	4.659	0.000	49	54133	10.0	9.72	
60 Trichloroethene	130	4.966	4.966	0.000	98	274615	10.0	9.54	
62 Methylcyclohexane	83	5.132	5.132	0.000	95	393884	10.0	10.5	
63 1,2-Dichloropropane	63	5.144	5.144	0.000	97	277420	10.0	13.7	
65 Dibromomethane	93	5.250	5.250	0.000	89	151948	10.0	10.7	
66 1,4-Dioxane	88	5.274	5.274	0.000	95	55361	200.0	170.0	
67 Dichlorobromomethane	83	5.380	5.380	0.000	99	317731	10.0	10.3	
69 2-Chloroethyl vinyl ether	63	5.653	5.653	0.000	91	296577	20.0	23.8	
70 cis-1,3-Dichloropropene	75	5.771	5.771	0.000	94	364068	10.0	11.3	
71 4-Methyl-2-pentanone (MIBK)	43	5.913	5.913	0.000	98	720037	20.0	30.9	
72 Toluene	91	6.067	6.067	0.000	98	993731	10.0	12.0	
73 trans-1,3-Dichloropropene	75	6.268	6.268	0.000	94	329818	10.0	12.2	
74 Ethyl methacrylate	69	6.351	6.351	0.000	93	303830	10.0	13.5	
75 1,1,2-Trichloroethane	97	6.422	6.422	0.000	93	197608	10.0	11.4	
76 Tetrachloroethene	164	6.552	6.552	0.000	93	145409	10.0	7.85	
77 1,3-Dichloropropane	76	6.575	6.575	0.000	93	371896	10.0	13.3	
78 2-Hexanone	43	6.658	6.658	0.000	98	538873	20.0	36.3	
80 Chlorodibromomethane	129	6.777	6.777	0.000	91	202319	10.0	10.0	
82 Ethylene Dibromide	107	6.871	6.871	0.000	98	191275	10.0	10.7	
84 Chlorobenzene	112	7.333	7.333	0.000	94	556715	10.0	10.0	
85 1,1,1,2-Tetrachloroethane	131	7.404	7.404	0.000	95	180686	10.0	8.48	
86 Ethylbenzene	106	7.439	7.439	0.000	99	279154	10.0	10.0	
87 m-Xylene & p-Xylene	106	7.546	7.546	0.000	97	339754	10.0	10.5	
88 o-Xylene	106	7.913	7.913	0.000	95	314389	10.0	10.0	
89 Styrene	104	7.924	7.924	0.000	95	561729	10.0	10.6	
90 Bromoform	173	8.090	8.090	0.000	94	121490	10.0	7.55	
91 Isopropylbenzene	105	8.268	8.268	0.000	96	793210	10.0	9.80	
95 Bromobenzene	156	8.540	8.540	0.000	95	180701	10.0	8.94	
94 1,1,2,2-Tetrachloroethane	83	8.552	8.552	0.000	93	310026	10.0	14.2	
96 1,2,3-Trichloropropane	110	8.587	8.587	0.000	86	83639	10.0	11.0	
97 trans-1,4-Dichloro-2-buten	53	8.611	8.611	0.000	83	103287	10.0	14.8	
98 N-Propylbenzene	120	8.658	8.658	0.000	100	222426	10.0	10.9	
99 2-Chlorotoluene	126	8.741	8.741	0.000	96	189119	10.0	10.3	
100 1,3,5-Trimethylbenzene	105	8.836	8.836	0.000	85	610353	10.0	11.0	
101 4-Chlorotoluene	126	8.836	8.836	0.000	98	194888	10.0	10.4	
102 tert-Butylbenzene	119	9.143	9.143	0.000	94	486818	10.0	10.2	
104 1,2,4-Trimethylbenzene	105	9.191	9.191	0.000	95	633389	10.0	11.3	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
105 sec-Butylbenzene	105	9.356	9.356	0.000	95	746380	10.0	11.0	
106 1,3-Dichlorobenzene	146	9.451	9.451	0.000	96	330083	10.0	8.98	
107 4-Isopropyltoluene	119	9.510	9.510	0.000	97	600696	10.0	10.0	
108 1,4-Dichlorobenzene	146	9.546	9.546	0.000	92	348425	10.0	8.84	
112 1,2-Dichlorobenzene	146	9.901	9.901	0.000	90	300760	10.0	8.75	
111 n-Butylbenzene	91	9.901	9.901	0.000	98	539247	10.0	11.0	
114 1,2-Dibromo-3-Chloropropan	157	10.670	10.670	0.000	77	49361	10.0	7.49	
116 1,2,4-Trichlorobenzene	180	11.486	11.486	0.000	93	141634	10.0	6.55	
117 Hexachlorobutadiene	225	11.675	11.675	0.000	95	56094	10.0	5.62	
118 Naphthalene	128	11.723	11.723	0.000	97	417559	10.0	7.26	
119 1,2,3-Trichlorobenzene	180	11.959	11.959	0.000	94	141000	10.0	6.95	
S 159 Total BTEX	1				0		50.0	53.6	
S 133 Trihalomethanes, Total	1				0		40.0	38.1	
S 132 Xylenes, Total	106				0		20.0	20.4	

Reagents:

VMRGAS_00285	Amount Added: 8.00	Units: uL	
VMAROLISTDW_00288	Amount Added: 8.00	Units: uL	
VMRPRIMW_00325	Amount Added: 8.00	Units: uL	
vm40ml_vials_00009	Amount Added: 0.00	Units:	Run Reagent
vmDist_H2o_00139	Amount Added: 0.00	Units:	Run Reagent
vm50ss_stk_00080	Amount Added: 2.00	Units: uL	Run Reagent
VM50IS_00075	Amount Added: 2.00	Units: uL	Run Reagent

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8779.D

Injection Date: 14-Mar-2019 10:07:30

Instrument ID: A3UX11

Operator ID: 43582

Lims ID: CCVIS L4 8260

Worklist Smp#: 2

Client ID:

Purge Vol: 5.000 mL

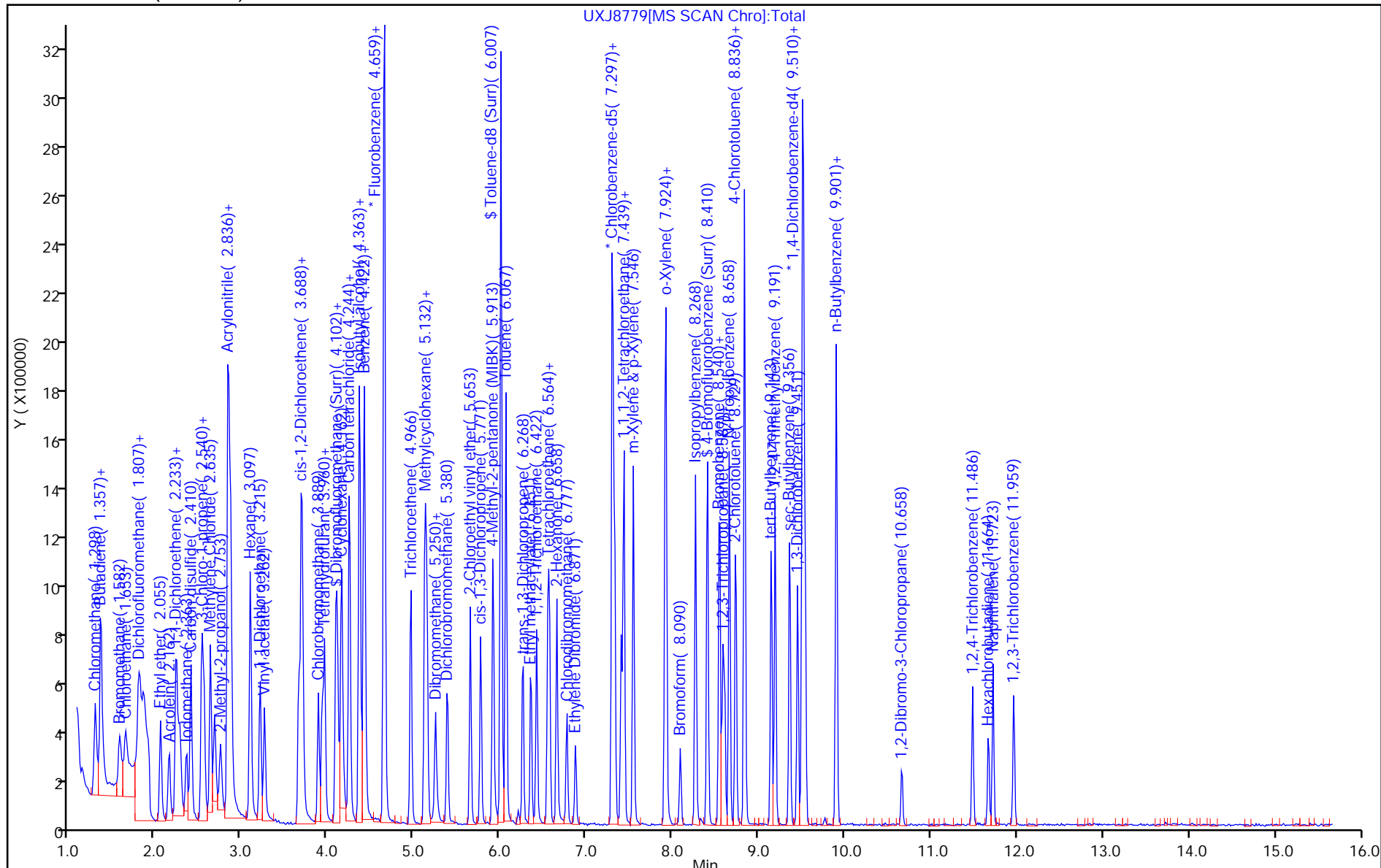
Dil. Factor: 1.0000

ALS Bottle#: 1

Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: ICV 240-364785/14 Calibration Date: 01/21/2019 13:48
 Instrument ID: A3UX15 Calib Start Date: 01/21/2019 11:37
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 01/21/2019 13:26
 Lab File ID: UXC8416.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Dichlorodifluoromethane	Ave	0.3072	0.3077		0.0100	0.0100	0.2	50.0
Chloromethane	Ave	0.3135	0.2939	0.1000	0.00937	0.0100	-6.3	50.0
Vinyl chloride	Ave	0.2885	0.2965		0.0103	0.0100	2.8	20.0
Butadiene	Ave	0.3794	0.3439		0.00906	0.0100	-9.4	50.0
Bromomethane	Ave	0.2308	0.2181		0.00945	0.0100	-5.5	50.0
Chloroethane	Ave	0.2041	0.1937		0.00949	0.0100	-5.1	50.0
Dichlorofluoromethane	Ave	0.5102	0.5038		0.00987	0.0100	-1.3	50.0
Trichlorofluoromethane	Ave	0.4337	0.4655		0.0107	0.0100	7.3	50.0
Ethyl ether	Ave	0.2310	0.2499		0.0108	0.0100	8.2	50.0
Acrolein	Ave	0.0317	0.0225		0.0354	0.0500	-29.1	50.0
1,1-Dichloroethene	Ave	0.3005	0.2950		0.00982	0.0100	-1.8	20.0
1,1,2-Trichloro-1,2,2-trifluoroethane	Ave	0.2468	0.2669		0.0108	0.0100	8.1	50.0
Acetone	Lin1		0.0260		0.0217	0.0200	8.7	50.0
Iodomethane	Ave	0.4675	0.5088		0.0109	0.0100	8.8	50.0
Carbon disulfide	Ave	0.7741	0.7832		0.0101	0.0100	1.2	50.0
3-Chloro-1-propene	Ave	0.1698	0.1835		0.0108	0.0100	8.0	50.0
Methyl acetate	Ave	0.2050	0.2103		0.0205	0.0200	2.6	50.0
Methylene Chloride	Lin1		0.3143		0.0102	0.0100	1.8	50.0
2-Methyl-2-propanol	Ave	0.0203	0.0193		0.0952	0.100	-4.8	50.0
Acrylonitrile	Ave	0.0962	0.0979		0.102	0.100	1.8	50.0
Methyl tert-butyl ether	Ave	0.7834	0.8124		0.0104	0.0100	3.7	50.0
trans-1,2-Dichloroethene	Ave	0.2795	0.2813		0.0101	0.0100	0.6	50.0
Hexane	Ave	0.0586	0.0579		0.00987	0.0100	-1.3	20.0
1,1-Dichloroethane	Ave	0.4517	0.4658	0.1000	0.0103	0.0100	3.1	50.0
Vinyl acetate	Ave	0.5198	0.5720		0.0110	0.0100	10.0	50.0
2,2-Dichloropropane	Ave	0.0624	0.0692		0.0111	0.0100	10.9	50.0
2-Butanone (MEK)	Ave	0.0333	0.0338		0.0203	0.0200	1.7	50.0
cis-1,2-Dichloroethene	Ave	0.3078	0.3026		0.00983	0.0100	-1.7	50.0
Chlorobromomethane	Ave	0.1531	0.1576		0.0103	0.0100	2.9	50.0
Tetrahydrofuran	Ave	0.0828	0.0821		0.0198	0.0200	-0.8	50.0
Chloroform	Ave	0.4606	0.4715		0.0102	0.0100	2.4	20.0
1,1,1-Trichloroethane	Ave	0.3546	0.3844		0.0108	0.0100	8.4	50.0
Cyclohexane	Ave	0.3993	0.4228		0.0106	0.0100	5.9	50.0
1,1-Dichloropropene	Ave	0.3748	0.3803		0.0101	0.0100	1.5	50.0
Carbon tetrachloride	Ave	0.3441	0.3608		0.0105	0.0100	4.9	50.0
Isobutyl alcohol	Ave	0.0077	0.0075		0.244	0.250	-2.5	50.0
Benzene	Ave	1.116	1.120		0.0100	0.0100	0.4	50.0
1,2-Dichloroethane	Ave	0.3820	0.3923		0.0103	0.0100	2.7	50.0
n-Heptane	Ave	0.0532	0.0516		0.00969	0.0100	-3.1	50.0
Trichloroethene	Ave	0.3240	0.3277		0.0101	0.0100	1.2	50.0

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: ICV 240-364785/14 Calibration Date: 01/21/2019 13:48
 Instrument ID: A3UX15 Calib Start Date: 01/21/2019 11:37
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 01/21/2019 13:26
 Lab File ID: UXC8416.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Methylcyclohexane	Ave	0.3524	0.3635		0.0103	0.0100	3.1	50.0
1,2-Dichloropropane	Ave	0.2590	0.2693		0.0104	0.0100	4.0	20.0
1,4-Dioxane	Lin1		0.0018		0.174	0.200	-13.1	50.0
Dibromomethane	Ave	0.1717	0.1693		0.00986	0.0100	-1.4	50.0
Dichlorobromomethane	Ave	0.3410	0.3425		0.0100	0.0100	0.5	50.0
2-Chloroethyl vinyl ether	Ave	0.1745	0.1743		0.00998	0.0100	-0.2	50.0
cis-1,3-Dichloropropene	Ave	0.3944	0.4210		0.0107	0.0100	6.7	50.0
4-Methyl-2-pentanone (MIBK)	Ave	0.2678	0.2819		0.0211	0.0200	5.3	50.0
Toluene	Ave	1.476	1.520		0.0103	0.0100	3.0	20.0
trans-1,3-Dichloropropene	Ave	0.4646	0.4706		0.0101	0.0100	1.3	50.0
Ethyl methacrylate	Ave	0.4442	0.4407		0.00992	0.0100	-0.8	50.0
1,1,2-Trichloroethane	Ave	0.2982	0.3097		0.0104	0.0100	3.9	50.0
Tetrachloroethene	Ave	0.3268	0.3359		0.0103	0.0100	2.8	50.0
1,3-Dichloropropane	Ave	0.5243	0.5353		0.0102	0.0100	2.1	50.0
2-Hexanone	Ave	0.2283	0.2452		0.0215	0.0200	7.4	50.0
Chlorodibromomethane	Ave	0.3116	0.3163		0.0102	0.0100	1.5	50.0
Ethylene Dibromide	Ave	0.3191	0.3265		0.0102	0.0100	2.3	50.0
Chlorobenzene	Ave	0.9777	0.9870	0.3000	0.0101	0.0100	1.0	50.0
1,1,1,2-Tetrachloroethane	Ave	0.3450	0.3654		0.0106	0.0100	5.9	50.0
Ethylbenzene	Ave	0.5145	0.5244		0.0102	0.0100	1.9	20.0
m-Xylene & p-Xylene	Ave	1.221	1.254		0.0103	0.0100	2.7	50.0
o-Xylene	Ave	0.6189	0.6401		0.0103	0.0100	3.4	50.0
Styrene	Ave	1.036	1.018		0.00982	0.0100	-1.8	50.0
Bromoform	Lin1		0.1992	0.1000	0.00846	0.0100	-15.4	50.0
Isopropylbenzene	Ave	1.463	1.540		0.0105	0.0100	5.3	50.0
1,1,2,2-Tetrachloroethane	Ave	0.7086	0.7292	0.3000	0.0103	0.0100	2.9	50.0
Bromobenzene	Ave	0.8274	0.8332		0.0101	0.0100	0.7	50.0
1,2,3-Trichloropropane	Ave	0.2468	0.2591		0.0105	0.0100	5.0	50.0
trans-1,4-Dichloro-2-butene	Ave	0.1014	0.0704		0.00695	0.0100	-30.5	50.0
N-Propylbenzene	Ave	0.7286	0.7393		0.0101	0.0100	1.5	50.0
2-Chlorotoluene	Ave	0.6821	0.6987		0.0102	0.0100	2.4	50.0
1,3,5-Trimethylbenzene	Ave	2.119	2.131		0.0101	0.0100	0.6	50.0
4-Chlorotoluene	Ave	2.205	2.165		0.00982	0.0100	-1.8	50.0
tert-Butylbenzene	Ave	1.897	1.872		0.00987	0.0100	-1.3	50.0
1,2,4-Trimethylbenzene	Ave	2.227	2.255		0.0101	0.0100	1.2	50.0
sec-Butylbenzene	Ave	2.327	2.325		0.00999	0.0100	-0.0	50.0
1,3-Dichlorobenzene	Ave	1.425	1.437		0.0101	0.0100	0.9	50.0
4-Isopropyltoluene	Ave	2.083	2.090		0.0100	0.0100	0.3	50.0
1,4-Dichlorobenzene	Ave	1.477	1.484		0.0100	0.0100	0.4	50.0
n-Butylbenzene	Ave	1.649	1.603		0.00972	0.0100	-2.8	50.0
1,2-Dichlorobenzene	Ave	1.387	1.408		0.0102	0.0100	1.5	50.0

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: ICV 240-364785/14 Calibration Date: 01/21/2019 13:48
 Instrument ID: A3UX15 Calib Start Date: 01/21/2019 11:37
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 01/21/2019 13:26
 Lab File ID: UXC8416.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,2-Dibromo-3-Chloropropane	Ave	0.1611	0.1660		0.0103	0.0100	3.0	50.0
1,2,4-Trichlorobenzene	Ave	0.8784	0.8494		0.00967	0.0100	-3.3	50.0
Hexachlorobutadiene	Qua		0.3722		0.00828	0.0100	-17.2	50.0
Naphthalene	Ave	2.209	2.304		0.0104	0.0100	4.3	50.0
1,2,3-Trichlorobenzene	Ave	0.7902	0.7849		0.00993	0.0100	-0.7	50.0
Dibromofluoromethane (Surr)	Ave	0.2366	0.2183		0.00735	0.00797	-7.7	50.0
1,2-Dichloroethane-d4 (Surr)	Ave	0.3042	0.2714		0.00711	0.00797	-10.8	50.0
Toluene-d8 (Surr)	Ave	1.214	1.116		0.00733	0.00797	-8.1	50.0
4-Bromofluorobenzene (Surr)	Ave	0.4341	0.4000		0.00734	0.00797	-7.9	50.0

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8416.D
 Lims ID: ICV
 Client ID:
 Sample Type: ICV
 Inject. Date: 21-Jan-2019 13:48:30 ALS Bottle#: 7 Worklist Smp#: 14
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0083824-014
 Misc. Info.: C90121A-IC,8260LLUX15,,43582
 Operator ID: Instrument ID: A3UX15
 Sublist:
 Method: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\8260_15.m
 Limit Group: MSV 8260B ICAL
 Last Update: 22-Jan-2019 09:10:13 Calib Date: 21-Jan-2019 19:15:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8431.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0303

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	5.196	5.196	0.000	99	1793251	10.0	10.0	
* 2 Chlorobenzene-d5	117	7.876	7.876	0.000	85	1418410	10.0	10.0	
* 3 1,4-Dichlorobenzene-d4	152	10.117	10.117	0.000	93	796911	10.0	10.0	
\$ 4 Dibromofluoromethane (Surr	113	4.639	4.639	0.000	95	311994	7.97	7.35	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.923	4.923	0.000	100	387861	7.97	7.11	
\$ 6 Toluene-d8 (Surr)	98	6.560	6.560	0.000	93	1261093	7.97	7.33	
\$ 7 4-Bromofluorobenzene (Surr	95	8.979	8.979	0.000	96	452132	7.97	7.34	
9 Dichlorodifluoromethane	85	1.532	1.532	0.000	99	551850	10.0	10.0	
10 Chloromethane	50	1.686	1.686	0.000	99	527044	10.0	9.37	
11 Vinyl chloride	62	1.781	1.781	0.000	97	531634	10.0	10.3	
12 Butadiene	54	1.793	1.793	-0.001	96	616742	10.0	9.06	
13 Bromomethane	94	2.053	2.054	-0.001	91	391096	10.0	9.45	
14 Chloroethane	64	2.136	2.137	-0.001	100	347264	10.0	9.49	
15 Dichlorofluoromethane	67	2.302	2.303	-0.001	97	903414	10.0	9.87	
16 Trichlorofluoromethane	101	2.362	2.362	0.000	99	834671	10.0	10.7	
17 Ethyl ether	59	2.575	2.563	0.012	90	448075	10.0	10.8	
18 Acrolein	56	2.694	2.694	0.000	99	201389	50.0	35.4	
19 1,1-Dichloroethene	96	2.789	2.777	0.012	98	529054	10.0	9.82	
20 1,1,2-Trichloro-1,2,2-trif	151	2.800	2.789	0.011	86	478556	10.0	10.8	
22 Acetone	58	2.812	2.812	0.000	100	93357	20.0	21.7	
23 Iodomethane	142	2.931	2.931	0.000	98	912367	10.0	10.9	
24 Carbon disulfide	76	2.990	2.990	0.000	100	1404467	10.0	10.1	
25 3-Chloro-1-propene	76	3.073	3.073	0.000	88	329005	10.0	10.8	
27 Methyl acetate	43	3.085	3.085	0.000	98	754191	20.0	20.5	
28 Methylene Chloride	84	3.180	3.180	0.000	93	563557	10.0	10.2	
29 2-Methyl-2-propanol	59	3.263	3.263	0.000	99	346305	100.0	95.2	
32 Acrylonitrile	53	3.393	3.394	-0.001	99	1755477	100.0	101.8	
30 trans-1,2-Dichloroethene	96	3.405	3.405	0.000	97	504418	10.0	10.1	
31 Methyl tert-butyl ether	73	3.405	3.405	0.000	95	1456881	10.0	10.4	
33 Hexane	86	3.619	3.619	0.000	94	103809	10.0	9.87	
34 1,1-Dichloroethane	63	3.761	3.761	0.000	96	835310	10.0	10.3	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
36 Vinyl acetate	43	3.785	3.785	0.000	97	1025753	10.0	11.0	
39 cis-1,2-Dichloroethene	96	4.247	4.235	0.012	81	542596	10.0	9.83	
41 2-Butanone (MEK)	72	4.247	4.247	0.000	98	121331	20.0	20.3	
40 2,2-Dichloropropane	97	4.247	4.247	0.000	65	124051	10.0	11.1	
45 Chlorobromomethane	128	4.449	4.449	0.000	88	282681	10.0	10.3	
46 Tetrahydrofuran	42	4.484	4.485	-0.001	86	294586	20.0	19.8	
47 Chloroform	83	4.496	4.496	0.000	93	845575	10.0	10.2	
48 1,1,1-Trichloroethane	97	4.662	4.662	0.000	98	689257	10.0	10.8	
49 Cyclohexane	56	4.710	4.710	0.000	90	758127	10.0	10.6	
51 1,1-Dichloropropene	75	4.793	4.793	0.000	94	681963	10.0	10.1	
50 Carbon tetrachloride	117	4.805	4.805	0.000	96	647000	10.0	10.5	
52 Isobutyl alcohol	41	4.864	4.864	0.000	95	335938	250.0	243.9	
53 Benzene	78	4.971	4.971	0.000	96	2009249	10.0	10.0	
54 1,2-Dichloroethane	62	4.982	4.983	-0.001	98	703449	10.0	10.3	
56 n-Heptane	100	5.160	5.160	0.000	93	92495	10.0	9.69	
58 Trichloroethene	130	5.504	5.504	0.000	95	587677	10.0	10.1	
60 Methylcyclohexane	83	5.670	5.670	0.000	90	651902	10.0	10.3	
61 1,2-Dichloropropane	63	5.706	5.706	0.000	94	482855	10.0	10.4	
63 Dibromomethane	93	5.813	5.813	0.000	88	303520	10.0	9.86	
64 1,4-Dioxane	88	5.813	5.813	0.000	33	65935	200.0	173.7	
65 Dichlorobromomethane	83	5.931	5.931	0.000	99	614221	10.0	10.0	
67 2-Chloroethyl vinyl ether	63	6.180	6.180	0.000	93	312475	10.0	9.98	
68 cis-1,3-Dichloropropene	75	6.322	6.323	-0.001	95	754933	10.0	10.7	
69 4-Methyl-2-pentanone (MIBK)	43	6.441	6.441	0.000	97	1011196	20.0	21.1	
70 Toluene	91	6.619	6.619	0.000	98	2155730	10.0	10.3	
71 trans-1,3-Dichloropropene	75	6.809	6.809	0.000	94	667505	10.0	10.1	
72 Ethyl methacrylate	69	6.868	6.868	0.000	89	625023	10.0	9.92	
73 1,1,2-Trichloroethane	97	6.975	6.975	0.000	91	439234	10.0	10.4	
74 Tetrachloroethene	164	7.117	7.117	0.000	96	476419	10.0	10.3	
75 1,3-Dichloropropane	76	7.129	7.129	0.000	93	759222	10.0	10.2	
76 2-Hexanone	43	7.188	7.188	0.000	97	695520	20.0	21.5	
79 Chlorodibromomethane	129	7.342	7.342	0.000	89	448610	10.0	10.2	
80 Ethylene Dibromide	107	7.461	7.461	0.000	98	463121	10.0	10.2	
82 Chlorobenzene	112	7.900	7.900	0.000	95	1399979	10.0	10.1	
83 1,1,1,2-Tetrachloroethane	131	7.971	7.971	0.000	94	518307	10.0	10.6	
84 Ethylbenzene	106	7.995	7.995	-0.001	98	743771	10.0	10.2	
85 m-Xylene & p-Xylene	91	8.101	8.101	0.000	93	1778779	10.0	10.3	
86 o-Xylene	106	8.481	8.481	0.000	95	907873	10.0	10.3	
87 Styrene	104	8.493	8.493	0.000	94	1443373	10.0	9.82	
88 Bromoform	173	8.682	8.682	0.000	98	282491	10.0	8.46	
89 Isopropylbenzene	105	8.825	8.825	0.000	95	2184869	10.0	10.5	
93 1,1,2,2-Tetrachloroethane	83	9.121	9.121	0.000	97	581082	10.0	10.3	
92 Bromobenzene	156	9.133	9.133	0.000	91	663996	10.0	10.1	
95 trans-1,4-Dichloro-2-buten	53	9.168	9.169	-0.001	67	56124	10.0	6.95	
94 1,2,3-Trichloropropane	110	9.168	9.169	-0.001	84	206463	10.0	10.5	
96 N-Propylbenzene	120	9.228	9.228	0.000	98	589187	10.0	10.1	
97 2-Chlorotoluene	126	9.323	9.323	0.000	97	556828	10.0	10.2	
98 1,3,5-Trimethylbenzene	105	9.394	9.394	0.000	94	1698481	10.0	10.1	
99 4-Chlorotoluene	91	9.418	9.418	0.000	98	1725613	10.0	9.82	
101 tert-Butylbenzene	119	9.714	9.714	0.000	92	1491725	10.0	9.87	
103 1,2,4-Trimethylbenzene	105	9.761	9.762	-0.001	94	1796823	10.0	10.1	
105 sec-Butylbenzene	105	9.927	9.928	-0.001	94	1852982	10.0	10.0	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
106 1,3-Dichlorobenzene	146	10.046	10.046	0.000	98	1145305	10.0	10.1	
107 4-Isopropyltoluene	119	10.070	10.070	0.000	97	1665303	10.0	10.0	
108 1,4-Dichlorobenzene	146	10.141	10.141	0.000	96	1182487	10.0	10.0	
111 n-Butylbenzene	91	10.473	10.473	0.000	97	1277066	10.0	9.72	
112 1,2-Dichlorobenzene	146	10.509	10.509	-0.001	99	1122403	10.0	10.2	
113 1,2-Dibromo-3-Chloropropan	157	11.291	11.291	0.000	89	132296	10.0	10.3	
115 1,2,4-Trichlorobenzene	180	12.109	12.110	-0.001	94	676924	10.0	9.67	
116 Hexachlorobutadiene	225	12.275	12.276	-0.001	93	296600	10.0	8.28	
117 Naphthalene	128	12.358	12.359	-0.001	96	1836241	10.0	10.4	
118 1,2,3-Trichlorobenzene	180	12.607	12.608	-0.001	95	625482	10.0	9.93	
S 129 Xylenes, Total	106				0		20.0	20.6	
S 157 Total BTEX	1		0.000				50.0	ND	
S 130 Trihalomethanes, Total	1				0		40.0	38.9	

Reagents:

VMFASPW_00285	Amount Added: 8.00	Units: uL	
VMFASAW_00264	Amount Added: 8.00	Units: uL	
VMFASGW_00293	Amount Added: 8.00	Units: uL	
VM50IS_00072	Amount Added: 1.00	Units: uL	
vmDist_H2o_00124	Amount Added: 0.00	Units:	Run Reagent
vm50ss_stk_00079	Amount Added: 0.80	Units: uL	Run Reagent

Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8416.D

Injection Date: 21-Jan-2019 13:48:30

Instrument ID: A3UX15

Operator ID:

Lims ID: ICV

Worklist Smp#: 14

Client ID:

Purge Vol: 5.000 mL

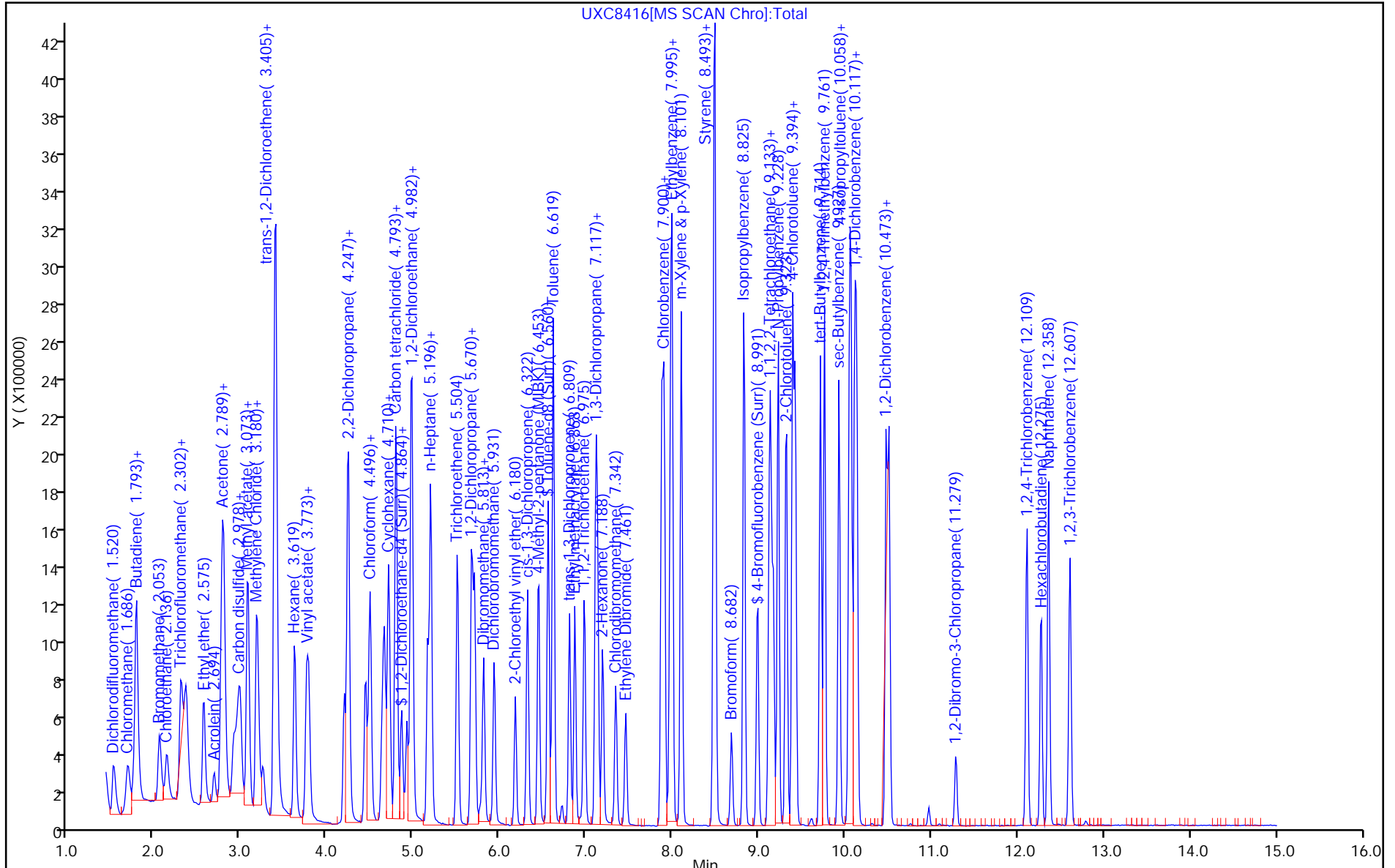
Dil. Factor: 1.0000

ALS Bottle#: 7

Method: 8260_15

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: ICV 240-364785/15 Calibration Date: 01/21/2019 16:21
 Instrument ID: A3UX15 Calib Start Date: 01/21/2019 11:37
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 01/21/2019 13:26
 Lab File ID: UXC8423.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Dibromofluoromethane (Surr)	Ave	0.2366	0.2218		0.00747	0.00797	-6.3	50.0
1,2-Dichloroethane-d4 (Surr)	Ave	0.3042	0.2777		0.00728	0.00797	-8.7	50.0
Toluene-d8 (Surr)	Ave	1.214	1.132		0.00743	0.00797	-6.7	50.0
4-Bromofluorobenzene (Surr)	Ave	0.4341	0.4058		0.00745	0.00797	-6.5	50.0

TestAmerica Canton
Target Compound Quantitation Report

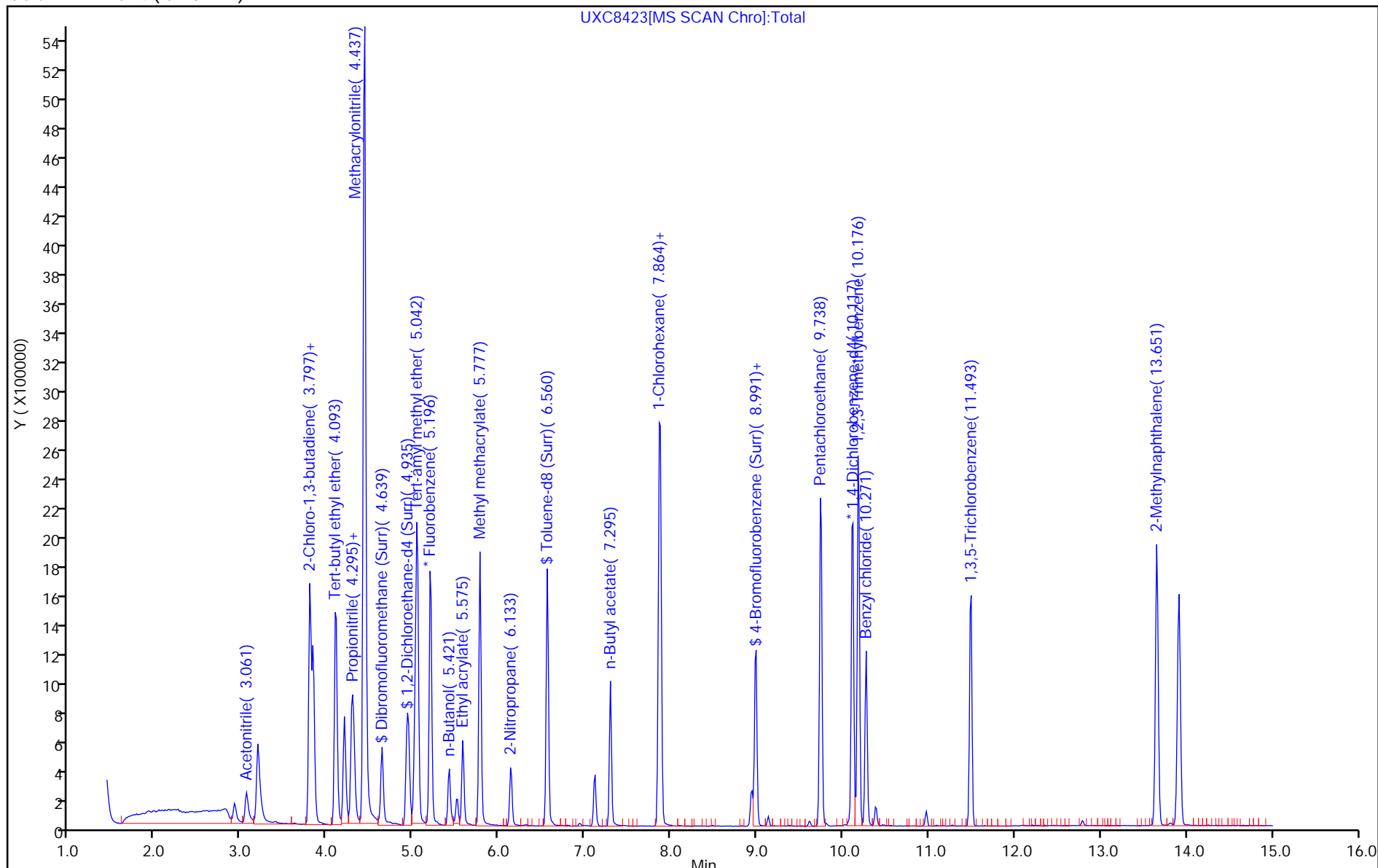
Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8423.D
 Lims ID: ICV A9
 Client ID:
 Sample Type: ICV
 Inject. Date: 21-Jan-2019 16:21:30 ALS Bottle#: 14 Worklist Smp#: 15
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0083824-015
 Misc. Info.: C90121A-IC,8260LLUX15,,43582
 Operator ID: Instrument ID: A3UX15
 Sublist:
 Method: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\8260_15.m
 Limit Group: MSV 8260B ICAL
 Last Update: 22-Jan-2019 09:10:13 Calib Date: 21-Jan-2019 19:15:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8431.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0303

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	5.196	5.196	0.000	99	1778453	10.0	10.0	
* 2 Chlorobenzene-d5	117	7.876	7.876	0.000	85	1420097	10.0	10.0	
* 3 1,4-Dichlorobenzene-d4	152	10.117	10.117	0.000	93	764446	10.0	10.0	
\$ 4 Dibromofluoromethane (Surr	113	4.639	4.639	0.000	95	314414	7.97	7.47	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.923	4.923	0.000	100	393596	7.97	7.28	
\$ 6 Toluene-d8 (Surr)	98	6.560	6.560	0.000	93	1281042	7.97	7.43	
\$ 7 4-Bromofluorobenzene (Surr	95	8.979	8.979	0.000	97	459309	7.97	7.45	
26 Acetonitrile	41	3.061	3.050	0.011	98	427448	100.0	104.0	
35 Isopropyl ether	87	3.797	3.797	0.000	92	442161	10.0	10.3	
37 2-Chloro-1,3-butadiene	53	3.832	3.832	0.000	93	736814	10.0	9.78	
38 Tert-butyl ethyl ether	59	4.093	4.093	0.000	97	1388893	10.0	10.0	
42 Ethyl acetate	43	4.283	4.283	0.000	99	937785	20.0	20.8	
43 Propionitrile	54	4.307	4.307	0.000	99	677621	100.0	106.8	
44 Methacrylonitrile	41	4.437	4.425	0.012	92	3212963	100.0	104.3	
55 Tert-amyl methyl ether	73	5.054	5.042	0.012	98	1404415	10.0	10.2	
57 n-Butanol	56	5.421	5.421	0.000	89	269615	250.0	249.7	
59 Ethyl acrylate	55	5.575	5.575	0.000	99	624490	10.0	10.5	
62 Methyl methacrylate	41	5.777	5.777	0.000	91	913802	20.0	21.1	
66 2-Nitropropane	41	6.133	6.133	0.000	99	271968	20.0	20.5	
77 n-Butyl acetate	43	7.295	7.295	0.000	97	736671	10.0	10.6	
81 1-Chlorohexane	91	7.864	7.864	0.000	98	551389	10.0	9.79	
91 Cyclohexanone	55	8.943	8.931	0.012	94	139432	100.0	114.7	
102 Pentachloroethane	167	9.750	9.750	0.000	95	563326	20.0	19.5	
109 1,2,3-Trimethylbenzene	105	10.176	10.176	0.000	98	1800683	10.0	10.3	
110 Benzyl chloride	126	10.271	10.271	0.000	98	227669	10.0	10.1	
114 1,3,5-Trichlorobenzene	180	11.493	11.493	0.000	97	692380	10.0	9.57	
119 2-Methylnaphthalene	142	13.651	13.651	0.000	93	1419962	20.0	17.3	

Reagents:

VMFASA9W_00216	Amount Added: 8.00	Units: uL	
VM50IS_00072	Amount Added: 1.00	Units: uL	
vmDist_H2o_00124	Amount Added: 0.00	Units:	Run Reagent
vm50ss_stk_00079	Amount Added: 0.80	Units: uL	Run Reagent

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FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: ICV 240-364785/15 Calibration Date: 01/21/2019 16:21
 Instrument ID: A3UX15 Calib Start Date: 01/21/2019 14:10
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 01/21/2019 15:59
 Lab File ID: UXC8423.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Acetonitrile	Ave	0.0231	0.0240		0.104	0.100	4.0	50.0
Isopropyl ether	Ave	0.2405	0.2486		0.0103	0.0100	3.4	50.0
2-Chloro-1,3-butadiene	Ave	0.4235	0.4143		0.00978	0.0100	-2.2	50.0
Tert-butyl ethyl ether	Ave	0.7799	0.7810		0.0100	0.0100	0.1	50.0
Ethyl acetate	Ave	0.2538	0.2637		0.0208	0.0200	3.9	50.0
Propionitrile	Ave	0.0357	0.0381		0.107	0.100	6.8	50.0
Methacrylonitrile	Ave	0.1733	0.1807		0.104	0.100	4.3	50.0
Tert-amyl methyl ether	Ave	0.7765	0.7897		0.0102	0.0100	1.7	50.0
n-Butanol	Ave	0.0061	0.0061		0.250	0.250	-0.1	50.0
Ethyl acrylate	Ave	0.3330	0.3511		0.0105	0.0100	5.4	50.0
Methyl methacrylate	Ave	0.2430	0.2569		0.0211	0.0200	5.7	50.0
2-Nitropropane	Ave	0.0745	0.0765		0.0205	0.0200	2.7	50.0
n-Butyl acetate	Ave	0.3918	0.4142		0.0106	0.0100	5.7	50.0
1-Chlorohexane	Ave	0.3968	0.3883		0.00979	0.0100	-2.1	50.0
Cyclohexanone	Lin1		0.0182		0.115	0.100	14.7	50.0
Pentachloroethane	Ave	0.2034	0.1983		0.0195	0.0200	-2.5	50.0
1,2,3-Trimethylbenzene	Ave	2.297	2.356		0.0103	0.0100	2.5	50.0
Benzyl chloride	Ave	0.2963	0.2978		0.0101	0.0100	0.5	50.0
1,3,5-Trichlorobenzene	Ave	0.9466	0.9057		0.00957	0.0100	-4.3	50.0
2-Methylnaphthalene	Ave	1.076	0.9288		0.0173	0.0200	-13.7	50.0

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8423.D
 Lims ID: ICV A9
 Client ID:
 Sample Type: ICV
 Inject. Date: 21-Jan-2019 16:21:30 ALS Bottle#: 14 Worklist Smp#: 15
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0083824-015
 Misc. Info.: C90121A-IC,8260LLUX15,,43582
 Operator ID: Instrument ID: A3UX15
 Sublist:
 Method: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\8260_15.m
 Limit Group: MSV 8260B ICAL
 Last Update: 22-Jan-2019 09:10:13 Calib Date: 21-Jan-2019 19:15:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8431.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0303

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	5.196	5.196	0.000	99	1778453	10.0	10.0	
* 2 Chlorobenzene-d5	117	7.876	7.876	0.000	85	1420097	10.0	10.0	
* 3 1,4-Dichlorobenzene-d4	152	10.117	10.117	0.000	93	764446	10.0	10.0	
\$ 4 Dibromofluoromethane (Surr	113	4.639	4.639	0.000	95	314414	7.97	7.47	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.923	4.923	0.000	100	393596	7.97	7.28	
\$ 6 Toluene-d8 (Surr)	98	6.560	6.560	0.000	93	1281042	7.97	7.43	
\$ 7 4-Bromofluorobenzene (Surr	95	8.979	8.979	0.000	97	459309	7.97	7.45	
26 Acetonitrile	41	3.061	3.050	0.011	98	427448	100.0	104.0	
35 Isopropyl ether	87	3.797	3.797	0.000	92	442161	10.0	10.3	
37 2-Chloro-1,3-butadiene	53	3.832	3.832	0.000	93	736814	10.0	9.78	
38 Tert-butyl ethyl ether	59	4.093	4.093	0.000	97	1388893	10.0	10.0	
42 Ethyl acetate	43	4.283	4.283	0.000	99	937785	20.0	20.8	
43 Propionitrile	54	4.307	4.307	0.000	99	677621	100.0	106.8	
44 Methacrylonitrile	41	4.437	4.425	0.012	92	3212963	100.0	104.3	
55 Tert-amyl methyl ether	73	5.054	5.042	0.012	98	1404415	10.0	10.2	
57 n-Butanol	56	5.421	5.421	0.000	89	269615	250.0	249.7	
59 Ethyl acrylate	55	5.575	5.575	0.000	99	624490	10.0	10.5	
62 Methyl methacrylate	41	5.777	5.777	0.000	91	913802	20.0	21.1	
66 2-Nitropropane	41	6.133	6.133	0.000	99	271968	20.0	20.5	
77 n-Butyl acetate	43	7.295	7.295	0.000	97	736671	10.0	10.6	
81 1-Chlorohexane	91	7.864	7.864	0.000	98	551389	10.0	9.79	
91 Cyclohexanone	55	8.943	8.931	0.012	94	139432	100.0	114.7	
102 Pentachloroethane	167	9.750	9.750	0.000	95	563326	20.0	19.5	
109 1,2,3-Trimethylbenzene	105	10.176	10.176	0.000	98	1800683	10.0	10.3	
110 Benzyl chloride	126	10.271	10.271	0.000	98	227669	10.0	10.1	
114 1,3,5-Trichlorobenzene	180	11.493	11.493	0.000	97	692380	10.0	9.57	
119 2-Methylnaphthalene	142	13.651	13.651	0.000	93	1419962	20.0	17.3	

Reagents:

VMFASA9W_00216	Amount Added: 8.00	Units: uL	
VM50IS_00072	Amount Added: 1.00	Units: uL	
vmDist_H2o_00124	Amount Added: 0.00	Units:	Run Reagent
vm50ss_stk_00079	Amount Added: 0.80	Units: uL	Run Reagent

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Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8423.D

Injection Date: 21-Jan-2019 16:21:30

Instrument ID: A3UX15

Operator ID:

Lims ID: ICV A9

Worklist Smp#: 15

Client ID:

Purge Vol: 5.000 mL

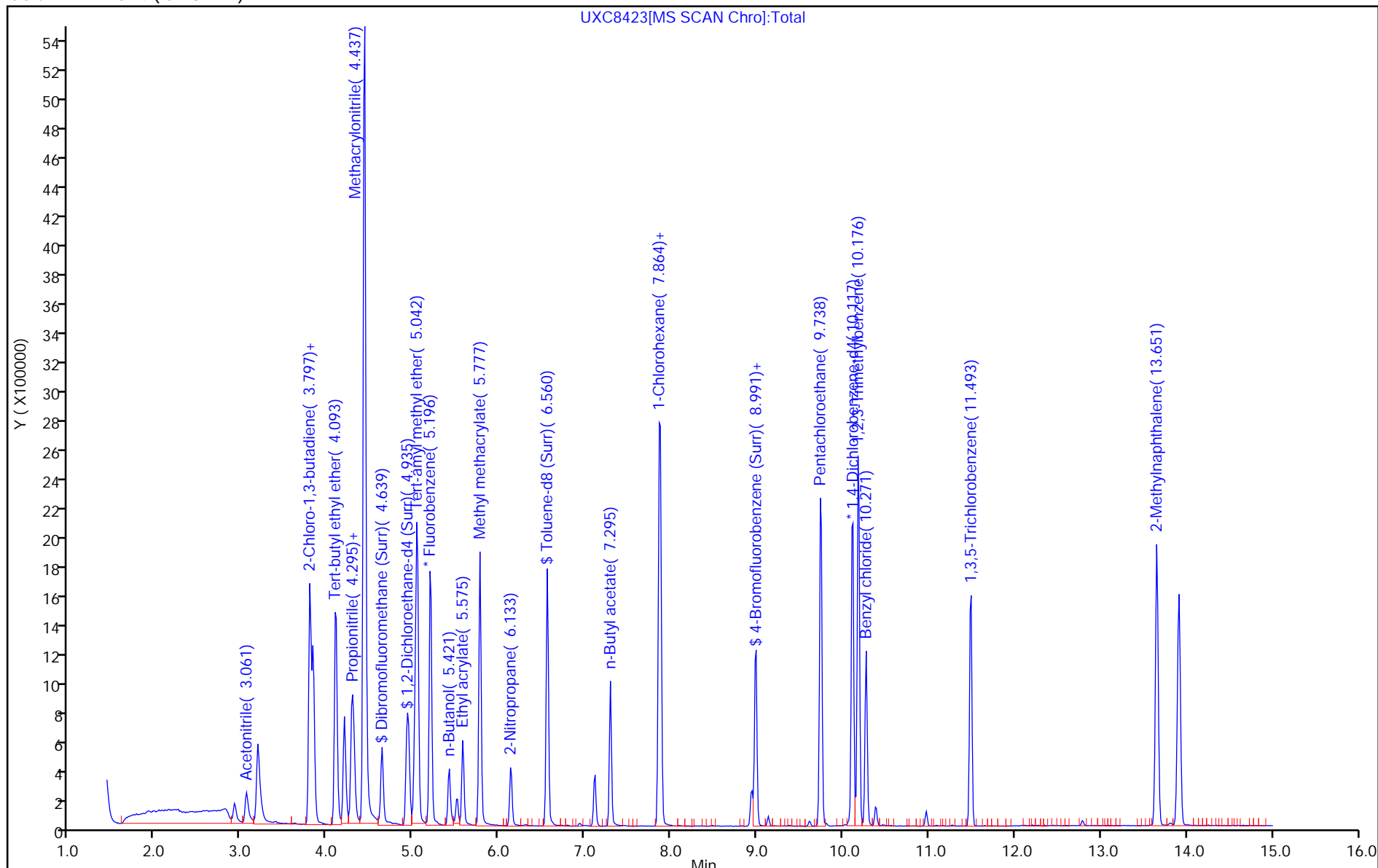
Dil. Factor: 1.0000

ALS Bottle#: 14

Method: 8260_15

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: CCVIS 240-372032/2 Calibration Date: 03/18/2019 10:59
 Instrument ID: A3UX15 Calib Start Date: 01/21/2019 11:37
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 01/21/2019 13:26
 Lab File ID: UXC9421.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Dichlorodifluoromethane	Ave	0.3072	0.2534		0.00825	0.0100	-17.5	50.0
Chloromethane	Ave	0.3135	0.2524	0.1000	0.00805	0.0100	-19.5	50.0
Vinyl chloride	Ave	0.2885	0.2492		0.00864	0.0100	-13.6	20.0
Butadiene	Ave	0.3794	0.2713		0.00715	0.0100	-28.5	50.0
Bromomethane	Ave	0.2308	0.1738		0.00753	0.0100	-24.7	50.0
Chloroethane	Ave	0.2041	0.1605		0.00786	0.0100	-21.4	50.0
Dichlorofluoromethane	Ave	0.5102	0.3989		0.00782	0.0100	-21.8	50.0
Trichlorofluoromethane	Ave	0.4337	0.4178		0.00963	0.0100	-3.7	50.0
Ethyl ether	Ave	0.2310	0.1949		0.00844	0.0100	-15.6	50.0
Acrolein	Ave	0.0317	0.0214		0.0337	0.0500	-32.6	50.0
1,1-Dichloroethene	Ave	0.3005	0.2609		0.00868	0.0100	-13.2	20.0
1,1,2-Trichloro-1,2,2-trifluoroethane	Ave	0.2468	0.2382		0.00965	0.0100	-3.5	50.0
Acetone	Lin1		0.0194		0.0156	0.0200	-22.2	50.0
Iodomethane	Ave	0.4675	0.4806		0.0103	0.0100	2.8	50.0
Carbon disulfide	Ave	0.7741	0.6312		0.00815	0.0100	-18.5	50.0
3-Chloro-1-propene	Ave	0.1698	0.1491		0.00878	0.0100	-12.2	50.0
Methyl acetate	Ave	0.2050	0.1377		0.0134	0.0200	-32.8	50.0
Methylene Chloride	Lin1		0.2934		0.00934	0.0100	-6.6	50.0
2-Methyl-2-propanol	Ave	0.0203	0.0118		0.0581	0.100	-41.9	50.0
Acrylonitrile	Ave	0.0962	0.0719		0.0747	0.100	-25.3	50.0
Methyl tert-butyl ether	Ave	0.7834	0.5704		0.00728	0.0100	-27.2	50.0
trans-1,2-Dichloroethene	Ave	0.2795	0.2935		0.0105	0.0100	5.0	50.0
Hexane	Ave	0.0586	0.0529		0.00902	0.0100	-9.8	20.0
1,1-Dichloroethane	Ave	0.4517	0.4870	0.1000	0.0108	0.0100	7.8	50.0
Vinyl acetate	Ave	0.5198	0.3392		0.00652	0.0100	-34.8	50.0
cis-1,2-Dichloroethene	Ave	0.3078	0.3347		0.0109	0.0100	8.7	50.0
2,2-Dichloropropane	Ave	0.0624	0.0711		0.0114	0.0100	13.9	50.0
2-Butanone (MEK)	Ave	0.0333	0.0218		0.0131	0.0200	-34.4	50.0
Chlorobromomethane	Ave	0.1531	0.1589		0.0104	0.0100	3.7	50.0
Tetrahydrofuran	Ave	0.0828	0.0496		0.0120	0.0200	-40.0	50.0
Chloroform	Ave	0.4606	0.5093		0.0111	0.0100	10.6	20.0
1,1,1-Trichloroethane	Ave	0.3546	0.4147		0.0117	0.0100	16.9	50.0
Cyclohexane	Ave	0.3993	0.3878		0.00971	0.0100	-2.9	50.0
1,1-Dichloropropene	Ave	0.3748	0.3738		0.00998	0.0100	-0.2	50.0
Carbon tetrachloride	Ave	0.3441	0.3694		0.0107	0.0100	7.4	50.0
Isobutyl alcohol	Ave	0.0077	0.0052		0.168	0.250	-32.6	50.0
Benzene	Ave	1.116	1.191		0.0107	0.0100	6.6	50.0
1,2-Dichloroethane	Ave	0.3820	0.3804		0.00996	0.0100	-0.4	50.0
n-Heptane	Ave	0.0532	0.0462		0.00869	0.0100	-13.1	50.0
Trichloroethene	Ave	0.3240	0.3329		0.0103	0.0100	2.7	50.0

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: CCVIS 240-372032/2 Calibration Date: 03/18/2019 10:59
 Instrument ID: A3UX15 Calib Start Date: 01/21/2019 11:37
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 01/21/2019 13:26
 Lab File ID: UXC9421.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Methylcyclohexane	Ave	0.3524	0.3304		0.00937	0.0100	-6.3	50.0
1,2-Dichloropropane	Ave	0.2590	0.2571		0.00993	0.0100	-0.7	20.0
1,4-Dioxane	Lin1		0.0019		0.181	0.200	-9.6	50.0
Dibromomethane	Ave	0.1717	0.1543		0.00898	0.0100	-10.2	50.0
Dichlorobromomethane	Ave	0.3410	0.3179		0.00932	0.0100	-6.8	50.0
2-Chloroethyl vinyl ether	Ave	0.1745	0.1036		0.0119	0.0200	-40.7	50.0
cis-1,3-Dichloropropene	Ave	0.3944	0.3153		0.00799	0.0100	-20.1	50.0
4-Methyl-2-pentanone (MIBK)	Ave	0.2678	0.1583		0.0118	0.0200	-40.9	50.0
Toluene	Ave	1.476	1.539		0.0104	0.0100	4.2	20.0
trans-1,3-Dichloropropene	Ave	0.4646	0.3300		0.00710	0.0100	-29.0	50.0
Ethyl methacrylate	Ave	0.4442	0.2904		0.00654	0.0100	-34.6	50.0
1,1,2-Trichloroethane	Ave	0.2982	0.2704		0.00907	0.0100	-9.3	50.0
Tetrachloroethene	Ave	0.3268	0.3644		0.0112	0.0100	11.5	50.0
1,3-Dichloropropane	Ave	0.5243	0.4702		0.00897	0.0100	-10.3	50.0
2-Hexanone	Ave	0.2283	0.1376		0.0121	0.0200	-39.7	50.0
Chlorodibromomethane	Ave	0.3116	0.2689		0.00863	0.0100	-13.7	50.0
Ethylene Dibromide	Ave	0.3191	0.2575		0.00807	0.0100	-19.3	50.0
Chlorobenzene	Ave	0.9777	1.014	0.3000	0.0104	0.0100	3.7	50.0
1,1,1,2-Tetrachloroethane	Ave	0.3450	0.3543		0.0103	0.0100	2.7	50.0
Ethylbenzene	Ave	0.5145	0.5150		0.0100	0.0100	0.1	20.0
m-Xylene & p-Xylene	Ave	1.221	1.230		0.0101	0.0100	0.8	50.0
o-Xylene	Ave	0.6189	0.6303		0.0102	0.0100	1.8	50.0
Styrene	Ave	1.036	1.055		0.0102	0.0100	1.9	50.0
Bromoform	Lin1		0.1439	0.1000	0.00629	0.0100	-37.1	50.0
Isopropylbenzene	Ave	1.463	1.473		0.0101	0.0100	0.7	50.0
1,1,2,2-Tetrachloroethane	Ave	0.7086	0.5014	0.3000	0.00708	0.0100	-29.2	50.0
Bromobenzene	Ave	0.8274	0.7803		0.00943	0.0100	-5.7	50.0
1,2,3-Trichloropropane	Ave	0.2468	0.1728		0.00700	0.0100	-30.0	50.0
trans-1,4-Dichloro-2-butene	Ave	0.1014	0.0667		0.00657	0.0100	-34.3	50.0
N-Propylbenzene	Ave	0.7286	0.6816		0.00935	0.0100	-6.5	50.0
2-Chlorotoluene	Ave	0.6821	0.6525		0.00957	0.0100	-4.3	50.0
1,3,5-Trimethylbenzene	Ave	2.119	2.006		0.00947	0.0100	-5.3	50.0
4-Chlorotoluene	Ave	2.205	2.125		0.00963	0.0100	-3.7	50.0
tert-Butylbenzene	Ave	1.897	1.631		0.00859	0.0100	-14.1	50.0
1,2,4-Trimethylbenzene	Ave	2.227	2.134		0.00958	0.0100	-4.2	50.0
sec-Butylbenzene	Ave	2.327	2.085		0.00896	0.0100	-10.4	50.0
1,3-Dichlorobenzene	Ave	1.425	1.355		0.00951	0.0100	-4.9	50.0
4-Isopropyltoluene	Ave	2.083	1.921		0.00922	0.0100	-7.8	50.0
1,4-Dichlorobenzene	Ave	1.477	1.407		0.00952	0.0100	-4.8	50.0
n-Butylbenzene	Ave	1.649	1.393		0.00845	0.0100	-15.5	50.0
1,2-Dichlorobenzene	Ave	1.387	1.321		0.00952	0.0100	-4.8	50.0

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Lab Sample ID: CCVIS 240-372032/2 Calibration Date: 03/18/2019 10:59
 Instrument ID: A3UX15 Calib Start Date: 01/21/2019 11:37
 GC Column: DB-624 ID: 0.18 (mm) Calib End Date: 01/21/2019 13:26
 Lab File ID: UXC9421.D Conc. Units: ng/uL Heated Purge: (Y/N) N

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,2-Dibromo-3-Chloropropane	Ave	0.1611	0.0745		0.00462	0.0100	-53.8*	50.0
1,2,4-Trichlorobenzene	Ave	0.8784	0.7191		0.00819	0.0100	-18.1	50.0
Hexachlorobutadiene	Qua		0.3827		0.00853	0.0100	-14.7	50.0
Naphthalene	Ave	2.209	1.235		0.00559	0.0100	-44.1	50.0
1,2,3-Trichlorobenzene	Ave	0.7902	0.6662		0.00843	0.0100	-15.7	50.0
Dibromofluoromethane (Surr)	Ave	0.2366	0.2045		0.00689	0.00797	-13.6	50.0
1,2-Dichloroethane-d4 (Surr)	Ave	0.3042	0.2416		0.00633	0.00797	-20.6	50.0
Toluene-d8 (Surr)	Ave	1.214	1.001		0.00657	0.00797	-17.5	50.0
4-Bromofluorobenzene (Surr)	Ave	0.4341	0.3465		0.00636	0.00797	-20.2	50.0

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX15\20190318-85248.b\UXC9421.D
 Lims ID: CCVIS L4 8260
 Client ID:
 Sample Type: CCVIS
 Inject. Date: 18-Mar-2019 10:59:30 ALS Bottle#: 1 Worklist Smp#: 2
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0085248-002
 Misc. Info.: C90318A,8260LLUX15,,43582
 Operator ID: Instrument ID: A3UX15
 Sublist: chrom-8260_15*sub83
 Method: \\chromna\Canton\ChromData\A3UX15\20190318-85248.b\8260_15.m
 Limit Group: MSV 8260B ICAL
 Last Update: 19-Mar-2019 08:39:32 Calib Date: 21-Jan-2019 19:15:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8431.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0304

First Level Reviewer: evansle

Date: 18-Mar-2019 13:40:38

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	5.196	5.196	0.000	99	2085639	10.0	10.0	
* 2 Chlorobenzene-d5	117	7.876	7.876	0.000	84	1609797	10.0	10.0	
* 3 1,4-Dichlorobenzene-d4	152	10.117	10.117	0.000	92	977106	10.0	10.0	
\$ 4 Dibromofluoromethane (Surr	113	4.639	4.639	0.000	95	339977	7.97	6.89	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.923	4.923	0.000	100	401650	7.97	6.33	
\$ 6 Toluene-d8 (Surr)	98	6.560	6.560	0.000	93	1284316	7.97	6.57	
\$ 7 4-Bromofluorobenzene (Surr	95	8.991	8.991	0.000	95	444515	7.97	6.36	
9 Dichlorodifluoromethane	85	1.520	1.520	0.000	99	528429	10.0	8.25	
10 Chloromethane	50	1.686	1.686	0.000	99	526505	10.0	8.05	
11 Vinyl chloride	62	1.769	1.769	0.000	98	519834	10.0	8.64	
12 Butadiene	54	1.781	1.781	0.000	99	565867	10.0	7.15	
13 Bromomethane	94	2.054	2.054	0.000	91	362488	10.0	7.53	
14 Chloroethane	64	2.125	2.125	0.000	99	334757	10.0	7.86	
15 Dichlorofluoromethane	67	2.303	2.303	0.000	97	832048	10.0	7.82	
16 Trichlorofluoromethane	101	2.350	2.350	0.000	97	871295	10.0	9.63	
17 Ethyl ether	59	2.564	2.564	0.000	90	406516	10.0	8.44	
18 Acrolein	56	2.682	2.682	0.000	99	222755	50.0	33.7	
19 1,1-Dichloroethene	96	2.777	2.777	0.000	98	544170	10.0	8.68	
20 1,1,2-Trichloro-1,2,2-trif	151	2.789	2.789	0.000	85	496817	10.0	9.65	
22 Acetone	58	2.813	2.813	0.000	99	80786	20.0	15.6	
23 Iodomethane	142	2.907	2.907	0.000	98	1002369	10.0	10.3	
24 Carbon disulfide	76	2.967	2.967	0.000	100	1316351	10.0	8.15	
25 3-Chloro-1-propene	76	3.073	3.073	0.000	89	311041	10.0	8.78	
27 Methyl acetate	43	3.085	3.085	0.000	98	574441	20.0	13.4	
28 Methylene Chloride	84	3.180	3.180	0.000	91	611898	10.0	9.34	
29 2-Methyl-2-propanol	59	3.263	3.263	0.000	90	245594	100.0	58.1	
32 Acrylonitrile	53	3.382	3.382	0.000	99	1499277	100.0	74.7	
31 Methyl tert-butyl ether	73	3.405	3.405	0.000	96	1189590	10.0	7.28	
30 trans-1,2-Dichloroethene	96	3.405	3.405	0.000	98	612223	10.0	10.5	
33 Hexane	86	3.619	3.619	0.000	94	110296	10.0	9.02	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
34 1,1-Dichloroethane	63	3.761	3.761	0.000	97	1015651	10.0	10.8	
36 Vinyl acetate	43	3.785	3.785	0.000	98	707360	10.0	6.52	
39 cis-1,2-Dichloroethene	96	4.236	4.236	0.000	80	697952	10.0	10.9	
41 2-Butanone (MEK)	72	4.247	4.247	0.000	98	91038	20.0	13.1	
40 2,2-Dichloropropane	97	4.247	4.247	0.000	82	148215	10.0	11.4	
45 Chlorobromomethane	128	4.437	4.437	0.000	91	331341	10.0	10.4	
46 Tetrahydrofuran	42	4.485	4.485	0.000	85	207071	20.0	12.0	
47 Chloroform	83	4.496	4.496	0.000	93	1062165	10.0	11.1	
48 1,1,1-Trichloroethane	97	4.662	4.662	0.000	99	864976	10.0	11.7	
49 Cyclohexane	56	4.710	4.710	0.000	91	808707	10.0	9.71	
51 1,1-Dichloropropene	75	4.793	4.793	0.000	94	779666	10.0	9.98	
50 Carbon tetrachloride	117	4.805	4.805	0.000	97	770426	10.0	10.7	
52 Isobutyl alcohol	41	4.864	4.864	0.000	89	269828	250.0	168.4	
53 Benzene	78	4.971	4.971	0.000	96	2483023	10.0	10.7	
54 1,2-Dichloroethane	62	4.983	4.983	0.000	98	793404	10.0	9.96	
56 n-Heptane	100	5.161	5.161	0.000	88	96418	10.0	8.69	
58 Trichloroethene	130	5.504	5.504	0.000	96	694244	10.0	10.3	
60 Methylcyclohexane	83	5.670	5.670	0.000	90	689042	10.0	9.37	
61 1,2-Dichloropropane	63	5.706	5.706	0.000	94	536224	10.0	9.93	
63 Dibromomethane	93	5.813	5.813	0.000	89	321708	10.0	8.98	
64 1,4-Dioxane	88	5.813	5.813	0.000	36	79951	200.0	180.9	
65 Dichlorobromomethane	83	5.931	5.931	0.000	99	663106	10.0	9.32	
67 2-Chloroethyl vinyl ether	63	6.180	6.180	0.000	94	432058	20.0	11.9	
68 cis-1,3-Dichloropropene	75	6.323	6.323	0.000	94	657632	10.0	7.99	
69 4-Methyl-2-pentanone (MIBK)	43	6.453	6.453	0.000	96	660185	20.0	11.8	
70 Toluene	91	6.619	6.619	0.000	98	2476981	10.0	10.4	
71 trans-1,3-Dichloropropene	75	6.809	6.809	0.000	95	531183	10.0	7.10	
72 Ethyl methacrylate	69	6.868	6.868	0.000	89	467554	10.0	6.54	
73 1,1,2-Trichloroethane	97	6.975	6.975	0.000	92	435239	10.0	9.07	
74 Tetrachloroethene	164	7.117	7.117	0.000	96	586613	10.0	11.2	
75 1,3-Dichloropropane	76	7.129	7.129	0.000	93	756851	10.0	8.97	
76 2-Hexanone	43	7.188	7.188	0.000	97	442908	20.0	12.1	
79 Chlorodibromomethane	129	7.343	7.343	0.000	89	432818	10.0	8.63	
80 Ethylene Dibromide	107	7.461	7.461	0.000	99	414517	10.0	8.07	
82 Chlorobenzene	112	7.900	7.900	0.000	95	1632775	10.0	10.4	
83 1,1,1,2-Tetrachloroethane	131	7.971	7.971	0.000	92	570358	10.0	10.3	
84 Ethylbenzene	106	7.995	7.995	0.000	98	829033	10.0	10.0	
85 m-Xylene & p-Xylene	91	8.101	8.101	0.000	93	1980348	10.0	10.1	
86 o-Xylene	106	8.481	8.481	0.000	95	1014701	10.0	10.2	
87 Styrene	104	8.493	8.493	0.000	92	1698792	10.0	10.2	
88 Bromoform	173	8.683	8.683	0.000	98	231585	10.0	6.29	
89 Isopropylbenzene	105	8.825	8.825	0.000	96	2371265	10.0	10.1	
93 1,1,2,2-Tetrachloroethane	83	9.121	9.121	0.000	98	489870	10.0	7.08	
92 Bromobenzene	156	9.145	9.145	0.000	88	762392	10.0	9.43	
95 trans-1,4-Dichloro-2-buten	53	9.169	9.169	0.000	67	65120	10.0	6.57	
94 1,2,3-Trichloropropane	110	9.169	9.169	0.000	84	168808	10.0	7.00	
96 N-Propylbenzene	120	9.228	9.228	0.000	98	665988	10.0	9.35	
97 2-Chlorotoluene	126	9.323	9.323	0.000	97	637518	10.0	9.57	
98 1,3,5-Trimethylbenzene	105	9.394	9.394	0.000	94	1960490	10.0	9.47	
99 4-Chlorotoluene	91	9.418	9.418	0.000	98	2076014	10.0	9.63	
101 tert-Butylbenzene	119	9.714	9.714	0.000	92	1593398	10.0	8.59	
103 1,2,4-Trimethylbenzene	105	9.762	9.762	0.000	94	2084740	10.0	9.58	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
105 sec-Butylbenzene	105	9.928	9.928	0.000	94	2037620	10.0	8.96	
106 1,3-Dichlorobenzene	146	10.046	10.046	0.000	99	1323675	10.0	9.51	
107 4-Isopropyltoluene	119	10.070	10.070	0.000	97	1876902	10.0	9.22	
108 1,4-Dichlorobenzene	146	10.141	10.141	0.000	96	1374527	10.0	9.52	
111 n-Butylbenzene	91	10.473	10.473	0.000	97	1360686	10.0	8.45	
112 1,2-Dichlorobenzene	146	10.509	10.509	0.000	99	1290266	10.0	9.52	
113 1,2-Dibromo-3-Chloropropan	157	11.291	11.291	0.000	92	72759	10.0	4.62	
115 1,2,4-Trichlorobenzene	180	12.110	12.110	0.000	93	702672	10.0	8.19	
116 Hexachlorobutadiene	225	12.276	12.276	0.000	91	373925	10.0	8.53	
117 Naphthalene	128	12.359	12.359	0.000	96	1206416	10.0	5.59	
118 1,2,3-Trichlorobenzene	180	12.608	12.608	0.000	95	650939	10.0	8.43	
S 129 Xylenes, Total	106				0		20.0	20.3	
S 157 Total BTEX	1		0.000				50.0	ND	
S 130 Trihalomethanes, Total	1				0		40.0	35.3	

Reagents:

VMRGAS_00285	Amount Added: 8.00	Units: uL	
VMRPRIMW_00325	Amount Added: 8.00	Units: uL	
VMAROLISTDW_00288	Amount Added: 8.00	Units: uL	
vm40ml_vials_00009	Amount Added: 0.00	Units:	Run Reagent
VM50IS_00075	Amount Added: 1.00	Units: uL	Run Reagent
vm50ss_stk_00080	Amount Added: 0.80	Units: uL	Run Reagent
vmDist_H2o_00140	Amount Added: 0.00	Units:	Run Reagent

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX15\20190318-85248.b\UXC9421.D

Injection Date: 18-Mar-2019 10:59:30

Instrument ID: A3UX15

Operator ID:

Lims ID: CCVIS L4 8260

Worklist Smp#: 2

Client ID:

Purge Vol: 5.000 mL

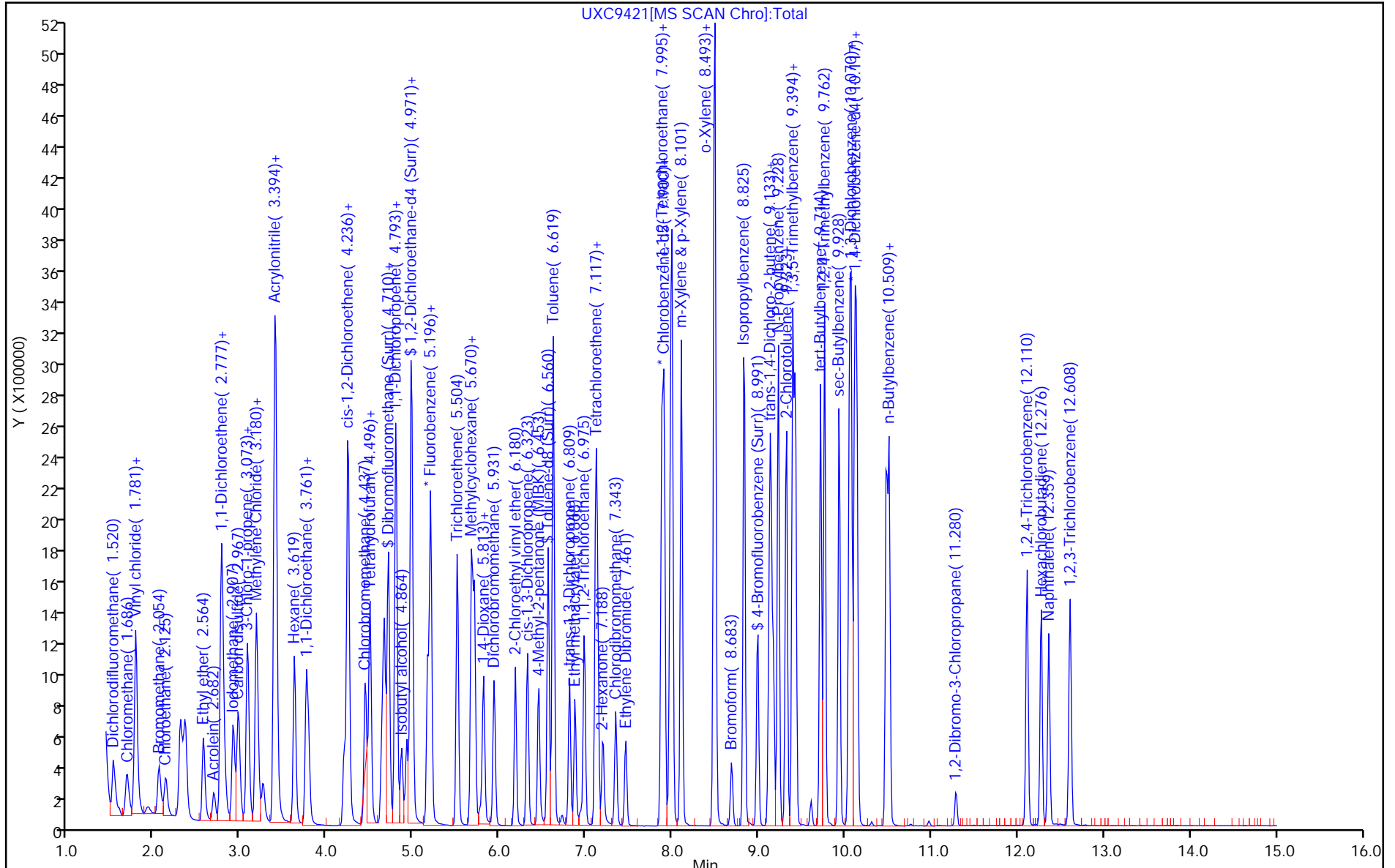
Dil. Factor: 1.0000

ALS Bottle#: 1

Method: 8260_15

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 367104

SDG No.: _____

Instrument ID: A3UX11 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/07/2019 13:09 Calibration End Date: 02/07/2019 15:00 Calibration ID: 49373

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD8260 240-367104/7	UXJ8026.D
Level 2	STD8260 240-367104/6	UXJ8025.D
Level 3	STD8260 240-367104/5	UXJ8024.D
Level 4	STD8260 240-367104/4	UXJ8023.D
Level 5	STD8260 240-367104/3	UXJ8022.D
Level 6	STD8260 240-367104/2	UXJ8021.D

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R ² OR COD	#	MIN R ² OR COD
	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
Dichlorodifluoromethane	0.2561 0.2965	0.3087	0.3159	0.2838	0.3181	Ave		0.2965			8.0		15.0				
Chloromethane	0.2639 0.2275	0.2688	0.2571	0.2172	0.2275	Ave		0.2437		0.1000	9.1		15.0				
Butadiene	0.2231 0.2211	0.2477	0.2327	0.1982	0.2411	Ave		0.2273			7.7		15.0				
Vinyl chloride	0.2701 0.2501	0.2708	0.2767	0.2462	0.2748	Ave		0.2648			5.0		15.0				
Bromomethane	0.1632 0.1581	0.1890	0.1875	0.1730	0.1786	Ave		0.1749			7.2		15.0				
Chloroethane	0.1115 0.1343	0.1496	0.1465	0.1424	0.1526	Ave		0.1395			10.8		15.0				
Dichlorofluoromethane	0.3612 0.3117	0.3753	0.3677	0.3417	0.3508	Ave		0.3514			6.5		15.0				
Trichlorofluoromethane	0.3705 0.3531	0.3652	0.4043	0.3445	0.3945	Ave		0.3720			6.3		15.0				
Ethyl ether	0.1930 0.1700	0.1835	0.1844	0.1896	0.1946	Ave		0.1858			4.8		15.0				
Acrolein	0.0439 0.0444	0.0477	0.0528	0.0475	0.0503	Ave		0.0478			7.1		15.0				
1,1-Dichloroethene	0.2250 0.2095	0.2341	0.2377	0.2326	0.2395	Ave		0.2297			4.8		15.0				
1,1,2-Trichloro-1,2,2-trifluoroethane	0.1811 0.1739	0.1966	0.2010	0.2060	0.1992	Ave		0.1930			6.5		15.0				
Acetone	0.1243 0.0933	0.1241	0.1066	0.0942	0.1010	Ave		0.1072			13.1		15.0				
Iodomethane	0.4087 0.3847	0.4067	0.4107	0.4206	0.4165	Ave		0.4080			3.1		15.0				
Carbon disulfide	0.8291 0.6716	0.8391	0.7182	0.7240	0.7387	Ave		0.7535			8.8		15.0				

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

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 367104

SDG No.: _____

Instrument ID: A3UX11 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/07/2019 13:09 Calibration End Date: 02/07/2019 15:00 Calibration ID: 49373

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R ² OR COD	#	MIN R ² OR COD
	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
3-Chloro-1-propene	0.1813 0.1508	0.1680	0.1668	0.1680	0.1816	Ave		0.1694			6.7		15.0				
Methyl acetate	0.2287 0.2144	0.2201	0.2459	0.2275	0.2383	Ave		0.2292			5.0		15.0				
Methylene Chloride	0.5096 0.2300	0.3875	0.2958	0.2714	0.2723	Lin1	0.2961	0.2352						0.9950		0.9900	
2-Methyl-2-propanol	0.0441 0.0356	0.0392	0.0377	0.0362	0.0395	Ave		0.0387			7.9		15.0				
Acrylonitrile	0.1160 0.1145	0.1217	0.1246	0.1171	0.1243	Ave		0.1197			3.7		15.0				
trans-1,2-Dichloroethene	0.2863 0.2479	0.2811	0.2728	0.2682	0.2837	Ave		0.2733			5.2		15.0				
Methyl tert-butyl ether	0.6061 0.5504	0.6345	0.6206	0.6201	0.6306	Ave		0.6104			5.1		15.0				
Hexane	0.0273 0.0675	0.0359	0.0643	0.0670	0.0711	Lin1	-0.048	0.0705						0.9980		0.9900	
1,1-Dichloroethane	0.4752 0.3990	0.4658	0.4289	0.4510	0.4670	Ave		0.4478		0.1000	6.5		15.0				
Vinyl acetate	0.3823 0.4460	0.4155	0.4245	0.4177	0.4413	Ave		0.4212			5.4		15.0				
2,2-Dichloropropane	0.0742 0.0561	0.0707	0.0639	0.0641	0.0671	Ave		0.0660			9.5		15.0				
cis-1,2-Dichloroethene	0.2779 0.2586	0.3025	0.2735	0.2928	0.3019	Ave		0.2845			6.2		15.0				
2-Butanone (MEK)	0.1145 0.1384	0.1443	0.1550	0.1256	0.1327	Ave		0.1351			10.5		15.0				
Chlorobromomethane	0.1533 0.1283	0.1594	0.1443	0.1431	0.1455	Ave		0.1457			7.2		15.0				
Tetrahydrofuran	0.0937 0.0892	0.0870	0.0940	0.0860	0.0939	Ave		0.0906			4.1		15.0				
Chloroform	0.4187 0.3891	0.4465	0.4523	0.4492	0.4466	Ave		0.4338			5.8		15.0				
1,1,1-Trichloroethane	0.4191 0.3332	0.3915	0.3584	0.3779	0.3844	Ave		0.3774			7.8		15.0				
Cyclohexane	0.2959 0.3516	0.3335	0.3536	0.3776	0.3932	Ave		0.3509			9.7		15.0				
1,1-Dichloropropene	0.3410 0.3388	0.3355	0.3288	0.3648	0.3649	Ave		0.3456			4.5		15.0				
Carbon tetrachloride	0.3771 0.3264	0.3538	0.3631	0.3545	0.3679	Ave		0.3572			4.9		15.0				

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

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 367104

SDG No.: _____

Instrument ID: A3UX11 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/07/2019 13:09 Calibration End Date: 02/07/2019 15:00 Calibration ID: 49373

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R ² OR COD	#	MIN R ² OR COD
	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
Isobutyl alcohol	0.0201 0.0189	0.0188	0.0203	0.0190	0.0209	Ave		0.0197			4.5		15.0				
Benzene	1.0033 0.9512	1.0488	0.9813	1.0084	1.0137	Ave		1.0011			3.3		15.0				
1,2-Dichloroethane	0.3636 0.3225	0.3571	0.3445	0.3564	0.3440	Ave		0.3480			4.2		15.0				
n-Heptane	0.0532 0.0561	0.0394	0.0663	0.0573	0.0616	Lin1	-0.009	0.0586						0.9930		0.9900	
Trichloroethene	0.3229 0.2740	0.3112	0.2971	0.2881	0.2932	Ave		0.2978			5.8		15.0				
Methylcyclohexane	0.3518 0.3869	0.3966	0.3694	0.3965	0.4296	Ave		0.3885			6.8		15.0				
1,2-Dichloropropane	0.1979 0.2052	0.2205	0.2101	0.2180	0.2083	Ave		0.2100			4.0		15.0				
Dibromomethane	0.1479 0.1403	0.1518	0.1482	0.1474	0.1493	Ave		0.1475			2.6		15.0				
1,4-Dioxane	0.0021 0.0037	0.0023	0.0036	0.0028	0.0033	Lin1	-0.039	0.0036						0.9910		0.9900	
Dichlorobromomethane	0.3397 0.3033	0.3234	0.3316	0.3140	0.3093	Ave		0.3202			4.3		15.0				
2-Chloroethyl vinyl ether	0.1068 0.1514	0.1286	0.1429	0.1184	0.1244	Ave		0.1287			12.6		15.0				
cis-1,3-Dichloropropene	0.3057 0.3545	0.3267	0.3461	0.3301	0.3306	Ave		0.3323			5.1		15.0				
4-Methyl-2-pentanone (MIBK)	0.2124 0.2572	0.2218	0.2653	0.2400	0.2479	Ave		0.2408			8.5		15.0				
Toluene	1.4332 1.4707	1.3622	1.4674	1.4473	1.4836	Ave		1.4441			3.0		15.0				
trans-1,3-Dichloropropene	0.4847 0.4784	0.4842	0.4783	0.4458	0.4500	Ave		0.4702			3.7		15.0				
Ethyl methacrylate	0.3475 0.4188	0.3758	0.3944	0.3988	0.4135	Ave		0.3915			6.7		15.0				
1,1,2-Trichloroethane	0.3624 0.2742	0.3027	0.3051	0.2949	0.2806	Ave		0.3033			10.3		15.0				
Tetrachloroethene	0.3266 0.3072	0.3132	0.3417	0.3262	0.3224	Ave		0.3229			3.7		15.0				
1,3-Dichloropropane	0.5058 0.4783	0.4774	0.4948	0.4817	0.4770	Ave		0.4858			2.4		15.0				
2-Hexanone	0.2241 0.2630	0.2662	0.2752	0.2548	0.2705	Ave		0.2590			7.1		15.0				

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

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 367104

SDG No.: _____

Instrument ID: A3UX11 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/07/2019 13:09 Calibration End Date: 02/07/2019 15:00 Calibration ID: 49373

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R ² OR COD	#	MIN R ² OR COD
	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
Chlorodibromomethane	0.3113 0.3375	0.3763	0.3706	0.3601	0.3570	Ave		0.3521			6.8		15.0				
Ethylene Dibromide	0.3621 0.3033	0.2878	0.3056	0.3062	0.3105	Ave		0.3126			8.2		15.0				
Chlorobenzene	1.0968 0.9164	0.9427	0.9575	0.9435	0.9510	Ave		0.9680		0.3000	6.7		15.0				
1,1,1,2-Tetrachloroethane	0.4421 0.3288	0.3620	0.3454	0.3664	0.3830	Ave		0.3713			10.6		15.0				
Ethylbenzene	0.5029 0.4881	0.4436	0.5010	0.4714	0.5084	Ave		0.4859			5.1		15.0				
m-Xylene & p-Xylene	0.5018 0.6004	0.4989	0.5786	0.5987	0.6211	Ave		0.5666			9.4		15.0				
o-Xylene	0.4845 0.5684	0.5086	0.5333	0.5884	0.6066	Ave		0.5483			8.7		15.0				
Styrene	0.7740 1.0043	0.8247	0.9335	0.9760	1.0322	Ave		0.9241			11.2		15.0				
Bromoform	0.2683 0.2637	0.2827	0.2965	0.2859	0.2852	Ave		0.2804		0.1000	4.3		15.0				
Isopropylbenzene	1.3016 1.4809	1.2282	1.3930	1.4757	1.5831	Ave		1.4104			9.2		15.0				
Bromobenzene	0.7752 0.7796	0.7507	0.7234	0.7177	0.7506	Ave		0.7495			3.4		15.0				
1,1,2,2-Tetrachloroethane	0.8549 0.8003	0.7913	0.8163	0.7838	0.8193	Ave		0.8110		0.3000	3.2		15.0				
1,2,3-Trichloropropane	0.2749 0.2776	0.3041	0.2874	0.2705	0.2749	Ave		0.2816			4.4		15.0				
trans-1,4-Dichloro-2-butene	0.2483 0.2604	0.2632	0.2915	0.2366	0.2526	Ave		0.2588			7.2		15.0				
N-Propylbenzene	0.6807 0.8566	0.6482	0.7462	0.7870	0.8370	Ave		0.7593			11.0		15.0				
2-Chlorotoluene	0.6217 0.7300	0.6045	0.6908	0.7052	0.7308	Ave		0.6805			8.0		15.0				
1,3,5-Trimethylbenzene	1.8263 2.3213	1.7650	2.0474	2.1790	2.2664	Ave		2.0676			11.2		15.0				
4-Chlorotoluene	0.6198 0.7574	0.6373	0.7346	0.7031	0.7299	Ave		0.6970			8.0		15.0				
tert-Butylbenzene	1.5194 2.0713	1.4757	1.8069	1.7744	1.9643	Ave		1.7687			13.4		15.0				
1,2,4-Trimethylbenzene	1.8163 2.3412	1.7839	2.0647	2.1935	2.3174	Ave		2.0862			11.6		15.0				

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

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 367104

SDG No.: _____

Instrument ID: A3UX11 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/07/2019 13:09 Calibration End Date: 02/07/2019 15:00 Calibration ID: 49373

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R ² OR COD	#	MIN R ² OR COD
	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
sec-Butylbenzene	2.2345 2.8648	2.2736	2.4455	2.4879	2.8048	Ave		2.5185			10.5		15.0				
1,3-Dichlorobenzene	1.3721 1.3792	1.3531	1.3430	1.3524	1.3792	Ave		1.3632			1.1		15.0				
4-Isopropyltoluene	1.8926 2.5053	1.9908	2.1641	2.3164	2.5054	Ave		2.2291			11.6		15.0				
1,4-Dichlorobenzene	1.7104 1.3790	1.5194	1.3950	1.3809	1.3892	Ave		1.4623			9.1		15.0				
1,2-Dichlorobenzene	1.2698 1.2958	1.2989	1.2234	1.2284	1.3312	Ave		1.2746			3.3		15.0				
n-Butylbenzene	1.7842 2.0152	1.5898	1.7143	1.8014	2.0030	Ave		1.8180			9.1		15.0				
1,2-Dibromo-3-Chloropropane	0.3467 0.2339	0.2240	0.2456	0.2445	0.2477	Lin1	0.0693	0.2354						0.9980		0.9900	
1,2,4-Trichlorobenzene	0.8504 0.8238	0.7825	0.7622	0.7359	0.8577	Ave		0.8021			6.2		15.0				
Hexachlorobutadiene	0.4299 0.3238	0.4004	0.3479	0.3514	0.3693	Ave		0.3705			10.4		15.0				
Naphthalene	2.0320 2.5297	1.9621	1.8141	1.9407	2.5249	Ave		2.1339			14.7		15.0				
1,2,3-Trichlorobenzene	0.7838 0.7673	0.7031	0.7101	0.7051	0.8441	Ave		0.7522			7.5		15.0				
Dibromofluoromethane (Surr)	0.2467 0.2099	0.2246	0.2361	0.2411	0.2415	Ave		0.2333			5.9		15.0				
1,2-Dichloroethane-d4 (Surr)	0.2992 0.2455	0.3054	0.2970	0.2790	0.2681	Ave		0.2824			8.1		15.0				
Toluene-d8 (Surr)	1.1868 1.1915	1.0838	1.1660	1.2217	1.2705	Ave		1.1867			5.2		15.0				
4-Bromofluorobenzene (Surr)	0.3769 0.3566	0.3761	0.4007	0.3903	0.3768	Ave		0.3796			3.9		15.0				

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

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 367104

SDG No.: _____

Instrument ID: A3UX11 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/07/2019 13:09 Calibration End Date: 02/07/2019 15:00 Calibration ID: 49373

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD8260 240-367104/7	UXJ8026.D
Level 2	STD8260 240-367104/6	UXJ8025.D
Level 3	STD8260 240-367104/5	UXJ8024.D
Level 4	STD8260 240-367104/4	UXJ8023.D
Level 5	STD8260 240-367104/3	UXJ8022.D
Level 6	STD8260 240-367104/2	UXJ8021.D

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/L)				
			LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5
Dichlorodifluoromethane	FB	Ave	14302 858734	33811	91261	166318	393607	1.00 40.0	2.00	5.00	10.0	20.0
Chloromethane	FB	Ave	14735 658748	29438	74271	127316	281547	1.00 40.0	2.00	5.00	10.0	20.0
Butadiene	FB	Ave	12457 640253	27131	67238	116135	298381	1.00 40.0	2.00	5.00	10.0	20.0
Vinyl chloride	FB	Ave	15085 724231	29665	79953	144260	339988	1.00 40.0	2.00	5.00	10.0	20.0
Bromomethane	FB	Ave	9111 457891	20704	54178	101367	220958	1.00 40.0	2.00	5.00	10.0	20.0
Chloroethane	FB	Ave	6225 388823	16382	42315	83470	188887	1.00 40.0	2.00	5.00	10.0	20.0
Dichlorofluoromethane	FB	Ave	20171 902632	41103	106239	200231	434053	1.00 40.0	2.00	5.00	10.0	20.0
Trichlorofluoromethane	FB	Ave	20689 1022513	40004	116801	201892	488108	1.00 40.0	2.00	5.00	10.0	20.0
Ethyl ether	FB	Ave	10777 492212	20099	53290	111089	240821	1.00 40.0	2.00	5.00	10.0	20.0
Acrolein	FB	Ave	12263 642909	26117	76307	139233	311245	5.00 200	10.0	25.0	50.0	100
1,1-Dichloroethene	FB	Ave	12563 606755	25640	68669	136302	296326	1.00 40.0	2.00	5.00	10.0	20.0
1,1,2-Trichloro-1,2,2-trifluoroethane	FB	Ave	10111 503676	21536	58077	120721	246455	1.00 40.0	2.00	5.00	10.0	20.0
Acetone	FB	Ave	13879 540581	27192	61588	110371	249848	2.00 80.0	4.00	10.0	20.0	40.0
Iodomethane	FB	Ave	22823 1113991	44542	118663	246503	515391	1.00 40.0	2.00	5.00	10.0	20.0
Carbon disulfide	FB	Ave	46298 1945109	91902	207524	424295	914051	1.00 40.0	2.00	5.00	10.0	20.0
3-Chloro-1-propene	FB	Ave	10125 436826	18405	48182	98437	224690	1.00 40.0	2.00	5.00	10.0	20.0

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 367104

SDG No.: _____

Instrument ID: A3UX11 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/07/2019 13:09 Calibration End Date: 02/07/2019 15:00 Calibration ID: 49373

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/L)				
			LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5
Methyl acetate	FB	Ave	25546 1242117	48220	142070	266658	589687	2.00 80.0	4.00	10.0	20.0	40.0
Methylene Chloride	FB	Lin1	28456 666164	42445	85471	159040	336958	1.00 40.0	2.00	5.00	10.0	20.0
2-Methyl-2-propanol	FB	Ave	24619 1030644	42922	108992	212100	489028	10.0 400	20.0	50.0	100	200
Acrylonitrile	FB	Ave	64753 3314967	133260	360130	686402	1538522	10.0 400	20.0	50.0	100	200
trans-1,2-Dichloroethene	FB	Ave	15989 717821	30791	78810	157177	351024	1.00 40.0	2.00	5.00	10.0	20.0
Methyl tert-butyl ether	FB	Ave	33845 1594022	69489	179324	363372	780350	1.00 40.0	2.00	5.00	10.0	20.0
Hexane	FB	Lin1	1522 195544	3932	18570	39284	87974	1.00 40.0	2.00	5.00	10.0	20.0
1,1-Dichloroethane	FB	Ave	26535 1155574	51015	123928	264327	577900	1.00 40.0	2.00	5.00	10.0	20.0
Vinyl acetate	FB	Ave	21348 1291718	45507	122657	244809	546119	1.00 40.0	2.00	5.00	10.0	20.0
2,2-Dichloropropane	FB	Ave	4143 162380	7740	18470	37577	83030	1.00 40.0	2.00	5.00	10.0	20.0
cis-1,2-Dichloroethene	FB	Ave	15518 748792	33131	79034	171580	373545	1.00 40.0	2.00	5.00	10.0	20.0
2-Butanone (MEK)	FB	Ave	12784 801645	31612	89550	147163	328459	2.00 80.0	4.00	10.0	20.0	40.0
Chlorobromomethane	FB	Ave	8562 371594	17457	41697	83844	180067	1.00 40.0	2.00	5.00	10.0	20.0
Tetrahydrofuran	FB	Ave	10465 516929	19047	54325	100847	232324	2.00 80.0	4.00	10.0	20.0	40.0
Chloroform	FB	Ave	23384 1126807	48908	130678	263272	552690	1.00 40.0	2.00	5.00	10.0	20.0
1,1,1-Trichloroethane	FB	Ave	23404 964948	42875	103539	221456	475687	1.00 40.0	2.00	5.00	10.0	20.0
Cyclohexane	FB	Ave	16522 1018178	36529	102175	221283	486587	1.00 40.0	2.00	5.00	10.0	20.0
1,1-Dichloropropene	FB	Ave	19040 981099	36750	95002	213777	451488	1.00 40.0	2.00	5.00	10.0	20.0
Carbon tetrachloride	FB	Ave	21059 945383	38755	104922	207777	455300	1.00 40.0	2.00	5.00	10.0	20.0
Isobutyl alcohol	CBNZ d5	Ave	17638 920911	35587	100691	178315	411928	25.0 1000	50.0	125	250	500
Benzene	FB	Ave	56029 2754879	114875	283539	590944	1254314	1.00 40.0	2.00	5.00	10.0	20.0

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 367104

SDG No.: _____

Instrument ID: A3UX11 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/07/2019 13:09 Calibration End Date: 02/07/2019 15:00 Calibration ID: 49373

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/L)				
			LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5
1,2-Dichloroethane	FB	Ave	20305 933881	39111	99528	208883	425614	1.00 40.0	2.00	5.00	10.0	20.0
n-Heptane	FB	Lin1	2973 162413	4317	19161	33582	76210	1.00 40.0	2.00	5.00	10.0	20.0
Trichloroethene	FB	Ave	18033 793578	34086	85827	168836	362851	1.00 40.0	2.00	5.00	10.0	20.0
Methylcyclohexane	FB	Ave	19643 1120599	43436	106738	232379	531548	1.00 40.0	2.00	5.00	10.0	20.0
1,2-Dichloropropane	FB	Ave	11053 594323	24147	60714	127764	257754	1.00 40.0	2.00	5.00	10.0	20.0
Dibromomethane	FB	Ave	8259 406414	16623	42824	86362	184789	1.00 40.0	2.00	5.00	10.0	20.0
1,4-Dioxane	FB	Lin1	2371 216961	5017	20834	33357	81643	20.0 800	40.0	100	200	400
Dichlorobromomethane	FB	Ave	18968 878356	35418	95815	184018	382766	1.00 40.0	2.00	5.00	10.0	20.0
2-Chloroethyl vinyl ether	FB	Ave	11930 876953	28161	82558	138767	307879	2.00 80.0	4.00	10.0	20.0	40.0
cis-1,3-Dichloropropene	FB	Ave	17074 1026562	35784	99988	193425	409067	1.00 40.0	2.00	5.00	10.0	20.0
4-Methyl-2-pentanone (MIBK)	FB	Ave	23723 1489645	48587	153307	281355	613490	2.00 80.0	4.00	10.0	20.0	40.0
Toluene	CBNZ d5	Ave	50368 2872531	103151	290746	543872	1170735	1.00 40.0	2.00	5.00	10.0	20.0
trans-1,3-Dichloropropene	CBNZ d5	Ave	17032 934338	36666	94776	167511	355082	1.00 40.0	2.00	5.00	10.0	20.0
Ethyl methacrylate	CBNZ d5	Ave	12213 818004	28459	78145	149850	326308	1.00 40.0	2.00	5.00	10.0	20.0
1,1,2-Trichloroethane	CBNZ d5	Ave	12734 535501	22922	60454	110808	221396	1.00 40.0	2.00	5.00	10.0	20.0
Tetrachloroethene	CBNZ d5	Ave	11477 600080	23718	67708	122571	254413	1.00 40.0	2.00	5.00	10.0	20.0
1,3-Dichloropropane	CBNZ d5	Ave	17776 934162	36153	98029	181024	376403	1.00 40.0	2.00	5.00	10.0	20.0
2-Hexanone	CBNZ d5	Ave	15753 1027463	40322	109052	191478	426962	2.00 80.0	4.00	10.0	20.0	40.0
Chlorodibromomethane	CBNZ d5	Ave	10939 659131	28493	73418	135325	281679	1.00 40.0	2.00	5.00	10.0	20.0
Ethylene Dibromide	CBNZ d5	Ave	12725 592368	21792	60552	115065	245052	1.00 40.0	2.00	5.00	10.0	20.0
Chlorobenzene	CBNZ d5	Ave	38544 1789883	71386	189707	354543	750415	1.00 40.0	2.00	5.00	10.0	20.0

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 367104

SDG No.: _____

Instrument ID: A3UX11 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/07/2019 13:09 Calibration End Date: 02/07/2019 15:00 Calibration ID: 49373

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/L)				
			LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5
1,1,1,2-Tetrachloroethane	CBNZ d5	Ave	15538 642189	27415	68426	137694	302221	1.00 40.0	2.00	5.00	10.0	20.0
Ethylbenzene	CBNZ d5	Ave	17672 953361	33590	99260	177149	401219	1.00 40.0	2.00	5.00	10.0	20.0
m-Xylene & p-Xylene	CBNZ d5	Ave	17633 1172678	37781	114642	225000	490117	1.00 40.0	2.00	5.00	10.0	20.0
o-Xylene	CBNZ d5	Ave	17027 1110191	38511	105654	221095	478708	1.00 40.0	2.00	5.00	10.0	20.0
Styrene	CBNZ d5	Ave	27201 1961423	62454	184963	366782	814512	1.00 40.0	2.00	5.00	10.0	20.0
Bromoform	CBNZ d5	Ave	9430 515031	21407	58743	107418	225037	1.00 40.0	2.00	5.00	10.0	20.0
Isopropylbenzene	CBNZ d5	Ave	45741 2892282	93008	276000	554538	1249226	1.00 40.0	2.00	5.00	10.0	20.0
Bromobenzene	DCBd 4	Ave	14505 744364	30472	77365	149296	318510	1.00 40.0	2.00	5.00	10.0	20.0
1,1,2,2-Tetrachloroethane	DCBd 4	Ave	15997 764142	32123	87302	163051	347666	1.00 40.0	2.00	5.00	10.0	20.0
1,2,3-Trichloropropane	DCBd 4	Ave	5144 265100	12345	30733	56270	116642	1.00 40.0	2.00	5.00	10.0	20.0
trans-1,4-Dichloro-2-butene	DCBd 4	Ave	4646 248656	10685	31177	49221	107185	1.00 40.0	2.00	5.00	10.0	20.0
N-Propylbenzene	DCBd 4	Ave	12737 817872	26311	79799	163725	355179	1.00 40.0	2.00	5.00	10.0	20.0
2-Chlorotoluene	DCBd 4	Ave	11634 697021	24540	73878	146693	310087	1.00 40.0	2.00	5.00	10.0	20.0
1,3,5-Trimethylbenzene	DCBd 4	Ave	34175 2216449	71649	218959	453276	961736	1.00 40.0	2.00	5.00	10.0	20.0
4-Chlorotoluene	DCBd 4	Ave	11598 723168	25871	78564	146255	309733	1.00 40.0	2.00	5.00	10.0	20.0
tert-Butylbenzene	DCBd 4	Ave	28431 1977723	59904	193232	369118	833511	1.00 40.0	2.00	5.00	10.0	20.0
1,2,4-Trimethylbenzene	DCBd 4	Ave	33988 2235448	72413	220805	456294	983336	1.00 40.0	2.00	5.00	10.0	20.0
sec-Butylbenzene	DCBd 4	Ave	41813 2735311	92291	261531	517546	1190193	1.00 40.0	2.00	5.00	10.0	20.0
1,3-Dichlorobenzene	DCBd 4	Ave	25676 1316887	54925	143622	281323	585237	1.00 40.0	2.00	5.00	10.0	20.0
4-Isopropyltoluene	DCBd 4	Ave	35416 2392076	80813	231439	481868	1063145	1.00 40.0	2.00	5.00	10.0	20.0
1,4-Dichlorobenzene	DCBd 4	Ave	32006 1316655	61678	149185	287252	589475	1.00 40.0	2.00	5.00	10.0	20.0

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 367104

SDG No.: _____

Instrument ID: A3UX11 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/07/2019 13:09 Calibration End Date: 02/07/2019 15:00 Calibration ID: 49373

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/L)				
			LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5
1,2-Dichlorobenzene	DCBd 4	Ave	23761 1237285	52725	130837	255531	564881	1.00 40.0	2.00	5.00	10.0	20.0
n-Butylbenzene	DCBd 4	Ave	33386 1924106	64537	183328	374733	849937	1.00 40.0	2.00	5.00	10.0	20.0
1,2-Dibromo-3-Chloropropane	DCBd 4	Lin1	6487 223296	9092	26261	50852	105103	1.00 40.0	2.00	5.00	10.0	20.0
1,2,4-Trichlorobenzene	DCBd 4	Ave	15913 786571	31765	81516	153092	363972	1.00 40.0	2.00	5.00	10.0	20.0
Hexachlorobutadiene	DCBd 4	Ave	8045 309212	16252	37204	73104	156722	1.00 40.0	2.00	5.00	10.0	20.0
Naphthalene	DCBd 4	Ave	38023 2415397	79647	194010	403718	1071425	1.00 40.0	2.00	5.00	10.0	20.0
1,2,3-Trichlorobenzene	DCBd 4	Ave	14666 732633	28542	75944	146668	358175	1.00 40.0	2.00	5.00	10.0	20.0
Dibromofluoromethane (Surr)	FB	Ave	13778 607914	24605	68214	141317	298829	1.00 40.0	2.00	5.00	10.0	20.0
1,2-Dichloroethane-d4 (Surr)	FB	Ave	16708 711134	33449	85806	163525	331727	1.00 40.0	2.00	5.00	10.0	20.0
Toluene-d8 (Surr)	CBNZ d5	Ave	41706 2327197	82074	231027	459092	1002587	1.00 40.0	2.00	5.00	10.0	20.0
4-Bromofluorobenzene (Surr)	CBNZ d5	Ave	13244 696421	28481	79386	146668	297359	1.00 40.0	2.00	5.00	10.0	20.0

Curve Type Legend:

Ave = Average ISTD
Lin1 = Linear 1/conc ISTD

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8021.D
 Lims ID: STD8260 L6
 Client ID:
 Sample Type: IC Calib Level: 6
 Inject. Date: 07-Feb-2019 13:09:30 ALS Bottle#: 1 Worklist Smp#: 2
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0084305-002
 Misc. Info.: J90207A-IC,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Sublist: chrom-8260_11*sub78
 Method: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 08-Feb-2019 09:14:31 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0324

First Level Reviewer: evansle

Date: 08-Feb-2019 08:04:40

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	4.647	4.659	-0.012	94	1448052	20.0	20.0	
* 2 Chlorobenzene-d5	117	7.297	7.309	-0.012	82	976557	20.0	20.0	
* 3 1,4-Dichlorobenzene-d4	152	9.522	9.522	0.000	91	477407	20.0	20.0	
\$ 4 Dibromofluoromethane (Surr	113	4.091	4.103	-0.012	95	607914	40.0	36.0	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.363	4.375	-0.012	80	711134	40.0	34.8	
\$ 6 Toluene-d8 (Surr)	98	6.008	6.008	0.000	92	2327197	40.0	40.2	
\$ 7 4-Bromofluorobenzene (Surr	95	8.410	8.410	0.000	86	696421	40.0	37.6	
9 Dichlorodifluoromethane	85	1.156	1.168	-0.012	100	858734	40.0	40.0	
10 Chloromethane	50	1.298	1.310	-0.012	100	658748	40.0	37.3	
13 Butadiene	54	1.357	1.369	-0.012	97	640253	40.0	38.9	
12 Vinyl chloride	62	1.369	1.393	-0.024	98	724231	40.0	37.8	
15 Bromomethane	94	1.570	1.606	-0.036	91	457891	40.0	36.2	
16 Chloroethane	64	1.641	1.665	-0.024	100	388823	40.0	38.5	
17 Dichlorofluoromethane	67	1.795	1.807	-0.012	98	902632	40.0	35.5	
18 Trichlorofluoromethane	101	1.843	1.831	0.012	99	1022513	40.0	38.0	
19 Ethyl ether	59	2.056	2.067	-0.011	89	492212	40.0	36.6	
20 Acrolein	56	2.150	2.162	-0.012	99	642909	200.0	185.9	
21 1,1-Dichloroethene	96	2.233	2.245	-0.012	98	606755	40.0	36.5	
22 1,1,2-Trichloro-1,2,2-trif	151	2.257	2.280	-0.023	90	503676	40.0	36.1	
23 Acetone	43	2.280	2.292	-0.012	100	540581	80.0	69.6	
25 Iodomethane	142	2.351	2.363	-0.012	96	1113991	40.0	37.7	
26 Carbon disulfide	76	2.399	2.411	-0.011	99	1945109	40.0	35.7	
28 3-Chloro-1-propene	76	2.529	2.541	-0.012	88	436826	40.0	35.6	
29 Methyl acetate	43	2.552	2.564	-0.012	97	1242117	80.0	74.9	
30 Methylene Chloride	84	2.635	2.647	-0.012	88	666164	40.0	37.9	
31 2-Methyl-2-propanol	59	2.765	2.754	0.011	97	1030644	400.0	367.7	
32 Acrylonitrile	53	2.836	2.848	-0.012	98	3314967	400.0	382.5	
34 trans-1,2-Dichloroethene	96	2.860	2.872	-0.012	98	717821	40.0	36.3	
33 Methyl tert-butyl ether	73	2.872	2.884	-0.012	94	1594022	40.0	36.1	
35 Hexane	86	3.097	3.109	-0.012	92	195544	40.0	39.0	

Compound	Sig	RT (min.)	Exp RT (min.)	Diff RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
36 1,1-Dichloroethane	63	3.203	3.215	-0.012	97	1155574	40.0	35.6	
37 Vinyl acetate	43	3.262	3.274	-0.012	97	1291718	40.0	42.4	
41 cis-1,2-Dichloroethene	96	3.688	3.700	-0.012	82	748792	40.0	36.3	
43 2,2-Dichloropropane	97	3.688	3.700	-0.012	59	162380	40.0	34.0	
42 2-Butanone (MEK)	43	3.712	3.712	0.000	98	801645	80.0	82.0	
47 Chlorobromomethane	128	3.890	3.890	0.000	90	371594	40.0	35.2	
48 Tetrahydrofuran	42	3.925	3.937	-0.012	84	516929	80.0	78.8	
49 Chloroform	83	3.961	3.961	0.000	94	1126807	40.0	35.9	
50 1,1,1-Trichloroethane	97	4.114	4.114	0.000	98	964948	40.0	35.3	
51 Cyclohexane	56	4.150	4.162	-0.012	88	1018178	40.0	40.1	
53 Carbon tetrachloride	117	4.245	4.256	-0.011	90	945383	40.0	36.6	
52 1,1-Dichloropropene	75	4.245	4.256	-0.011	95	981099	40.0	39.2	
54 Isobutyl alcohol	41	4.363	4.363	0.000	93	920911	1000.0	959.6	
55 Benzene	78	4.422	4.422	0.000	95	2754879	40.0	38.0	
56 1,2-Dichloroethane	62	4.434	4.434	0.000	98	933881	40.0	37.1	
58 n-Heptane	100	4.659	4.659	0.000	71	162413	40.0	38.5	
60 Trichloroethene	130	4.966	4.966	0.000	94	793578	40.0	36.8	
62 Methylcyclohexane	83	5.120	5.132	-0.012	89	1120599	40.0	39.8	
63 1,2-Dichloropropane	63	5.144	5.156	-0.012	95	594323	40.0	39.1	
65 Dibromomethane	93	5.250	5.250	0.000	90	406414	40.0	38.1	
66 1,4-Dioxane	88	5.274	5.274	0.000	91	216961	800.0	843.3	
67 Dichlorobromomethane	83	5.381	5.392	-0.011	99	878356	40.0	37.9	
69 2-Chloroethyl vinyl ether	63	5.653	5.653	0.000	92	876953	80.0	94.1	
70 cis-1,3-Dichloropropene	75	5.771	5.771	0.000	96	1026562	40.0	42.7	
71 4-Methyl-2-pentanone (MIBK)	43	5.913	5.913	0.000	96	1489645	80.0	85.5	
72 Toluene	91	6.067	6.067	0.000	99	2872531	40.0	40.7	
73 trans-1,3-Dichloropropene	75	6.268	6.268	0.000	92	934338	40.0	40.7	
74 Ethyl methacrylate	69	6.351	6.363	-0.012	88	818004	40.0	42.8	
75 1,1,2-Trichloroethane	97	6.422	6.422	0.000	92	535501	40.0	36.2	
76 Tetrachloroethene	164	6.552	6.552	0.000	98	600080	40.0	38.1	
77 1,3-Dichloropropane	76	6.576	6.576	0.000	88	934162	40.0	39.4	
78 2-Hexanone	43	6.659	6.659	0.000	95	1027463	80.0	81.3	
80 Chlorodibromomethane	129	6.777	6.777	0.000	90	659131	40.0	38.3	
82 Ethylene Dibromide	107	6.872	6.872	0.000	99	592368	40.0	38.8	
84 Chlorobenzene	112	7.333	7.333	0.000	98	1789883	40.0	37.9	
85 1,1,1,2-Tetrachloroethane	131	7.404	7.404	0.000	96	642189	40.0	35.4	
86 Ethylbenzene	106	7.439	7.439	0.000	97	953361	40.0	40.2	
87 m-Xylene & p-Xylene	106	7.546	7.546	0.000	97	1172678	40.0	42.4	
88 o-Xylene	106	7.913	7.913	0.000	96	1110191	40.0	41.5	
89 Styrene	104	7.925	7.925	0.000	94	1961423	40.0	43.5	
90 Bromoform	173	8.090	8.090	0.000	98	515031	40.0	37.6	
91 Isopropylbenzene	105	8.268	8.268	0.000	95	2892282	40.0	42.0	
95 Bromobenzene	156	8.540	8.540	0.000	88	744364	40.0	41.6	
94 1,1,2,2-Tetrachloroethane	83	8.552	8.552	0.000	94	764142	40.0	39.5	
96 1,2,3-Trichloropropane	110	8.587	8.587	0.000	86	265100	40.0	39.4	
97 trans-1,4-Dichloro-2-buten	53	8.611	8.611	0.000	80	248656	40.0	40.3	
98 N-Propylbenzene	120	8.658	8.658	0.000	99	817872	40.0	45.1	
99 2-Chlorotoluene	126	8.741	8.741	0.000	98	697021	40.0	42.9	
100 1,3,5-Trimethylbenzene	105	8.836	8.836	0.000	96	2216449	40.0	44.9	
101 4-Chlorotoluene	126	8.848	8.848	0.000	98	723168	40.0	43.5	
102 tert-Butylbenzene	119	9.143	9.155	-0.012	91	1977723	40.0	46.8	
104 1,2,4-Trimethylbenzene	105	9.191	9.191	0.000	95	2235448	40.0	44.9	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
105 sec-Butylbenzene	105	9.356	9.356	0.000	93	2735311	40.0	45.5	
106 1,3-Dichlorobenzene	146	9.451	9.451	0.000	99	1316887	40.0	40.5	
107 4-Isopropyltoluene	119	9.510	9.510	0.000	97	2392076	40.0	45.0	
108 1,4-Dichlorobenzene	146	9.546	9.546	0.000	97	1316655	40.0	37.7	
111 n-Butylbenzene	91	9.901	9.901	0.000	97	1924106	40.0	44.3	
112 1,2-Dichlorobenzene	146	9.901	9.901	0.000	98	1237285	40.0	40.7	
114 1,2-Dibromo-3-Chloropropan	157	10.670	10.670	0.000	93	223296	40.0	39.4	
116 1,2,4-Trichlorobenzene	180	11.486	11.486	0.000	93	786571	40.0	41.1	
117 Hexachlorobutadiene	225	11.676	11.676	0.000	96	309212	40.0	35.0	
118 Naphthalene	128	11.723	11.723	0.000	96	2415397	40.0	47.4	
119 1,2,3-Trichlorobenzene	180	11.960	11.960	0.000	94	732633	40.0	40.8	
S 133 Trihalomethanes, Total	1				0		160.0	149.7	
S 159 Total BTEX	1				0		200.0	202.8	
S 130 1,2-Dichloroethene, Total	96				0			72.6	
S 131 1,3-Dichloropropene, Total	75				0			83.4	
S 132 Xylenes, Total	106				0		80.0	83.9	

Reagents:

VMRGAS_00281	Amount Added: 32.00	Units: uL
vm50ss_stk_00079	Amount Added: 32.00	Units: uL
vm50is_stk_A_00002	Amount Added: 2.00	Units: uL
VMRPRIMW_00320	Amount Added: 32.00	Units: uL
VMAROLISTDW_00283	Amount Added: 32.00	Units: uL

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8021.D

Injection Date: 07-Feb-2019 13:09:30

Instrument ID: A3UX11

Operator ID: 43582

Lims ID: STD8260 L6

Worklist Smp#: 2

Client ID:

Purge Vol: 5.000 mL

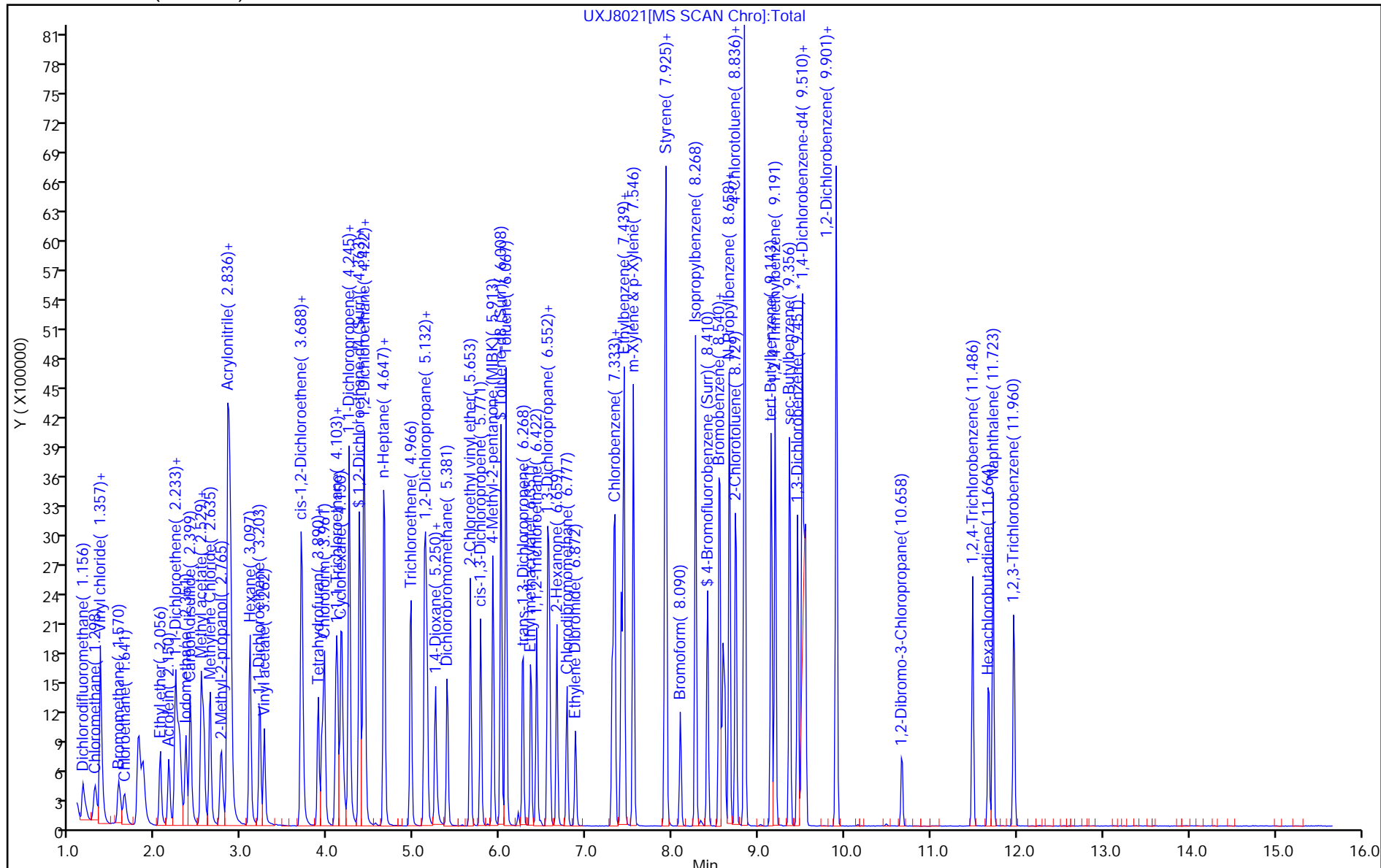
Dil. Factor: 1.0000

ALS Bottle#: 1

Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8022.D
 Lims ID: STD8260 L5
 Client ID:
 Sample Type: IC Calib Level: 5
 Inject. Date: 07-Feb-2019 13:31:30 ALS Bottle#: 2 Worklist Smp#: 3
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0084305-003
 Misc. Info.: J90207A-IC,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Sublist: chrom-8260_11*sub78
 Method: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 08-Feb-2019 09:14:37 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0324

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	4.659	4.659	0.000	97	1237419	20.0	20.0	
* 2 Chlorobenzene-d5	117	7.309	7.309	0.000	82	789116	20.0	20.0	
* 3 1,4-Dichlorobenzene-d4	152	9.522	9.522	0.000	94	424336	20.0	20.0	
\$ 4 Dibromofluoromethane (Surr	113	4.091	4.103	-0.012	97	298829	20.0	20.7	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.375	4.375	0.000	95	331727	20.0	19.0	
\$ 6 Toluene-d8 (Surr)	98	6.008	6.008	0.000	92	1002587	20.0	21.4	
\$ 7 4-Bromofluorobenzene (Surr	95	8.410	8.410	0.000	87	297359	20.0	19.9	
9 Dichlorodifluoromethane	85	1.168	1.168	0.000	99	393607	20.0	21.5	
10 Chloromethane	50	1.322	1.310	0.012	99	281547	20.0	18.7	
13 Butadiene	54	1.369	1.369	0.000	97	298381	20.0	21.2	
12 Vinyl chloride	62	1.393	1.393	0.000	98	339988	20.0	20.8	
15 Bromomethane	94	1.594	1.606	-0.012	90	220958	20.0	20.4	
16 Chloroethane	64	1.653	1.665	-0.012	99	188887	20.0	21.9	
17 Dichlorofluoromethane	67	1.807	1.807	0.000	99	434053	20.0	20.0	
18 Trichlorofluoromethane	101	1.830	1.831	-0.001	99	488108	20.0	21.2	
19 Ethyl ether	59	2.067	2.067	0.000	89	240821	20.0	20.9	
20 Acrolein	56	2.162	2.162	0.000	99	311245	100.0	105.3	
21 1,1-Dichloroethene	96	2.245	2.245	0.000	98	296326	20.0	20.8	
22 1,1,2-Trichloro-1,2,2-trif	151	2.280	2.280	0.000	90	246455	20.0	20.6	
23 Acetone	43	2.292	2.292	0.000	99	249848	40.0	37.7	
25 Iodomethane	142	2.363	2.363	0.000	98	515391	20.0	20.4	
26 Carbon disulfide	76	2.410	2.411	0.000	99	914051	20.0	19.6	
28 3-Chloro-1-propene	76	2.540	2.541	-0.001	88	224690	20.0	21.4	
29 Methyl acetate	43	2.564	2.564	0.000	97	589687	40.0	41.6	
30 Methylene Chloride	84	2.635	2.647	-0.012	87	336958	20.0	21.9	
31 2-Methyl-2-propanol	59	2.753	2.754	-0.001	95	489028	200.0	204.2	
32 Acrylonitrile	53	2.848	2.848	0.000	99	1538522	200.0	207.7	
34 trans-1,2-Dichloroethene	96	2.872	2.872	0.000	98	351024	20.0	20.8	
33 Methyl tert-butyl ether	73	2.884	2.884	0.000	96	780350	20.0	20.7	
35 Hexane	86	3.108	3.109	-0.001	91	87974	20.0	20.8	
36 1,1-Dichloroethane	63	3.215	3.215	0.000	96	577900	20.0	20.9	

Compound	Sig	RT (min.)	Exp RT (min.)	Diff RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
37 Vinyl acetate	43	3.262	3.274	-0.012	97	546119	20.0	21.0	
43 2,2-Dichloropropane	97	3.700	3.700	0.000	61	83030	20.0	20.3	
41 cis-1,2-Dichloroethene	96	3.700	3.700	0.000	81	373545	20.0	21.2	
42 2-Butanone (MEK)	43	3.712	3.712	0.000	98	328459	40.0	39.3	
47 Chlorobromomethane	128	3.889	3.890	-0.001	92	180067	20.0	20.0	
48 Tetrahydrofuran	42	3.937	3.937	0.000	85	232324	40.0	41.4	
49 Chloroform	83	3.960	3.961	-0.001	94	552690	20.0	20.6	
50 1,1,1-Trichloroethane	97	4.114	4.114	0.000	98	475687	20.0	20.4	
51 Cyclohexane	56	4.162	4.162	0.000	87	486587	20.0	22.4	
52 1,1-Dichloropropene	75	4.256	4.256	0.000	96	451488	20.0	21.1	
53 Carbon tetrachloride	117	4.256	4.256	0.000	91	455300	20.0	20.6	
54 Isobutyl alcohol	41	4.363	4.363	0.000	94	411928	500.0	531.2	
55 Benzene	78	4.422	4.422	0.000	95	1254314	20.0	20.3	
56 1,2-Dichloroethane	62	4.434	4.434	0.000	98	425614	20.0	19.8	
58 n-Heptane	100	4.659	4.659	0.000	53	76210	20.0	21.2	
60 Trichloroethene	130	4.966	4.966	0.000	94	362851	20.0	19.7	
62 Methylcyclohexane	83	5.132	5.132	0.000	89	531548	20.0	22.1	
63 1,2-Dichloropropane	63	5.156	5.156	0.000	94	257754	20.0	19.8	
65 Dibromomethane	93	5.250	5.250	0.000	90	184789	20.0	20.3	
66 1,4-Dioxane	88	5.274	5.274	0.000	89	81643	400.0	377.5	
67 Dichlorobromomethane	83	5.392	5.392	0.000	99	382766	20.0	19.3	
69 2-Chloroethyl vinyl ether	63	5.653	5.653	0.000	92	307879	40.0	38.7	
70 cis-1,3-Dichloropropene	75	5.771	5.771	0.000	96	409067	20.0	19.9	
71 4-Methyl-2-pentanone (MIBK)	43	5.913	5.913	0.000	96	613490	40.0	41.2	
72 Toluene	91	6.067	6.067	0.000	99	1170735	20.0	20.5	
73 trans-1,3-Dichloropropene	75	6.268	6.268	0.000	91	355082	20.0	19.1	
74 Ethyl methacrylate	69	6.363	6.363	-0.001	88	326308	20.0	21.1	
75 1,1,2-Trichloroethane	97	6.422	6.422	0.000	91	221396	20.0	18.5	
76 Tetrachloroethene	164	6.552	6.552	0.000	97	254413	20.0	20.0	
77 1,3-Dichloropropane	76	6.575	6.576	-0.001	89	376403	20.0	19.6	
78 2-Hexanone	43	6.658	6.659	-0.001	97	426962	40.0	41.8	
80 Chlorodibromomethane	129	6.777	6.777	0.000	89	281679	20.0	20.3	
82 Ethylene Dibromide	107	6.871	6.872	-0.001	96	245052	20.0	19.9	
84 Chlorobenzene	112	7.333	7.333	0.000	97	750415	20.0	19.6	
85 1,1,1,2-Tetrachloroethane	131	7.404	7.404	0.000	96	302221	20.0	20.6	
86 Ethylbenzene	106	7.439	7.439	0.000	98	401219	20.0	20.9	
87 m-Xylene & p-Xylene	106	7.546	7.546	0.000	98	490117	20.0	21.9	
88 o-Xylene	106	7.913	7.913	0.000	96	478708	20.0	22.1	
89 Styrene	104	7.924	7.925	-0.001	93	814512	20.0	22.3	
90 Bromoform	173	8.090	8.090	0.000	98	225037	20.0	20.3	
91 Isopropylbenzene	105	8.268	8.268	0.000	95	1249226	20.0	22.4	
95 Bromobenzene	156	8.552	8.540	0.012	95	318510	20.0	20.0	
94 1,1,2,2-Tetrachloroethane	83	8.552	8.552	0.000	96	347666	20.0	20.2	
96 1,2,3-Trichloropropane	110	8.587	8.587	0.000	85	116642	20.0	19.5	
97 trans-1,4-Dichloro-2-buten	53	8.611	8.611	0.000	80	107185	20.0	19.5	
98 N-Propylbenzene	120	8.658	8.658	0.000	99	355179	20.0	22.0	
99 2-Chlorotoluene	126	8.741	8.741	0.000	98	310087	20.0	21.5	
100 1,3,5-Trimethylbenzene	105	8.836	8.836	0.000	90	961736	20.0	21.9	
101 4-Chlorotoluene	126	8.836	8.848	-0.012	94	309733	20.0	20.9	
102 tert-Butylbenzene	119	9.143	9.155	-0.012	91	833511	20.0	22.2	
104 1,2,4-Trimethylbenzene	105	9.191	9.191	0.000	96	983336	20.0	22.2	
105 sec-Butylbenzene	105	9.356	9.356	0.000	93	1190193	20.0	22.3	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
106 1,3-Dichlorobenzene	146	9.451	9.451	0.000	97	585237	20.0	20.2	
107 4-Isopropyltoluene	119	9.510	9.510	0.000	97	1063145	20.0	22.5	
108 1,4-Dichlorobenzene	146	9.546	9.546	0.000	96	589475	20.0	19.0	
112 1,2-Dichlorobenzene	146	9.901	9.901	0.000	97	564881	20.0	20.9	
111 n-Butylbenzene	91	9.901	9.901	0.000	97	849937	20.0	22.0	
114 1,2-Dibromo-3-Chloropropan	157	10.670	10.670	0.000	90	105103	20.0	20.8	
116 1,2,4-Trichlorobenzene	180	11.486	11.486	0.000	93	363972	20.0	21.4	
117 Hexachlorobutadiene	225	11.675	11.676	-0.001	95	156722	20.0	19.9	
118 Naphthalene	128	11.723	11.723	0.000	96	1071425	20.0	23.7	
119 1,2,3-Trichlorobenzene	180	11.971	11.960	0.011	95	358175	20.0	22.4	
S 159 Total BTEX	1				0		100.0	105.8	
S 133 Trihalomethanes, Total	1				0		80.0	80.5	
S 130 1,2-Dichloroethene, Total	96				0			42.0	
S 131 1,3-Dichloropropene, Total	75				0			39.0	
S 132 Xylenes, Total	106				0		40.0	44.1	

Reagents:

VMRGAS_00281	Amount Added: 16.00	Units: uL
vm50ss_stk_00079	Amount Added: 16.00	Units: uL
vm50is_stk_A_00002	Amount Added: 2.00	Units: uL
VMRPRIMW_00320	Amount Added: 16.00	Units: uL
VMAROLISTDW_00283	Amount Added: 16.00	Units: uL



TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8022.D

Injection Date: 07-Feb-2019 13:31:30

Instrument ID: A3UX11

Operator ID: 43582

Lims ID: STD8260 L5

Worklist Smp#: 3

Client ID:

Purge Vol: 5.000 mL

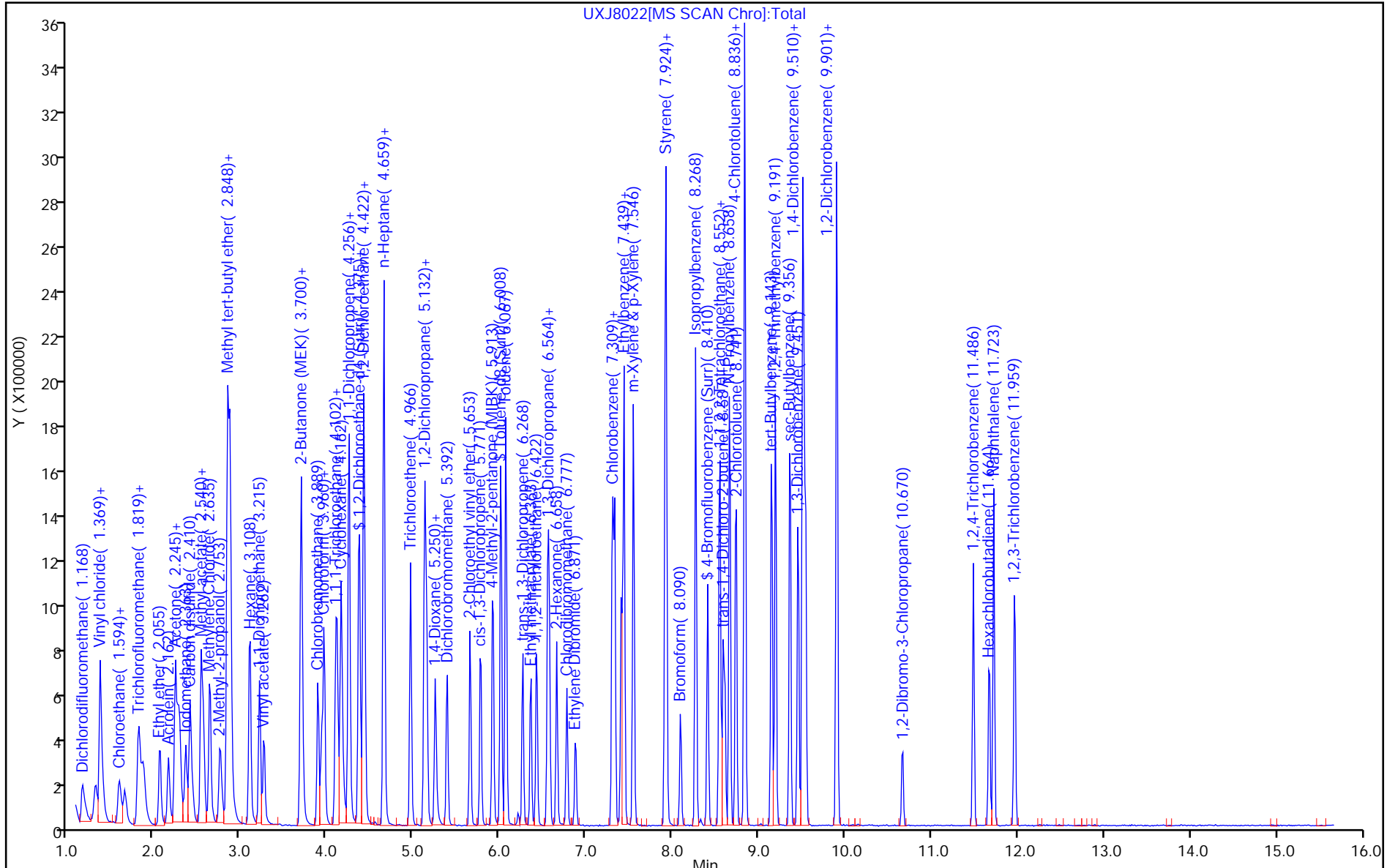
Dil. Factor: 1.0000

ALS Bottle#: 2

Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8023.D
 Lims ID: STD8260 L4
 Client ID:
 Sample Type: ICIS Calib Level: 4
 Inject. Date: 07-Feb-2019 13:53:30 ALS Bottle#: 3 Worklist Smp#: 4
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0084305-004
 Misc. Info.: J90207A-IC,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Sublist: chrom-8260_11*sub78
 Method: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 08-Feb-2019 09:14:43 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0324

First Level Reviewer: evansle

Date: 08-Feb-2019 07:55:40

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	4.659	4.659	0.000	99	1172070	20.0	20.0	
* 2 Chlorobenzene-d5	117	7.309	7.309	0.000	77	751566	20.0	20.0	
* 3 1,4-Dichlorobenzene-d4	152	9.522	9.522	0.000	92	416048	20.0	20.0	
\$ 4 Dibromofluoromethane (Surr	113	4.103	4.103	0.000	93	141317	10.0	10.3	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.375	4.375	0.000	74	163525	10.0	9.88	
\$ 6 Toluene-d8 (Surr)	98	6.008	6.008	0.000	81	459092	10.0	10.3	
\$ 7 4-Bromofluorobenzene (Surr	95	8.410	8.410	0.000	86	146668	10.0	10.3	
9 Dichlorodifluoromethane	85	1.168	1.168	0.000	96	166318	10.0	9.57	
10 Chloromethane	50	1.310	1.310	0.000	69	127316	10.0	8.92	
13 Butadiene	54	1.369	1.369	0.000	92	116135	10.0	8.72	
12 Vinyl chloride	62	1.393	1.393	0.000	81	144260	10.0	9.30	
15 Bromomethane	94	1.606	1.606	0.000	84	101367	10.0	9.89	
16 Chloroethane	64	1.665	1.665	0.000	85	83470	10.0	10.2	
17 Dichlorofluoromethane	67	1.807	1.807	0.000	95	200231	10.0	9.72	
18 Trichlorofluoromethane	101	1.831	1.831	0.000	86	201892	10.0	9.26	
19 Ethyl ether	59	2.067	2.067	0.000	85	111089	10.0	10.2	
20 Acrolein	56	2.162	2.162	0.000	88	139233	50.0	49.7	
21 1,1-Dichloroethene	96	2.245	2.245	0.000	93	136302	10.0	10.1	
22 1,1,2-Trichloro-1,2,2-trif	151	2.280	2.280	0.000	84	120721	10.0	10.7	
23 Acetone	43	2.292	2.292	0.000	94	110371	20.0	17.6	
25 Iodomethane	142	2.363	2.363	0.000	95	246503	10.0	10.3	
26 Carbon disulfide	76	2.411	2.411	0.000	98	424295	10.0	9.61	
28 3-Chloro-1-propene	76	2.541	2.541	0.000	79	98437	10.0	9.91	
29 Methyl acetate	43	2.564	2.564	0.000	97	266658	20.0	19.9	
30 Methylene Chloride	84	2.647	2.647	0.000	79	159040	10.0	10.3	
31 2-Methyl-2-propanol	59	2.754	2.754	0.000	88	212100	100.0	93.5	
32 Acrylonitrile	53	2.848	2.848	0.000	100	686402	100.0	97.9	
34 trans-1,2-Dichloroethene	96	2.872	2.872	0.000	92	157177	10.0	9.81	
33 Methyl tert-butyl ether	73	2.884	2.884	0.000	89	363372	10.0	10.2	
35 Hexane	86	3.109	3.109	0.000	90	39284	10.0	10.2	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
36 1,1-Dichloroethane	63	3.215	3.215	0.000	96	264327	10.0	10.1	
37 Vinyl acetate	43	3.274	3.274	0.000	95	244809	10.0	9.92	
41 cis-1,2-Dichloroethene	96	3.700	3.700	0.000	81	171580	10.0	10.3	
43 2,2-Dichloropropane	97	3.700	3.700	0.000	60	37577	10.0	9.71	
42 2-Butanone (MEK)	43	3.712	3.712	0.000	98	147163	20.0	18.6	
47 Chlorobromomethane	128	3.890	3.890	0.000	72	83844	10.0	9.82	
48 Tetrahydrofuran	42	3.937	3.937	0.000	84	100847	20.0	19.0	
49 Chloroform	83	3.961	3.961	0.000	93	263272	10.0	10.4	
50 1,1,1-Trichloroethane	97	4.114	4.114	0.000	56	221456	10.0	10.0	
51 Cyclohexane	56	4.162	4.162	0.000	87	221283	10.0	10.8	
53 Carbon tetrachloride	117	4.256	4.256	0.000	80	207777	10.0	9.93	
52 1,1-Dichloropropene	75	4.256	4.256	0.000	93	213777	10.0	10.6	
54 Isobutyl alcohol	41	4.363	4.363	0.000	91	178315	250.0	241.4	
55 Benzene	78	4.422	4.422	0.000	90	590944	10.0	10.1	
56 1,2-Dichloroethane	62	4.434	4.434	0.000	74	208883	10.0	10.2	
58 n-Heptane	100	4.659	4.659	0.000	44	33582	10.0	9.94	
60 Trichloroethene	130	4.966	4.966	0.000	90	168836	10.0	9.68	
62 Methylcyclohexane	83	5.132	5.132	0.000	87	232379	10.0	10.2	
63 1,2-Dichloropropane	63	5.156	5.156	0.000	91	127764	10.0	10.4	
65 Dibromomethane	93	5.250	5.250	0.000	82	86362	10.0	10.0	
66 1,4-Dioxane	88	5.274	5.274	0.000	91	33357	200.0	169.0	
67 Dichlorobromomethane	83	5.392	5.392	0.000	98	184018	10.0	9.81	
69 2-Chloroethyl vinyl ether	63	5.653	5.653	0.000	90	138767	20.0	18.4	
70 cis-1,3-Dichloropropene	75	5.771	5.771	0.000	86	193425	10.0	9.93	
71 4-Methyl-2-pentanone (MIBK)	43	5.913	5.913	0.000	91	281355	20.0	19.9	
72 Toluene	91	6.067	6.067	0.000	91	543872	10.0	10.0	
73 trans-1,3-Dichloropropene	75	6.268	6.268	0.000	90	167511	10.0	9.48	
74 Ethyl methacrylate	69	6.363	6.363	0.000	84	149850	10.0	10.2	
75 1,1,2-Trichloroethane	97	6.422	6.422	0.000	89	110808	10.0	9.72	
76 Tetrachloroethene	164	6.552	6.552	0.000	97	122571	10.0	10.1	
77 1,3-Dichloropropane	76	6.576	6.576	0.000	87	181024	10.0	9.92	
78 2-Hexanone	43	6.659	6.659	0.000	94	191478	20.0	19.7	
80 Chlorodibromomethane	129	6.777	6.777	0.000	89	135325	10.0	10.2	
82 Ethylene Dibromide	107	6.872	6.872	0.000	59	115065	10.0	9.80	
84 Chlorobenzene	112	7.333	7.333	0.000	97	354543	10.0	9.75	
85 1,1,1,2-Tetrachloroethane	131	7.404	7.404	0.000	94	137694	10.0	9.87	
86 Ethylbenzene	106	7.439	7.439	0.000	98	177149	10.0	9.70	
87 m-Xylene & p-Xylene	106	7.546	7.546	0.000	97	225000	10.0	10.6	
88 o-Xylene	106	7.913	7.913	0.000	96	221095	10.0	10.7	
89 Styrene	104	7.925	7.925	0.000	91	366782	10.0	10.6	
90 Bromoform	173	8.090	8.090	0.000	95	107418	10.0	10.2	
91 Isopropylbenzene	105	8.268	8.268	0.000	95	554538	10.0	10.5	
95 Bromobenzene	156	8.540	8.540	0.000	81	149296	10.0	9.58	
94 1,1,2,2-Tetrachloroethane	83	8.552	8.552	0.000	96	163051	10.0	9.66	
96 1,2,3-Trichloropropane	110	8.587	8.587	0.000	53	56270	10.0	9.61	
97 trans-1,4-Dichloro-2-buten	53	8.611	8.611	0.000	80	49221	10.0	9.14	
98 N-Propylbenzene	120	8.658	8.658	0.000	92	163725	10.0	10.4	
99 2-Chlorotoluene	126	8.741	8.741	0.000	95	146693	10.0	10.4	
100 1,3,5-Trimethylbenzene	105	8.836	8.836	0.000	88	453276	10.0	10.5	
101 4-Chlorotoluene	126	8.848	8.848	0.000	93	146255	10.0	10.1	
102 tert-Butylbenzene	119	9.155	9.155	0.000	91	369118	10.0	10.0	a
104 1,2,4-Trimethylbenzene	105	9.191	9.191	0.000	70	456294	10.0	10.5	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
105 sec-Butylbenzene	105	9.356	9.356	0.000	89	517546	10.0	9.88	
106 1,3-Dichlorobenzene	146	9.451	9.451	0.000	78	281323	10.0	9.92	
107 4-Isopropyltoluene	119	9.510	9.510	0.000	93	481868	10.0	10.4	
108 1,4-Dichlorobenzene	146	9.546	9.546	0.000	96	287252	10.0	9.44	
111 n-Butylbenzene	91	9.901	9.901	0.000	96	374733	10.0	9.91	
112 1,2-Dichlorobenzene	146	9.901	9.901	0.000	85	255531	10.0	9.64	
114 1,2-Dibromo-3-Chloropropan	157	10.670	10.670	0.000	81	50852	10.0	10.1	
116 1,2,4-Trichlorobenzene	180	11.486	11.486	0.000	93	153092	10.0	9.18	
117 Hexachlorobutadiene	225	11.676	11.676	0.000	80	73104	10.0	9.49	
118 Naphthalene	128	11.723	11.723	0.000	96	403718	10.0	9.09	
119 1,2,3-Trichlorobenzene	180	11.960	11.960	0.000	92	146668	10.0	9.37	
S 133 Trihalomethanes, Total	1				0		40.0	40.6	
S 159 Total BTEX	1				0		50.0	51.1	
S 132 Xylenes, Total	106				0		20.0	21.3	

QC Flag Legend

Review Flags

a - User Assigned ID

Reagents:

VMRGAS_00281	Amount Added: 8.00	Units: uL
vm50ss_stk_00079	Amount Added: 8.00	Units: uL
vm50is_stk_A_00002	Amount Added: 2.00	Units: uL
VMRPRIMW_00320	Amount Added: 8.00	Units: uL
VMAROLISTDW_00283	Amount Added: 8.00	Units: uL

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8023.D

Injection Date: 07-Feb-2019 13:53:30

Instrument ID: A3UX11

Operator ID: 43582

Lims ID: STD8260 L4

Worklist Smp#: 4

Client ID:

Purge Vol: 5.000 mL

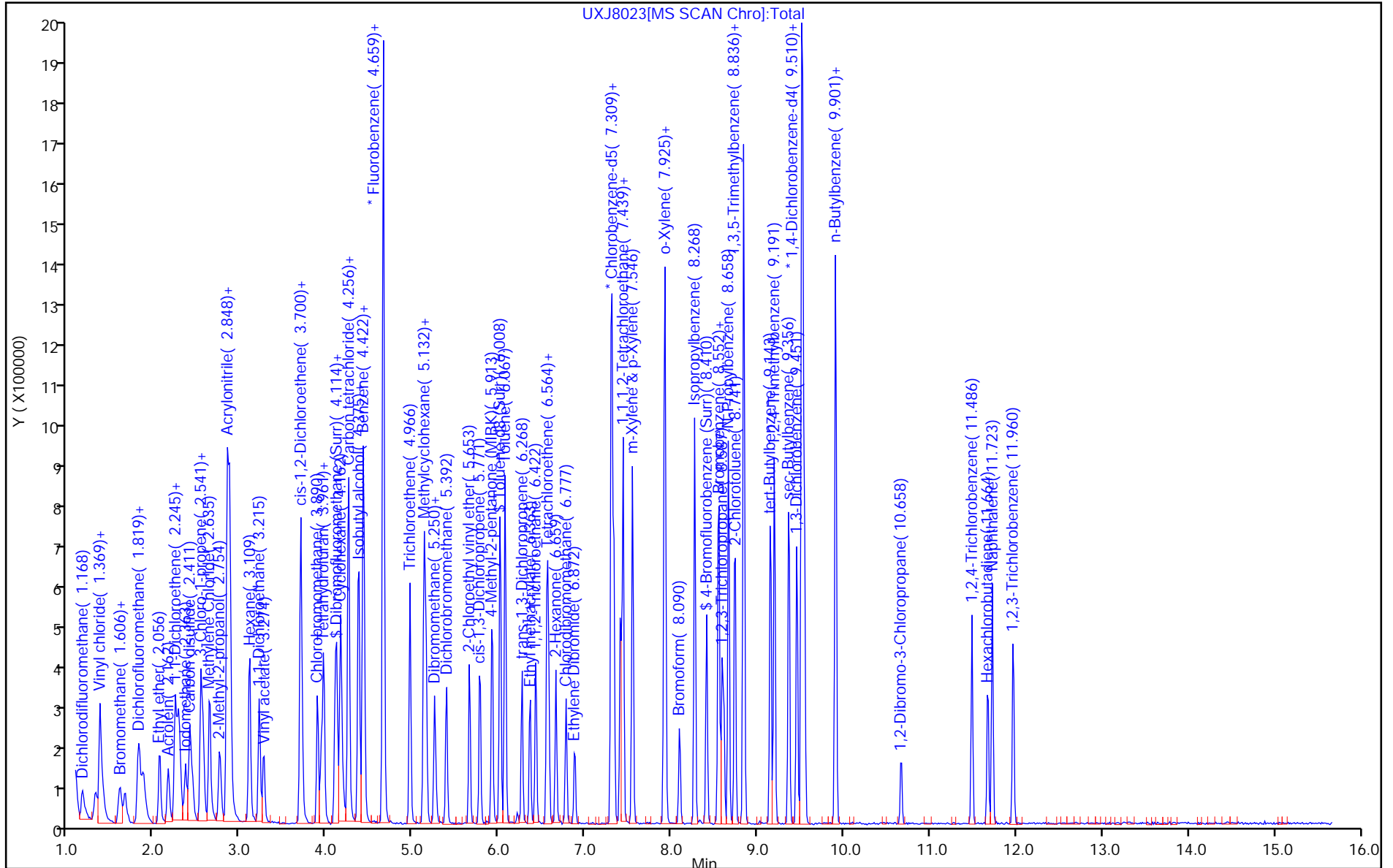
Dil. Factor: 1.0000

ALS Bottle#: 3

Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton

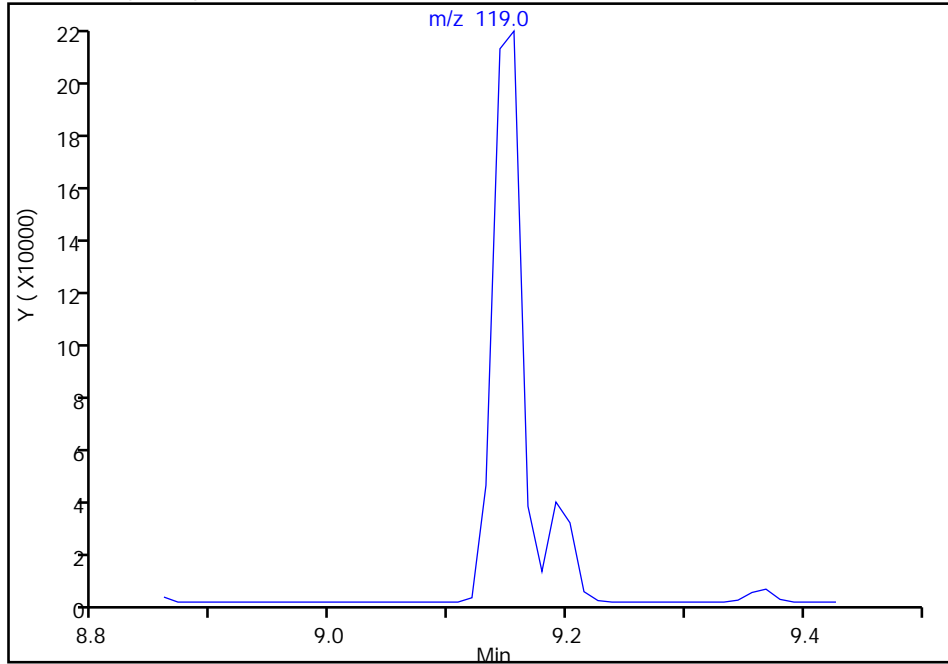
Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8023.D
Injection Date: 07-Feb-2019 13:53:30 Instrument ID: A3UX11
Lims ID: STD8260 L4
Client ID:
Operator ID: 43582 ALS Bottle#: 3 Worklist Smp#: 4
Purge Vol: 5.000 mL Dil. Factor: 1.0000
Method: 8260_11 Limit Group: MSV 8260B ICAL
Column: DB-624 (0.18 mm) Detector MS SCAN

102 tert-Butylbenzene, CAS: 98-06-6

Signal: 1

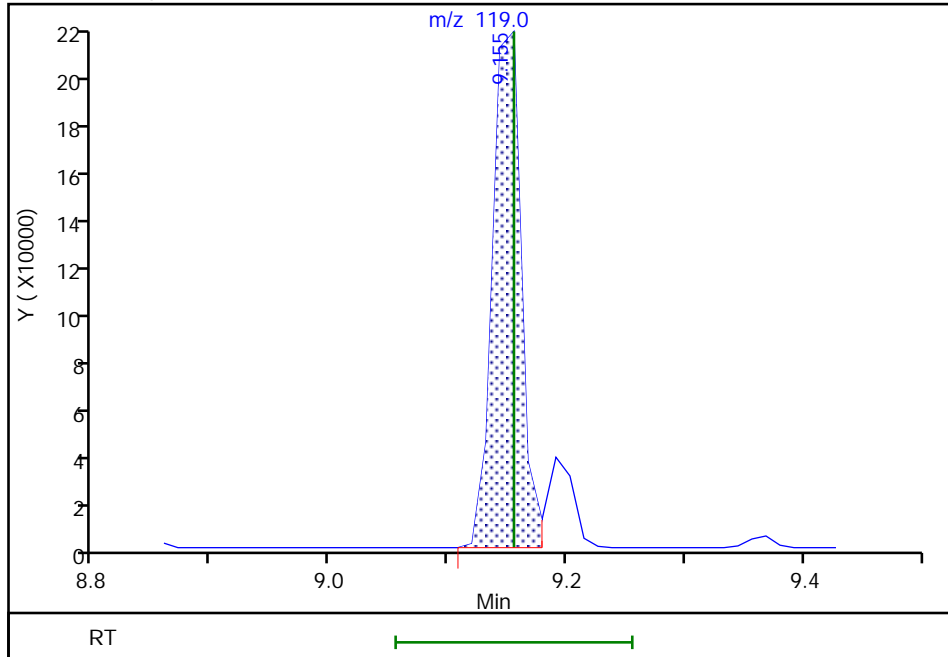
Not Detected
Expected RT: 9.16

Processing Integration Results



Manual Integration Results

RT: 9.16
Area: 369118
Amount: 10.032482
Amount Units: ug/l



Reviewer: evansle, 08-Feb-2019 08:24:18

Audit Action: Assigned Compound ID

Audit Reason: Peak assignment corrected

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8024.D
 Lims ID: STD8260 L3
 Client ID:
 Sample Type: IC Calib Level: 3
 Inject. Date: 07-Feb-2019 14:15:30 ALS Bottle#: 4 Worklist Smp#: 5
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0084305-005
 Misc. Info.: J90207A-IC,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Sublist: chrom-8260_11*sub78
 Method: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 08-Feb-2019 09:14:48 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0324

First Level Reviewer: evansle

Date: 08-Feb-2019 08:27:27

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	4.659	4.659	0.000	99	1155721	20.0	20.0	
* 2 Chlorobenzene-d5	117	7.297	7.309	-0.012	83	792526	20.0	20.0	
* 3 1,4-Dichlorobenzene-d4	152	9.522	9.522	0.000	92	427773	20.0	20.0	
\$ 4 Dibromofluoromethane (Surr	113	4.091	4.103	-0.012	94	68214	5.00	5.06	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.375	4.375	0.000	97	85806	5.00	5.26	
\$ 6 Toluene-d8 (Surr)	98	6.007	6.008	-0.001	91	231027	5.00	4.91	
\$ 7 4-Bromofluorobenzene (Surr	95	8.410	8.410	0.000	88	79386	5.00	5.28	
9 Dichlorodifluoromethane	85	1.156	1.168	-0.012	98	91261	5.00	5.33	
10 Chloromethane	50	1.298	1.310	-0.012	99	74271	5.00	5.27	
13 Butadiene	54	1.357	1.369	-0.012	95	67238	5.00	5.12	
12 Vinyl chloride	62	1.381	1.393	-0.012	98	79953	5.00	5.23	
15 Bromomethane	94	1.582	1.606	-0.024	88	54178	5.00	5.36	
16 Chloroethane	64	1.641	1.665	-0.024	99	42315	5.00	5.25	
17 Dichlorofluoromethane	67	1.807	1.807	0.000	99	106239	5.00	5.23	
18 Trichlorofluoromethane	101	1.854	1.831	0.023	90	116801	5.00	5.43	
19 Ethyl ether	59	2.055	2.067	-0.012	87	53290	5.00	4.96	
20 Acrolein	56	2.162	2.162	0.000	99	76307	25.0	27.6	
21 1,1-Dichloroethene	96	2.233	2.245	-0.012	98	68669	5.00	5.17	
22 1,1,2-Trichloro-1,2,2-trif	151	2.268	2.280	-0.012	88	58077	5.00	5.21	
23 Acetone	43	2.280	2.292	-0.012	98	61588	10.0	9.94	
25 Iodomethane	142	2.351	2.363	-0.012	97	118663	5.00	5.03	
26 Carbon disulfide	76	2.410	2.411	0.000	99	207524	5.00	4.77	
28 3-Chloro-1-propene	76	2.540	2.541	-0.001	89	48182	5.00	4.92	
29 Methyl acetate	43	2.564	2.564	0.000	97	142070	10.0	10.7	
30 Methylene Chloride	84	2.635	2.647	-0.012	89	85471	5.00	5.03	
31 2-Methyl-2-propanol	59	2.753	2.754	-0.001	94	108992	50.0	48.7	
32 Acrylonitrile	53	2.836	2.848	-0.012	98	360130	50.0	52.1	
34 trans-1,2-Dichloroethene	96	2.872	2.872	0.000	97	78810	5.00	4.99	
33 Methyl tert-butyl ether	73	2.872	2.884	-0.012	96	179324	5.00	5.08	
35 Hexane	86	3.097	3.109	-0.012	92	18570	5.00	5.24	

Compound	Sig	RT (min.)	Exp RT (min.)	Diff RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
36 1,1-Dichloroethane	63	3.215	3.215	0.000	96	123928	5.00	4.79	
37 Vinyl acetate	43	3.262	3.274	-0.012	98	122657	5.00	5.04	
43 2,2-Dichloropropane	97	3.688	3.700	-0.012	62	18470	5.00	4.84	
41 cis-1,2-Dichloroethene	96	3.700	3.700	0.000	82	79034	5.00	4.81	
42 2-Butanone (MEK)	43	3.712	3.712	0.000	99	89550	10.0	11.5	
47 Chlorobromomethane	128	3.889	3.890	-0.001	91	41697	5.00	4.95	
48 Tetrahydrofuran	42	3.937	3.937	0.000	86	54325	10.0	10.4	
49 Chloroform	83	3.960	3.961	-0.001	94	130678	5.00	5.21	
50 1,1,1-Trichloroethane	97	4.114	4.114	0.000	99	103539	5.00	4.75	
51 Cyclohexane	56	4.162	4.162	0.000	87	102175	5.00	5.04	
52 1,1-Dichloropropene	75	4.244	4.256	-0.012	94	95002	5.00	4.76	
53 Carbon tetrachloride	117	4.256	4.256	0.000	97	104922	5.00	5.08	
54 Isobutyl alcohol	41	4.363	4.363	0.000	92	100691	125.0	129.3	
55 Benzene	78	4.422	4.422	0.000	95	283539	5.00	4.90	
56 1,2-Dichloroethane	62	4.434	4.434	0.000	98	99528	5.00	4.95	
58 n-Heptane	100	4.659	4.659	0.000	40	19161	5.00	5.82	
60 Trichloroethene	130	4.966	4.966	0.000	97	85827	5.00	4.99	
62 Methylcyclohexane	83	5.132	5.132	0.000	88	106738	5.00	4.75	
63 1,2-Dichloropropane	63	5.156	5.156	0.000	95	60714	5.00	5.00	
65 Dibromomethane	93	5.250	5.250	0.000	91	42824	5.00	5.02	
66 1,4-Dioxane	88	5.274	5.274	0.000	87	20834	100.0	111.1	
67 Dichlorobromomethane	83	5.380	5.392	-0.012	99	95815	5.00	5.18	
69 2-Chloroethyl vinyl ether	63	5.653	5.653	-0.001	92	82558	10.0	11.1	
70 cis-1,3-Dichloropropene	75	5.771	5.771	0.000	96	99988	5.00	5.21	
71 4-Methyl-2-pentanone (MIBK)	43	5.913	5.913	0.000	95	153307	10.0	11.0	
72 Toluene	91	6.067	6.067	0.000	99	290746	5.00	5.08	
73 trans-1,3-Dichloropropene	75	6.268	6.268	0.000	91	94776	5.00	5.09	
74 Ethyl methacrylate	69	6.362	6.363	-0.001	86	78145	5.00	5.04	
75 1,1,2-Trichloroethane	97	6.422	6.422	0.000	90	60454	5.00	5.03	
76 Tetrachloroethene	164	6.552	6.552	0.000	98	67708	5.00	5.29	
77 1,3-Dichloropropane	76	6.575	6.576	-0.001	90	98029	5.00	5.09	
78 2-Hexanone	43	6.658	6.659	-0.001	93	109052	10.0	10.6	
80 Chlorodibromomethane	129	6.777	6.777	0.000	88	73418	5.00	5.26	
82 Ethylene Dibromide	107	6.871	6.872	-0.001	97	60552	5.00	4.89	
84 Chlorobenzene	112	7.333	7.333	0.000	98	189707	5.00	4.95	
85 1,1,1,2-Tetrachloroethane	131	7.404	7.404	0.000	95	68426	5.00	4.65	
86 Ethylbenzene	106	7.439	7.439	0.000	97	99260	5.00	5.16	
87 m-Xylene & p-Xylene	106	7.546	7.546	0.000	95	114642	5.00	5.11	
88 o-Xylene	106	7.913	7.913	0.000	96	105654	5.00	4.86	
89 Styrene	104	7.924	7.925	-0.001	93	184963	5.00	5.05	
90 Bromoform	173	8.090	8.090	0.000	96	58743	5.00	5.29	
91 Isopropylbenzene	105	8.268	8.268	0.000	95	276000	5.00	4.94	
95 Bromobenzene	156	8.540	8.540	0.000	94	77365	5.00	4.83	
94 1,1,2,2-Tetrachloroethane	83	8.552	8.552	0.000	94	87302	5.00	5.03	
96 1,2,3-Trichloropropane	110	8.587	8.587	0.000	85	30733	5.00	5.10	
97 trans-1,4-Dichloro-2-buten	53	8.611	8.611	0.000	80	31177	5.00	5.63	
98 N-Propylbenzene	120	8.658	8.658	0.000	98	79799	5.00	4.91	
99 2-Chlorotoluene	126	8.741	8.741	0.000	98	73878	5.00	5.08	
100 1,3,5-Trimethylbenzene	105	8.836	8.836	0.000	94	218959	5.00	4.95	
101 4-Chlorotoluene	126	8.847	8.848	-0.001	97	78564	5.00	5.27	
102 tert-Butylbenzene	119	9.143	9.155	-0.012	90	193232	5.00	5.11	
104 1,2,4-Trimethylbenzene	105	9.191	9.191	0.000	96	220805	5.00	4.95	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
105 sec-Butylbenzene	105	9.356	9.356	0.000	94	261531	5.00	4.86	
106 1,3-Dichlorobenzene	146	9.451	9.451	0.000	97	143622	5.00	4.93	
107 4-Isopropyltoluene	119	9.510	9.510	0.000	97	231439	5.00	4.85	
108 1,4-Dichlorobenzene	146	9.546	9.546	0.000	95	149185	5.00	4.77	
112 1,2-Dichlorobenzene	146	9.901	9.901	0.000	96	130837	5.00	4.80	
111 n-Butylbenzene	91	9.901	9.901	0.000	97	183328	5.00	4.71	
114 1,2-Dibromo-3-Chloropropan	157	10.670	10.670	0.000	91	26261	5.00	4.92	
116 1,2,4-Trichlorobenzene	180	11.486	11.486	0.000	92	81516	5.00	4.75	
117 Hexachlorobutadiene	225	11.675	11.676	-0.001	93	37204	5.00	4.70	
118 Naphthalene	128	11.723	11.723	0.000	96	194010	5.00	4.25	
119 1,2,3-Trichlorobenzene	180	11.959	11.960	-0.001	96	75944	5.00	4.72	
S 159 Total BTEX	1				0		25.0	25.1	
S 133 Trihalomethanes, Total	1				0		20.0	20.9	
S 130 1,2-Dichloroethene, Total	96				0			9.80	
S 131 1,3-Dichloropropene, Total	75				0			10.3	
S 132 Xylenes, Total	106				0		10.0	9.97	

Reagents:

VMRGAS_00281	Amount Added: 4.00	Units: uL
vm50ss_stk_00079	Amount Added: 4.00	Units: uL
vm50is_stk_A_00002	Amount Added: 2.00	Units: uL
VMRPRIMW_00320	Amount Added: 4.00	Units: uL
VMAROLISTDW_00283	Amount Added: 4.00	Units: uL

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8024.D

Injection Date: 07-Feb-2019 14:15:30

Instrument ID: A3UX11

Operator ID: 43582

Lims ID: STD8260 L3

Worklist Smp#: 5

Client ID:

Purge Vol: 5.000 mL

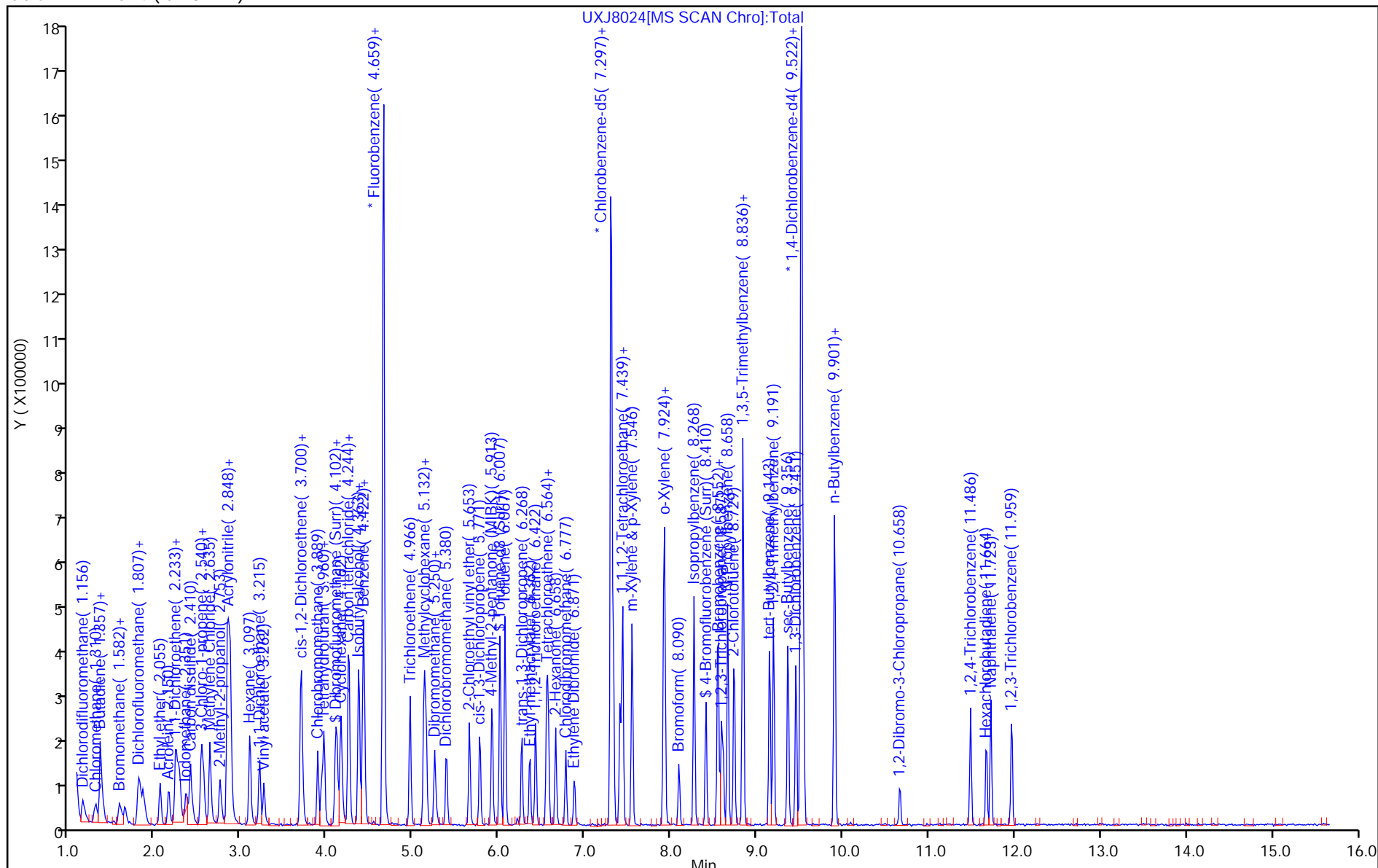
Dil. Factor: 1.0000

ALS Bottle#: 4

Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8025.D
 Lims ID: STD8260 L2
 Client ID:
 Sample Type: IC Calib Level: 2
 Inject. Date: 07-Feb-2019 14:38:30 ALS Bottle#: 5 Worklist Smp#: 6
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0084305-006
 Misc. Info.: J90207A-IC,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Sublist: chrom-8260_11*sub78
 Method: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 08-Feb-2019 09:14:54 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0324

First Level Reviewer: evansle

Date: 08-Feb-2019 08:07:56

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	4.659	4.659	0.000	99	1095260	20.0	20.0	
* 2 Chlorobenzene-d5	117	7.309	7.309	0.000	83	757263	20.0	20.0	
* 3 1,4-Dichlorobenzene-d4	152	9.522	9.522	0.000	93	405932	20.0	20.0	
\$ 4 Dibromofluoromethane (Surr	113	4.091	4.103	-0.012	97	24605	2.00	1.93	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.375	4.375	0.000	81	33449	2.00	2.16	
\$ 6 Toluene-d8 (Surr)	98	6.008	6.008	0.000	93	82074	2.00	1.83	
\$ 7 4-Bromofluorobenzene (Surr	95	8.410	8.410	0.000	86	28481	2.00	1.98	
9 Dichlorodifluoromethane	85	1.156	1.168	-0.012	94	33811	2.00	2.08	
10 Chloromethane	50	1.298	1.310	-0.012	99	29438	2.00	2.21	
13 Butadiene	54	1.357	1.369	-0.012	96	27131	2.00	2.18	
12 Vinyl chloride	62	1.381	1.393	-0.012	39	29665	2.00	2.05	
15 Bromomethane	94	1.582	1.606	-0.024	87	20704	2.00	2.16	
16 Chloroethane	64	1.641	1.665	-0.024	96	16382	2.00	2.14	
17 Dichlorofluoromethane	67	1.795	1.807	-0.012	97	41103	2.00	2.14	
18 Trichlorofluoromethane	101	1.819	1.831	-0.012	82	40004	2.00	1.96	
19 Ethyl ether	59	2.056	2.067	-0.011	86	20099	2.00	1.97	
20 Acrolein	56	2.150	2.162	-0.012	97	26117	10.0	9.98	
21 1,1-Dichloroethene	96	2.233	2.245	-0.012	96	25640	2.00	2.04	
22 1,1,2-Trichloro-1,2,2-trif	151	2.268	2.280	-0.012	89	21536	2.00	2.04	
23 Acetone	43	2.280	2.292	-0.012	97	27192	4.00	4.63	
25 Iodomethane	142	2.351	2.363	-0.012	97	44542	2.00	1.99	
26 Carbon disulfide	76	2.410	2.411	0.000	99	91902	2.00	2.23	
28 3-Chloro-1-propene	76	2.529	2.541	-0.012	89	18405	2.00	1.98	
29 Methyl acetate	43	2.552	2.564	-0.012	96	48220	4.00	3.84	
30 Methylene Chloride	84	2.635	2.647	-0.012	90	42445	2.00	2.04	
31 2-Methyl-2-propanol	59	2.754	2.754	0.000	96	42922	20.0	20.2	
32 Acrylonitrile	53	2.836	2.848	-0.012	96	133260	20.0	20.3	
34 trans-1,2-Dichloroethene	96	2.860	2.872	-0.012	96	30791	2.00	2.06	
33 Methyl tert-butyl ether	73	2.872	2.884	-0.012	93	69489	2.00	2.08	
35 Hexane	86	3.097	3.109	-0.012	84	3932	2.00	1.70	

Compound	Sig	RT (min.)	Exp RT (min.)	Diff RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
36 1,1-Dichloroethane	63	3.215	3.215	0.000	96	51015	2.00	2.08	
37 Vinyl acetate	43	3.262	3.274	-0.012	96	45507	2.00	1.97	
41 cis-1,2-Dichloroethene	96	3.688	3.700	-0.012	80	33131	2.00	2.13	
43 2,2-Dichloropropane	97	3.688	3.700	-0.012	60	7740	2.00	2.14	
42 2-Butanone (MEK)	43	3.712	3.712	0.000	99	31612	4.00	4.27	
47 Chlorobromomethane	128	3.890	3.890	0.000	89	17457	2.00	2.19	
48 Tetrahydrofuran	42	3.925	3.937	-0.012	77	19047	4.00	3.84	
49 Chloroform	83	3.961	3.961	0.000	93	48908	2.00	2.06	
50 1,1,1-Trichloroethane	97	4.114	4.114	0.000	97	42875	2.00	2.07	
51 Cyclohexane	56	4.162	4.162	0.000	89	36529	2.00	1.90	
53 Carbon tetrachloride	117	4.245	4.256	-0.011	95	38755	2.00	1.98	
52 1,1-Dichloropropene	75	4.245	4.256	-0.011	91	36750	2.00	1.94	
54 Isobutyl alcohol	41	4.363	4.363	0.000	90	35587	50.0	47.8	
55 Benzene	78	4.422	4.422	0.000	94	114875	2.00	2.10	
56 1,2-Dichloroethane	62	4.434	4.434	0.000	98	39111	2.00	2.05	
58 n-Heptane	100	4.647	4.659	-0.012	33	4317	2.00	1.50	
60 Trichloroethene	130	4.966	4.966	0.000	89	34086	2.00	2.09	
62 Methylcyclohexane	83	5.132	5.132	0.000	86	43436	2.00	2.04	
63 1,2-Dichloropropane	63	5.156	5.156	0.000	94	24147	2.00	2.10	
65 Dibromomethane	93	5.250	5.250	0.000	89	16623	2.00	2.06	
66 1,4-Dioxane	88	5.274	5.274	0.000	45	5017	40.0	36.4	
67 Dichlorobromomethane	83	5.381	5.392	-0.011	99	35418	2.00	2.02	
69 2-Chloroethyl vinyl ether	63	5.653	5.653	0.000	92	28161	4.00	3.99	
70 cis-1,3-Dichloropropene	75	5.771	5.771	0.000	96	35784	2.00	1.97	
71 4-Methyl-2-pentanone (MIBK)	43	5.913	5.913	0.000	94	48587	4.00	3.68	
72 Toluene	91	6.067	6.067	0.000	96	103151	2.00	1.89	
73 trans-1,3-Dichloropropene	75	6.268	6.268	0.000	91	36666	2.00	2.06	
74 Ethyl methacrylate	69	6.351	6.363	-0.012	87	28459	2.00	1.92	
75 1,1,2-Trichloroethane	97	6.422	6.422	0.000	84	22922	2.00	2.00	
76 Tetrachloroethene	164	6.552	6.552	0.000	95	23718	2.00	1.94	
77 1,3-Dichloropropane	76	6.576	6.576	0.000	90	36153	2.00	1.97	
78 2-Hexanone	43	6.659	6.659	-0.001	98	40322	4.00	4.11	
80 Chlorodibromomethane	129	6.777	6.777	0.000	85	28493	2.00	2.14	
82 Ethylene Dibromide	107	6.871	6.872	-0.001	95	21792	2.00	1.84	
84 Chlorobenzene	112	7.333	7.333	0.000	93	71386	2.00	1.95	
85 1,1,1,2-Tetrachloroethane	131	7.404	7.404	0.000	94	27415	2.00	1.95	
86 Ethylbenzene	106	7.439	7.439	0.000	99	33590	2.00	1.83	
87 m-Xylene & p-Xylene	106	7.546	7.546	0.000	98	37781	2.00	1.76	
88 o-Xylene	106	7.913	7.913	0.000	96	38511	2.00	1.86	
89 Styrene	104	7.925	7.925	0.000	93	62454	2.00	1.78	
90 Bromoform	173	8.090	8.090	0.000	95	21407	2.00	2.02	
91 Isopropylbenzene	105	8.268	8.268	0.000	95	93008	2.00	1.74	
95 Bromobenzene	156	8.540	8.540	0.000	91	30472	2.00	2.00	
94 1,1,2,2-Tetrachloroethane	83	8.552	8.552	0.000	92	32123	2.00	1.95	
96 1,2,3-Trichloropropane	110	8.587	8.587	0.000	84	12345	2.00	2.16	
97 trans-1,4-Dichloro-2-buten	53	8.611	8.611	0.000	67	10685	2.00	2.03	
98 N-Propylbenzene	120	8.658	8.658	0.000	99	26311	2.00	1.71	
99 2-Chlorotoluene	126	8.741	8.741	0.000	97	24540	2.00	1.78	
100 1,3,5-Trimethylbenzene	105	8.836	8.836	0.000	97	71649	2.00	1.71	
101 4-Chlorotoluene	126	8.836	8.848	-0.012	98	25871	2.00	1.83	
102 tert-Butylbenzene	119	9.143	9.155	-0.012	90	59904	2.00	1.67	
104 1,2,4-Trimethylbenzene	105	9.191	9.191	0.000	92	72413	2.00	1.71	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
105 sec-Butylbenzene	105	9.356	9.356	0.000	93	92291	2.00	1.81	
106 1,3-Dichlorobenzene	146	9.451	9.451	0.000	96	54925	2.00	1.99	
107 4-Isopropyltoluene	119	9.510	9.510	0.000	96	80813	2.00	1.79	
108 1,4-Dichlorobenzene	146	9.546	9.546	0.000	93	61678	2.00	2.08	
111 n-Butylbenzene	91	9.901	9.901	0.000	97	64537	2.00	1.75	
112 1,2-Dichlorobenzene	146	9.901	9.901	0.000	95	52725	2.00	2.04	
114 1,2-Dibromo-3-Chloropropan	157	10.658	10.670	-0.012	83	9092	2.00	1.61	
116 1,2,4-Trichlorobenzene	180	11.486	11.486	0.000	90	31765	2.00	1.95	
117 Hexachlorobutadiene	225	11.676	11.676	0.000	95	16252	2.00	2.16	
118 Naphthalene	128	11.723	11.723	0.000	95	79647	2.00	1.84	
119 1,2,3-Trichlorobenzene	180	11.971	11.960	0.011	90	28542	2.00	1.87	
S 133 Trihalomethanes, Total	1				0		8.00	8.23	
S 159 Total BTEX	1				0		10.0	9.42	
S 130 1,2-Dichloroethene, Total	96				0			4.18	
S 131 1,3-Dichloropropene, Total	75				0			4.03	
S 132 Xylenes, Total	106				0		4.00	3.62	

Reagents:

VMRGAS_00281	Amount Added: 1.60	Units: uL
vm50ss_stk_00079	Amount Added: 1.60	Units: uL
vm50is_stk_A_00002	Amount Added: 2.00	Units: uL
VMRPRIMW_00320	Amount Added: 1.60	Units: uL
VMAROLISTDW_00283	Amount Added: 1.60	Units: uL

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8025.D

Injection Date: 07-Feb-2019 14:38:30

Instrument ID: A3UX11

Operator ID: 43582

Lims ID: STD8260 L2

Worklist Smp#: 6

Client ID:

Purge Vol: 5.000 mL

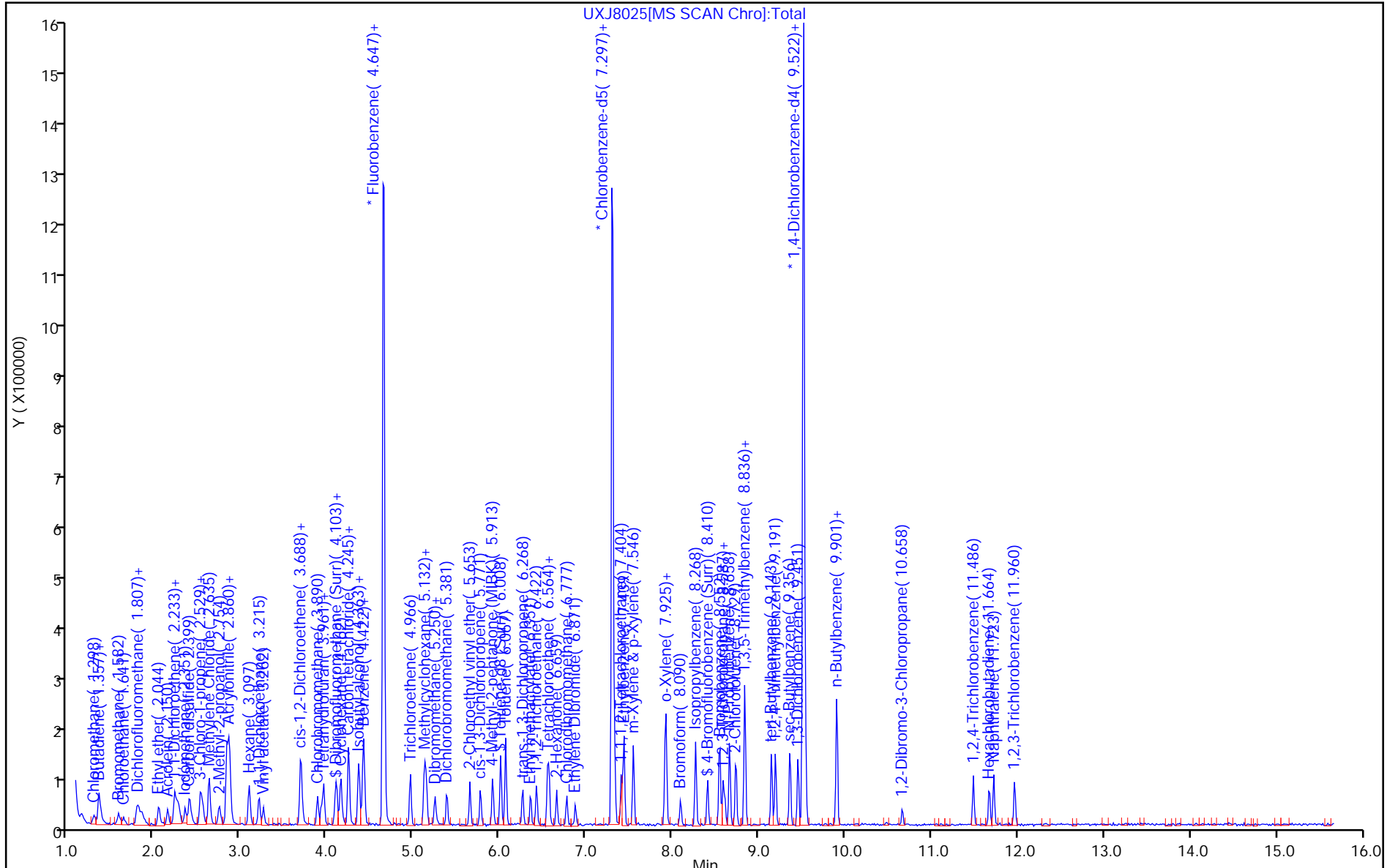
Dil. Factor: 1.0000

ALS Bottle#: 5

Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8026.D
 Lims ID: STD8260 L1
 Client ID:
 Sample Type: IC Calib Level: 1
 Inject. Date: 07-Feb-2019 15:00:30 ALS Bottle#: 6 Worklist Smp#: 7
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0084305-007
 Misc. Info.: J90207A-IC,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Sublist: chrom-8260_11*sub78
 Method: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 08-Feb-2019 09:15:00 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0324

First Level Reviewer: evansle

Date: 08-Feb-2019 08:08:43

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	4.659	4.659	0.000	99	1116865	20.0	20.0	
* 2 Chlorobenzene-d5	117	7.309	7.309	0.000	83	702853	20.0	20.0	
* 3 1,4-Dichlorobenzene-d4	152	9.522	9.522	0.000	92	374248	20.0	20.0	
\$ 4 Dibromofluoromethane (Surr	113	4.091	4.103	-0.012	94	13778	1.00	1.06	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.375	4.375	0.000	96	16708	1.00	1.06	
\$ 6 Toluene-d8 (Surr)	98	6.008	6.008	0.000	92	41706	1.00	1.00	
\$ 7 4-Bromofluorobenzene (Surr	95	8.410	8.410	0.000	89	13244	1.00	0.99	
9 Dichlorodifluoromethane	85	1.168	1.168	0.000	94	14302	1.00	0.8637	
10 Chloromethane	50	1.322	1.310	0.012	1	14735	1.00	1.08	a
13 Butadiene	54	1.369	1.369	0.000	92	12457	1.00	0.9813	
12 Vinyl chloride	62	1.393	1.393	0.000	93	15085	1.00	1.02	
15 Bromomethane	94	1.594	1.606	-0.012	85	9111	1.00	0.9329	
16 Chloroethane	64	1.653	1.665	-0.012	39	6225	1.00	0.7992	
17 Dichlorofluoromethane	67	1.807	1.807	0.000	95	20171	1.00	1.03	
18 Trichlorofluoromethane	101	1.854	1.831	0.023	84	20689	1.00	1.00	
19 Ethyl ether	59	2.055	2.067	-0.012	81	10777	1.00	1.04	
20 Acrolein	56	2.150	2.162	-0.012	98	12263	5.00	4.60	
21 1,1-Dichloroethene	96	2.245	2.245	0.000	94	12563	1.00	0.9793	
22 1,1,2-Trichloro-1,2,2-trif	151	2.280	2.280	0.000	63	10111	1.00	0.9383	
23 Acetone	43	2.280	2.292	-0.012	98	13879	2.00	2.32	
25 Iodomethane	142	2.351	2.363	-0.012	99	22823	1.00	1.00	
26 Carbon disulfide	76	2.410	2.411	0.000	98	46298	1.00	1.10	
28 3-Chloro-1-propene	76	2.541	2.541	0.000	88	10125	1.00	1.07	
29 Methyl acetate	43	2.564	2.564	0.000	97	25546	2.00	2.00	
30 Methylene Chloride	84	2.635	2.647	-0.012	93	28456	1.00	0.9073	
31 2-Methyl-2-propanol	59	2.754	2.754	0.000	95	24619	10.0	11.4	
32 Acrylonitrile	53	2.836	2.848	-0.012	97	64753	10.0	9.69	
34 trans-1,2-Dichloroethene	96	2.872	2.872	0.000	97	15989	1.00	1.05	
33 Methyl tert-butyl ether	73	2.872	2.884	-0.012	94	33845	1.00	0.99	
35 Hexane	86	3.097	3.109	-0.012	83	1522	1.00	1.07	

Compound	Sig	RT (min.)	Exp RT (min.)	Diff RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
36 1,1-Dichloroethane	63	3.215	3.215	0.000	95	26535	1.00	1.06	
37 Vinyl acetate	43	3.262	3.274	-0.012	95	21348	1.00	0.9075	
43 2,2-Dichloropropane	97	3.688	3.700	-0.012	54	4143	1.00	1.12	
41 cis-1,2-Dichloroethene	96	3.700	3.700	0.000	83	15518	1.00	0.9767	
42 2-Butanone (MEK)	43	3.712	3.712	0.000	69	12784	2.00	1.69	
47 Chlorobromomethane	128	3.901	3.890	0.011	94	8562	1.00	1.05	
48 Tetrahydrofuran	42	3.937	3.937	0.000	84	10465	2.00	2.07	
49 Chloroform	83	3.960	3.961	-0.001	96	23384	1.00	0.9654	
50 1,1,1-Trichloroethane	97	4.114	4.114	0.000	96	23404	1.00	1.11	
51 Cyclohexane	56	4.162	4.162	0.000	85	16522	1.00	0.8432	
52 1,1-Dichloropropene	75	4.244	4.256	-0.012	93	19040	1.00	0.9865	
53 Carbon tetrachloride	117	4.244	4.256	-0.012	96	21059	1.00	1.06	
54 Isobutyl alcohol	41	4.375	4.363	0.012	83	17638	25.0	25.5	
55 Benzene	78	4.422	4.422	0.000	94	56029	1.00	1.00	
56 1,2-Dichloroethane	62	4.434	4.434	0.000	95	20305	1.00	1.04	
58 n-Heptane	100	4.659	4.659	0.000	37	2973	1.00	1.07	
60 Trichloroethene	130	4.966	4.966	0.000	93	18033	1.00	1.08	
62 Methylcyclohexane	83	5.132	5.132	0.000	86	19643	1.00	0.9055	
63 1,2-Dichloropropane	63	5.156	5.156	0.000	91	11053	1.00	0.9425	
65 Dibromomethane	93	5.250	5.250	0.000	87	8259	1.00	1.00	
66 1,4-Dioxane	88	5.274	5.274	0.000	23	2371	20.0	22.7	
67 Dichlorobromomethane	83	5.392	5.392	0.000	96	18968	1.00	1.06	
69 2-Chloroethyl vinyl ether	63	5.653	5.653	0.000	94	11930	2.00	1.66	
70 cis-1,3-Dichloropropene	75	5.771	5.771	0.000	94	17074	1.00	0.9202	
71 4-Methyl-2-pentanone (MIBK)	43	5.913	5.913	0.000	93	23723	2.00	1.76	
72 Toluene	91	6.067	6.067	0.000	98	50368	1.00	0.99	
73 trans-1,3-Dichloropropene	75	6.268	6.268	0.000	94	17032	1.00	1.03	
74 Ethyl methacrylate	69	6.363	6.363	0.000	86	12213	1.00	0.8877	
75 1,1,2-Trichloroethane	97	6.422	6.422	0.000	84	12734	1.00	1.19	
76 Tetrachloroethene	164	6.552	6.552	0.000	97	11477	1.00	1.01	
77 1,3-Dichloropropane	76	6.576	6.576	0.000	90	17776	1.00	1.04	
78 2-Hexanone	43	6.658	6.659	-0.001	96	15753	2.00	1.73	
80 Chlorodibromomethane	129	6.777	6.777	0.000	93	10939	1.00	0.8840	
82 Ethylene Dibromide	107	6.871	6.872	-0.001	94	12725	1.00	1.16	a
84 Chlorobenzene	112	7.333	7.333	0.000	97	38544	1.00	1.13	
85 1,1,1,2-Tetrachloroethane	131	7.404	7.404	0.000	93	15538	1.00	1.19	
86 Ethylbenzene	106	7.439	7.439	0.000	96	17672	1.00	1.03	
87 m-Xylene & p-Xylene	106	7.546	7.546	0.000	98	17633	1.00	0.8856	
88 o-Xylene	106	7.913	7.913	0.000	95	17027	1.00	0.8837	
89 Styrene	104	7.924	7.925	-0.001	91	27201	1.00	0.8376	
90 Bromoform	173	8.090	8.090	0.000	95	9430	1.00	0.9571	
91 Isopropylbenzene	105	8.268	8.268	0.000	94	45741	1.00	0.9228	
95 Bromobenzene	156	8.552	8.540	0.012	85	14505	1.00	1.03	
94 1,1,2,2-Tetrachloroethane	83	8.552	8.552	0.000	93	15997	1.00	1.05	
96 1,2,3-Trichloropropane	110	8.587	8.587	0.000	82	5144	1.00	0.9763	
97 trans-1,4-Dichloro-2-buten	53	8.611	8.611	0.000	67	4646	1.00	0.9595	
98 N-Propylbenzene	120	8.658	8.658	0.000	98	12737	1.00	0.8965	
99 2-Chlorotoluene	126	8.729	8.741	-0.012	97	11634	1.00	0.9136	
100 1,3,5-Trimethylbenzene	105	8.836	8.836	0.000	89	34175	1.00	0.8833	
101 4-Chlorotoluene	126	8.847	8.848	-0.001	97	11598	1.00	0.8892	
102 tert-Butylbenzene	119	9.143	9.155	-0.012	89	28431	1.00	0.8591	
104 1,2,4-Trimethylbenzene	105	9.191	9.191	0.000	95	33988	1.00	0.8707	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
105 sec-Butylbenzene	105	9.356	9.356	0.000	93	41813	1.00	0.8872	
106 1,3-Dichlorobenzene	146	9.451	9.451	0.000	96	25676	1.00	1.01	
107 4-Isopropyltoluene	119	9.510	9.510	0.000	95	35416	1.00	0.8491	
108 1,4-Dichlorobenzene	146	9.546	9.546	0.000	92	32006	1.00	1.17	
112 1,2-Dichlorobenzene	146	9.901	9.901	0.000	88	23761	1.00	1.00	
111 n-Butylbenzene	91	9.901	9.901	0.000	95	33386	1.00	0.9814	
114 1,2-Dibromo-3-Chloropropan	157	10.670	10.670	0.000	90	6487	1.00	1.18	
116 1,2,4-Trichlorobenzene	180	11.486	11.486	0.000	92	15913	1.00	1.06	
117 Hexachlorobutadiene	225	11.676	11.676	0.000	92	8045	1.00	1.16	
118 Naphthalene	128	11.723	11.723	0.000	96	38023	1.00	0.9522	
119 1,2,3-Trichlorobenzene	180	11.971	11.960	0.011	92	14666	1.00	1.04	
S 159 Total BTEX	1				0		5.00	4.80	
S 133 Trihalomethanes, Total	1				0		4.00	3.87	
S 130 1,2-Dichloroethene, Total	96				0			2.02	
S 131 1,3-Dichloropropene, Total	75				0			1.95	
S 132 Xylenes, Total	106				0		2.00	1.77	

QC Flag Legend

Review Flags

a - User Assigned ID

Reagents:

VMRGAS_00281	Amount Added: 0.80	Units: uL
vm50ss_stk_00079	Amount Added: 0.80	Units: uL
vm50is_stk_A_00002	Amount Added: 2.00	Units: uL
VMRPRIMW_00320	Amount Added: 0.80	Units: uL
VMAROLISTDW_00283	Amount Added: 0.80	Units: uL



TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8026.D

Injection Date: 07-Feb-2019 15:00:30

Instrument ID: A3UX11

Operator ID: 43582

Lims ID: STD8260 L1

Worklist Smp#: 7

Client ID:

Purge Vol: 5.000 mL

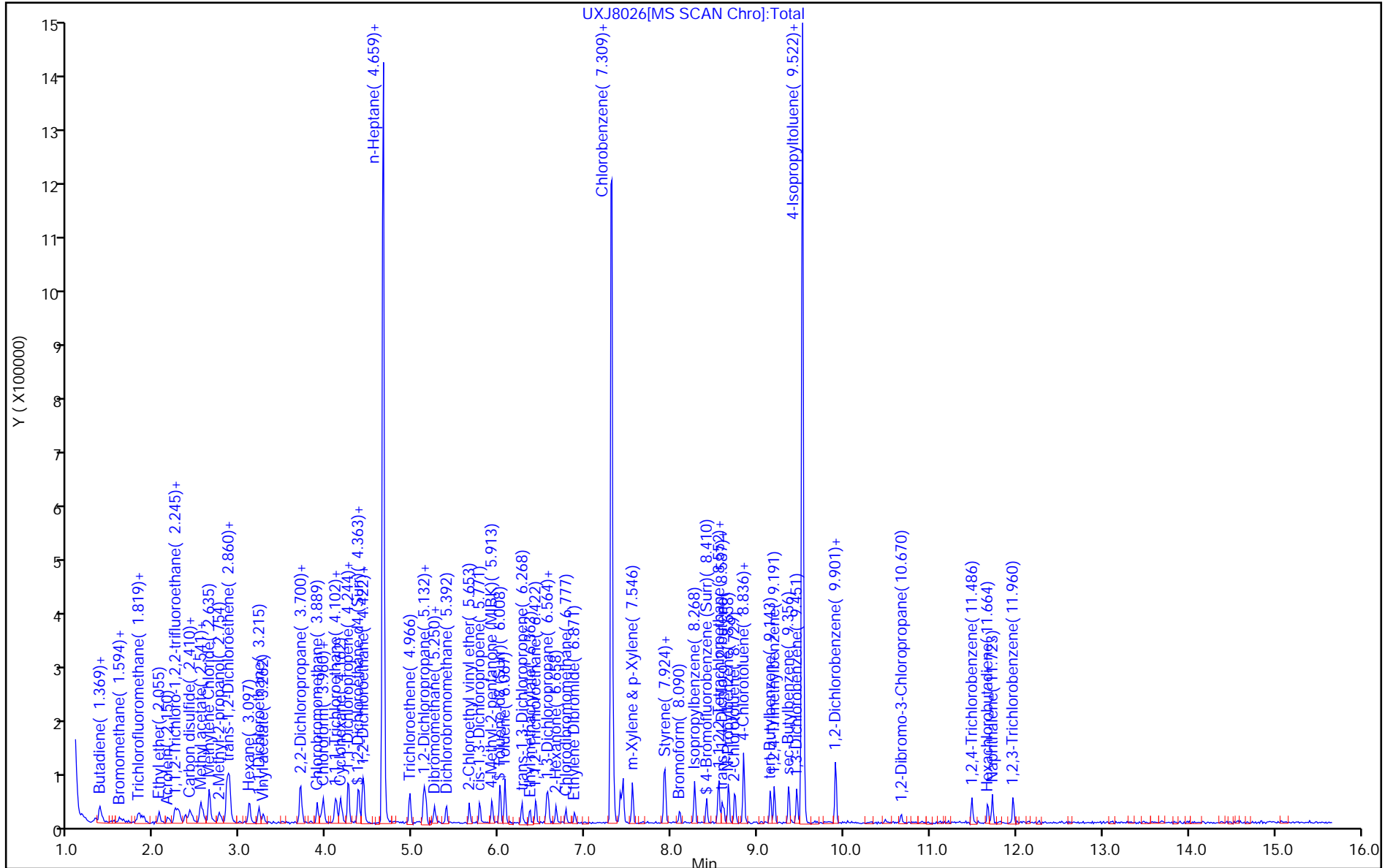
Dil. Factor: 1.0000

ALS Bottle#: 6

Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton

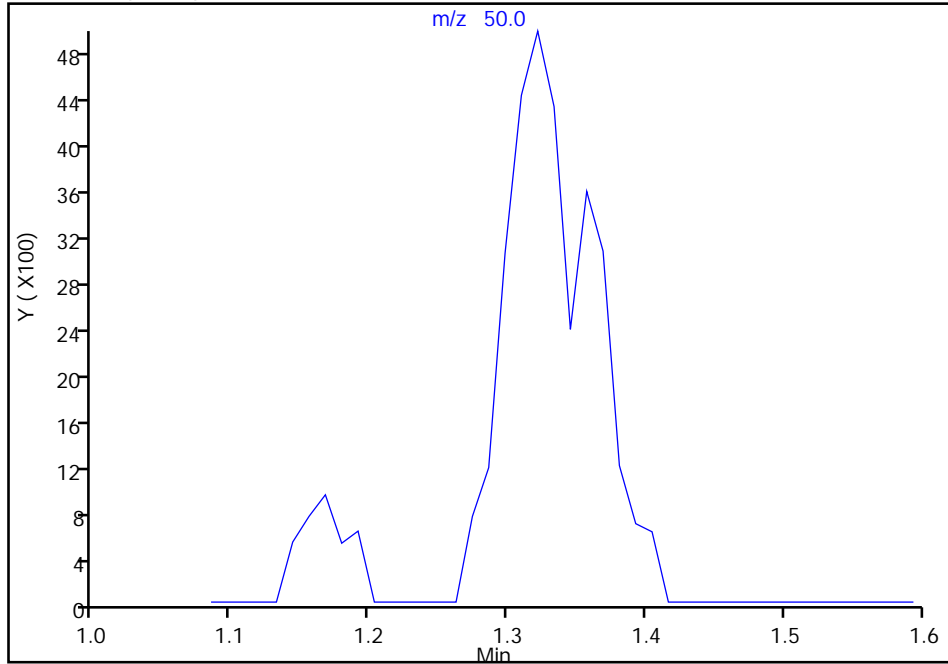
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Injection Date:	07-Feb-2019 15:00:30	Instrument ID:	A3UX11
Lims ID:	STD8260 L1		
Client ID:			
Operator ID:	43582	ALS Bottle#:	6
Purge Vol:	5.000 mL	Dil. Factor:	1.0000
Method:	8260_11	Limit Group:	MSV 8260B ICAL
Column:	DB-624 (0.18 mm)	Detector:	MS SCAN
		Worklist Smp#:	7

10 Chloromethane, CAS: 74-87-3

Signal: 1

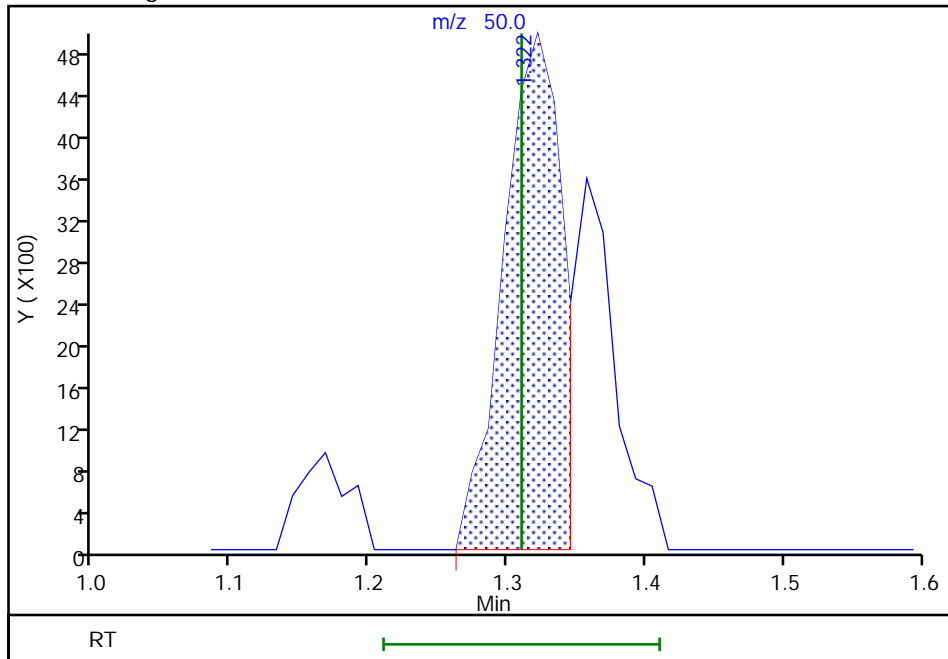
Not Detected
Expected RT: 1.31

Processing Integration Results



RT: 1.32
 Area: 14735
 Amount: 1.082937
 Amount Units: ug/l

Manual Integration Results



Reviewer: evansle, 08-Feb-2019 08:08:26

Audit Action: Assigned Compound ID

Audit Reason: Peak assignment corrected



TestAmerica Canton

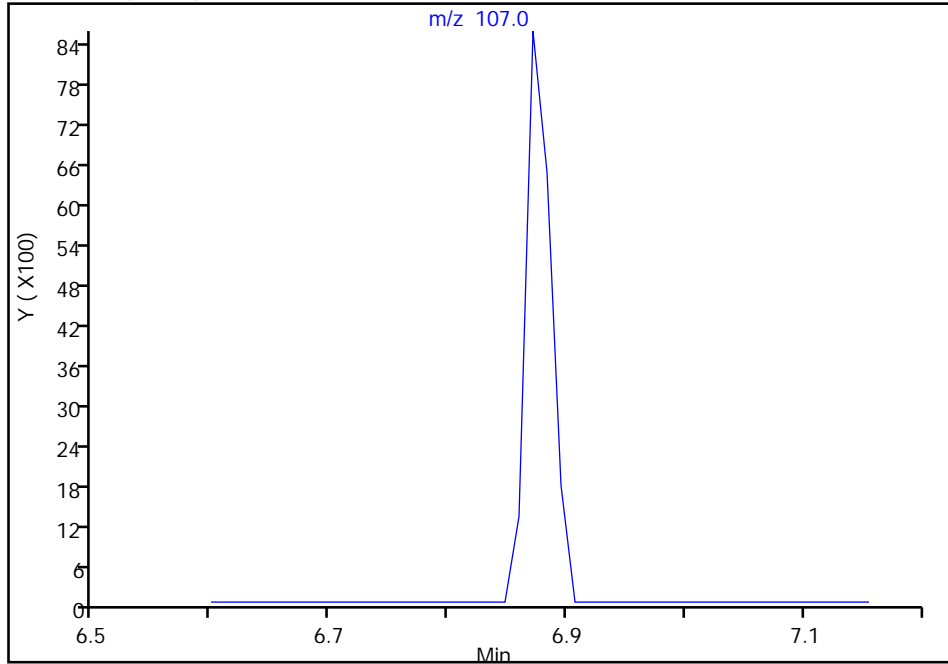
Data File:	\\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8026.D		
Injection Date:	07-Feb-2019 15:00:30	Instrument ID:	A3UX11
Lims ID:	STD8260 L1		
Client ID:			
Operator ID:	43582	ALS Bottle#:	6
Purge Vol:	5.000 mL	Dil. Factor:	1.0000
Method:	8260_11	Limit Group:	MSV 8260B ICAL
Column:	DB-624 (0.18 mm)	Detector:	MS SCAN
		Worklist Smp#:	7

82 Ethylene Dibromide, CAS: 106-93-4

Signal: 1

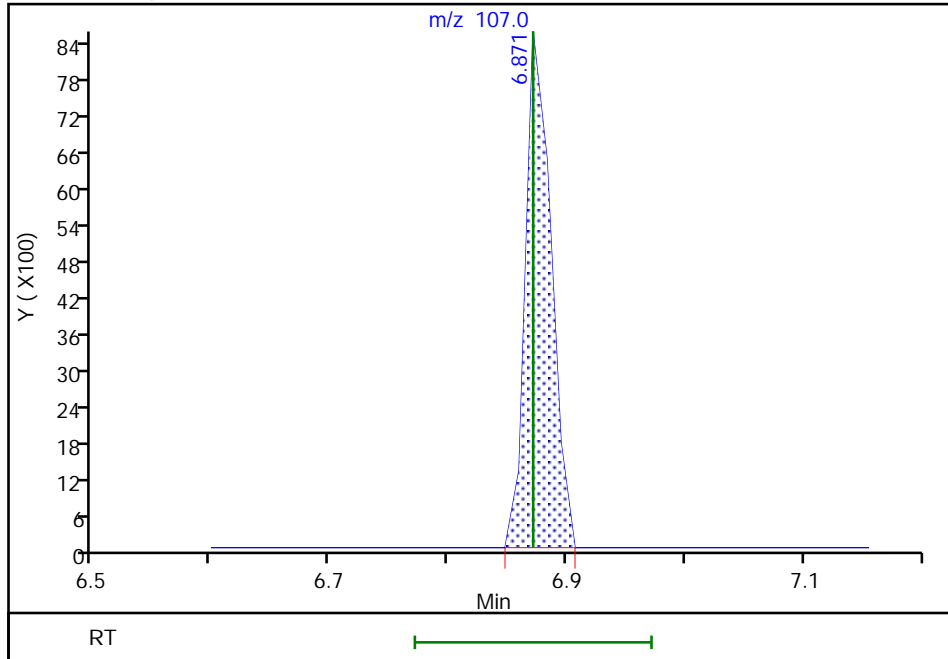
Not Detected
Expected RT: 6.87

Processing Integration Results



RT: 6.87
 Area: 12725
 Amount: 1.158386
 Amount Units: ug/l

Manual Integration Results



Reviewer: evansle, 08-Feb-2019 08:23:09
Audit Action: Assigned Compound ID

Audit Reason: Peak assignment corrected



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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 364785

SDG No.: _____

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 01/21/2019 11:37 Calibration End Date: 01/21/2019 13:26 Calibration ID: 49063

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD8260 240-364785/7	UXC8415.D
Level 2	STD8260 240-364785/6	UXC8414.D
Level 3	STD8260 240-364785/5	UXC8413.D
Level 4	STD8260 240-364785/4	UXC8412.D
Level 5	STD8260 240-364785/3	UXC8411.D
Level 6	STD8260 240-364785/2	UXC8410.D

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R ² OR COD	#	MIN R ² OR COD
	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
Dichlorodifluoromethane	0.2704 0.3071	0.3084	0.3222	0.3169	0.3182	Ave		0.3072			6.2		15.0				
Chloromethane	0.3289 0.2976	0.3274	0.3004	0.3124	0.3145	Ave		0.3135		0.1000	4.2		15.0				
Vinyl chloride	0.3261 0.2600	0.2864	0.2844	0.3065	0.2674	Ave		0.2885			8.5		15.0				
Butadiene	0.4028 0.3420	0.4066	0.3934	0.3729	0.3588	Ave		0.3794			6.8		15.0				
Bromomethane	0.2686 0.1933	0.2748	0.2163	0.2258	0.2060	Ave		0.2308			14.5		15.0				
Chloroethane	0.2298 0.1805	0.2264	0.1983	0.1989	0.1909	Ave		0.2041			9.7		15.0				
Dichlorofluoromethane	0.5874 0.4462	0.5508	0.5133	0.4857	0.4776	Ave		0.5102			10.1		15.0				
Trichlorofluoromethane	0.3885 0.4373	0.4314	0.4444	0.4568	0.4435	Ave		0.4337			5.5		15.0				
Ethyl ether	0.2527 0.2122	0.2393	0.2313	0.2239	0.2264	Ave		0.2310			6.0		15.0				
Acrolein	0.0365 0.0295	0.0303	0.0315	0.0310	0.0314	Ave		0.0317			7.8		15.0				
1,1-Dichloroethene	0.3235 0.2800	0.3292	0.3015	0.2758	0.2929	Ave		0.3005			7.3		15.0				
1,1,2-Trichloro-1,2,2-trifluoroethane	0.2411 0.2414	0.2509	0.2485	0.2467	0.2520	Ave		0.2468			1.9		15.0				
Acetone	0.0487 0.0213	0.0318	0.0276	0.0238	0.0247	Lin1	0.0519	0.0216						0.9960		0.9900	
Iodomethane	0.5303 0.4178	0.4271	0.4860	0.4853	0.4584	Ave		0.4675			9.0		15.0				
Carbon disulfide	0.9104 0.6640	0.7771	0.8074	0.7715	0.7145	Ave		0.7741			10.9		15.0				

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

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 364785

SDG No.: _____

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 01/21/2019 11:37 Calibration End Date: 01/21/2019 13:26 Calibration ID: 49063

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R ² OR COD	#	MIN R ² OR COD
	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
3-Chloro-1-propene	0.1530 0.1820	0.1600	0.1683	0.1693	0.1865	Ave		0.1698			7.5		15.0				
Methyl acetate	0.2259 0.1968	0.2056	0.2072	0.1929	0.2018	Ave		0.2050			5.6		15.0				
Methylene Chloride	0.8777 0.2608	0.5429	0.3681	0.3100	0.2858	Lin1	0.6169	0.2481						0.9990		0.9900	
2-Methyl-2-propanol	0.0205 0.0194	0.0225	0.0201	0.0175	0.0217	Ave		0.0203			8.7		15.0				
Acrylonitrile	0.0992 0.0981	0.0964	0.0945	0.0907	0.0981	Ave		0.0962			3.3		15.0				
Methyl tert-butyl ether	0.7930 0.7799	0.7813	0.7775	0.7682	0.8007	Ave		0.7834			1.5		15.0				
trans-1,2-Dichloroethene	0.3101 0.2660	0.2823	0.2736	0.2669	0.2784	Ave		0.2795			5.8		15.0				
Hexane	0.0608 0.0585	0.0607	0.0559	0.0560	0.0600	Ave		0.0586			3.8		15.0				
1,1-Dichloroethane	0.4596 0.4425	0.4568	0.4457	0.4441	0.4617	Ave		0.4517		0.1000	1.9		15.0				
Vinyl acetate	0.4987 0.5397	0.4861	0.5183	0.5294	0.5465	Ave		0.5198			4.5		15.0				
cis-1,2-Dichloroethene	0.3231 0.2906	0.3188	0.3032	0.3017	0.3096	Ave		0.3078			3.9		15.0				
2,2-Dichloropropane	0.0662 0.0571	0.0600	0.0637	0.0645	0.0628	Ave		0.0624			5.3		15.0				
2-Butanone (MEK)	0.0354 0.0328	0.0334	0.0325	0.0320	0.0334	Ave		0.0333			3.6		15.0				
Chlorobromomethane	0.1617 0.1477	0.1552	0.1523	0.1470	0.1549	Ave		0.1531			3.6		15.0				
Tetrahydrofuran	0.0906 0.0816	0.0853	0.0820	0.0745	0.0826	Ave		0.0828			6.4		15.0				
Chloroform	0.4755 0.4441	0.4649	0.4595	0.4523	0.4670	Ave		0.4606			2.4		15.0				
1,1,1-Trichloroethane	0.3615 0.3251	0.3567	0.3634	0.3678	0.3533	Ave		0.3546			4.3		15.0				
Cyclohexane	0.3905 0.3986	0.3949	0.4008	0.3961	0.4147	Ave		0.3993			2.1		15.0				
1,1-Dichloropropene	0.3764 0.3614	0.3858	0.3737	0.3673	0.3839	Ave		0.3748			2.5		15.0				
Carbon tetrachloride	0.3226 0.3473	0.3407	0.3468	0.3462	0.3607	Ave		0.3441			3.6		15.0				

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

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 364785

SDG No.: _____

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 01/21/2019 11:37 Calibration End Date: 01/21/2019 13:26 Calibration ID: 49063

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R ² OR COD	#	MIN R ² OR COD
	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
Isobutyl alcohol	0.0072 0.0077	0.0080	0.0077	0.0069	0.0086	Ave		0.0077			7.7		15.0				
Benzene	1.1739 1.0767	1.1296	1.1053	1.0802	1.1333	Ave		1.1165			3.3		15.0				
1,2-Dichloroethane	0.3855 0.3692	0.3897	0.3815	0.3777	0.3886	Ave		0.3820			2.0		15.0				
n-Heptane	0.0499 0.0533	0.0523	0.0527	0.0543	0.0568	Ave		0.0532			4.3		15.0				
Trichloroethene	0.3350 0.3084	0.3298	0.3223	0.3165	0.3318	Ave		0.3240			3.1		15.0				
Methylcyclohexane	0.3407 0.3494	0.3484	0.3548	0.3494	0.3719	Ave		0.3524			3.0		15.0				
1,2-Dichloropropane	0.2476 0.2542	0.2676	0.2609	0.2580	0.2658	Ave		0.2590			2.9		15.0				
Dibromomethane	0.1763 0.1613	0.1789	0.1712	0.1695	0.1730	Ave		0.1717			3.6		15.0				
1,4-Dioxane	0.0015 0.0021	0.0019	0.0021	0.0019	0.0024	Lin1	-0.013	0.0022						0.9940		0.9900	
Dichlorobromomethane	0.3251 0.3470	0.3335	0.3329	0.3402	0.3670	Ave		0.3410			4.3		15.0				
2-Chloroethyl vinyl ether	0.1741 0.1794	0.1688	0.1738	0.1710	0.1802	Ave		0.1745			2.6		15.0				
cis-1,3-Dichloropropene	0.3518 0.4202	0.3782	0.3877	0.3983	0.4304	Ave		0.3944			7.3		15.0				
4-Methyl-2-pentanone (MIBK)	0.2785 0.2686	0.2562	0.2693	0.2591	0.2753	Ave		0.2678			3.3		15.0				
Toluene	1.5012 1.4288	1.4730	1.4635	1.4625	1.5286	Ave		1.4763			2.3		15.0				
trans-1,3-Dichloropropene	0.4131 0.5013	0.4307	0.4524	0.4755	0.5145	Ave		0.4646			8.6		15.0				
Ethyl methacrylate	0.4397 0.4538	0.4440	0.4290	0.4357	0.4632	Ave		0.4442			2.8		15.0				
1,1,2-Trichloroethane	0.2996 0.2912	0.3009	0.2966	0.2943	0.3064	Ave		0.2982			1.8		15.0				
Tetrachloroethene	0.3214 0.3151	0.3301	0.3272	0.3250	0.3419	Ave		0.3268			2.8		15.0				
1,3-Dichloropropane	0.5170 0.5210	0.5290	0.5125	0.5245	0.5419	Ave		0.5243			2.0		15.0				
2-Hexanone	0.2098 0.2323	0.2329	0.2285	0.2242	0.2423	Ave		0.2283			4.8		15.0				

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

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 364785

SDG No.: _____

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 01/21/2019 11:37 Calibration End Date: 01/21/2019 13:26 Calibration ID: 49063

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R ² OR COD	#	MIN R ² OR COD
	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
Chlorodibromomethane	0.2647 0.3480	0.2850	0.3021	0.3162	0.3534	Ave		0.3116			11.2		15.0				
Ethylene Dibromide	0.3243 0.3131	0.3134	0.3221	0.3160	0.3258	Ave		0.3191			1.8		15.0				
Chlorobenzene	0.9888 0.9546	0.9840	0.9675	0.9702	1.0011	Ave		0.9777		0.3000	1.7		15.0				
1,1,1,2-Tetrachloroethane	0.3344 0.3512	0.3269	0.3433	0.3517	0.3627	Ave		0.3450			3.8		15.0				
Ethylbenzene	0.5276 0.5075	0.5044	0.5075	0.5104	0.5293	Ave		0.5145			2.1		15.0				
m-Xylene & p-Xylene	1.2031 1.2384	1.2096	1.1828	1.2237	1.2677	Ave		1.2209			2.4		15.0				
o-Xylene	0.6311 0.6050	0.6086	0.6092	0.6181	0.6414	Ave		0.6189			2.3		15.0				
Styrene	0.9861 1.0619	0.9946	1.0226	1.0439	1.1061	Ave		1.0359			4.3		15.0				
Bromoform	0.1422 0.2592	0.1774	0.1876	0.2065	0.2510	Lin1	-0.162	0.2546		0.1000				0.9940		0.9900	
Isopropylbenzene	1.4236 1.4623	1.4332	1.4499	1.4809	1.5299	Ave		1.4633			2.6		15.0				
1,1,2,2-Tetrachloroethane	0.7263 0.6975	0.7108	0.7013	0.6946	0.7209	Ave		0.7086		0.3000	1.8		15.0				
Bromobenzene	0.8425 0.8141	0.8308	0.8169	0.8158	0.8444	Ave		0.8274			1.7		15.0				
1,2,3-Trichloropropane	0.2495 0.2424	0.2553	0.2455	0.2417	0.2464	Ave		0.2468			2.0		15.0				
trans-1,4-Dichloro-2-butene	0.0434 0.1642	0.1010	0.0590	0.0798	0.1609	Ave		0.1014			50.5	*	15.0				
N-Propylbenzene	0.7265 0.7228	0.7398	0.7179	0.7231	0.7419	Ave		0.7286			1.3		15.0				
2-Chlorotoluene	0.7109 0.6635	0.6787	0.6802	0.6673	0.6918	Ave		0.6821			2.5		15.0				
1,3,5-Trimethylbenzene	2.1272 2.0859	2.0953	2.0772	2.1235	2.2036	Ave		2.1188			2.2		15.0				
4-Chlorotoluene	2.2568 2.1504	2.2467	2.1776	2.1619	2.2388	Ave		2.2054			2.1		15.0				
tert-Butylbenzene	1.8677 1.9925	1.8939	1.8513	1.8658	1.9133	Ave		1.8974			2.7		15.0				
1,2,4-Trimethylbenzene	2.2391 2.1885	2.2675	2.1657	2.2226	2.2813	Ave		2.2275			2.0		15.0				

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

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 364785

SDG No.: _____

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 01/21/2019 11:37 Calibration End Date: 01/21/2019 13:26 Calibration ID: 49063

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R ² OR COD	#	MIN R ² OR COD
	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
sec-Butylbenzene	2.2915 2.2504	2.3538	2.3171	2.3275	2.4198	Ave		2.3267			2.5		15.0				
1,3-Dichlorobenzene	1.4983 1.3538	1.4434	1.4148	1.4043	1.4340	Ave		1.4248			3.4		15.0				
4-Isopropyltoluene	2.0430 2.0131	2.1052	2.0529	2.1055	2.1774	Ave		2.0829			2.8		15.0				
1,4-Dichlorobenzene	1.5786 1.4104	1.5222	1.4387	1.4418	1.4716	Ave		1.4772			4.2		15.0				
n-Butylbenzene	1.6870 1.5530	1.6578	1.6297	1.6645	1.6997	Ave		1.6486			3.2		15.0				
1,2-Dichlorobenzene	1.4437 1.3061	1.4168	1.3751	1.3937	1.3869	Ave		1.3871			3.4		15.0				
1,2-Dibromo-3-Chloropropane	0.1400 0.1672	0.1533	0.1616	0.1661	0.1785	Ave		0.1611			8.2		15.0				
1,2,4-Trichlorobenzene	0.9724 0.7297	0.9185	0.8808	0.8895	0.8794	Ave		0.8784			9.2		15.0				
Hexachlorobutadiene	0.5543 0.3117	0.4931	0.4340	0.4326	0.4056	Qua	-0.024	0.4891	-0.004405					0.9990		0.9900	
Naphthalene	2.2857 1.9486	2.1983	2.2894	2.2740	2.2589	Ave		2.2091			6.0		15.0				
1,2,3-Trichlorobenzene	0.8662 0.6533	0.8274	0.8022	0.7980	0.7943	Ave		0.7902			9.1		15.0				
Dibromofluoromethane (Surr)	0.2371 0.2318	0.2202	0.2449	0.2390	0.2468	Ave		0.2366			4.1		15.0				
1,2-Dichloroethane-d4 (Surr)	0.3138 0.2896	0.2914	0.3140	0.3062	0.3101	Ave		0.3042			3.6		15.0				
Toluene-d8 (Surr)	1.2471 1.1640	1.1461	1.2369	1.2229	1.2656	Ave		1.2138			3.9		15.0				
4-Bromofluorobenzene (Surr)	0.4328 0.4271	0.4150	0.4450	0.4370	0.4478	Ave		0.4341			2.8		15.0				

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

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 364785

SDG No.: _____

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 01/21/2019 11:37 Calibration End Date: 01/21/2019 13:26 Calibration ID: 49063

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD8260 240-364785/7	UXC8415.D
Level 2	STD8260 240-364785/6	UXC8414.D
Level 3	STD8260 240-364785/5	UXC8413.D
Level 4	STD8260 240-364785/4	UXC8412.D
Level 5	STD8260 240-364785/3	UXC8411.D
Level 6	STD8260 240-364785/2	UXC8410.D

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/L)				
			LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5
Dichlorodifluoromethane	FB	Ave	47729 2392918	112693	296202	593761	1222220	1.00 40.0	2.00	5.00	10.0	20.0
Chloromethane	FB	Ave	58046 2319242	119628	276149	585358	1207742	1.00 40.0	2.00	5.00	10.0	20.0
Vinyl chloride	FB	Ave	57558 2025909	104647	261470	574312	1027051	1.00 40.0	2.00	5.00	10.0	20.0
Butadiene	FB	Ave	71088 2665435	148568	361713	698736	1378009	1.00 40.0	2.00	5.00	10.0	20.0
Bromomethane	FB	Ave	47411 1506749	100403	198900	423055	791041	1.00 40.0	2.00	5.00	10.0	20.0
Chloroethane	FB	Ave	40558 1406432	82718	182320	372694	733356	1.00 40.0	2.00	5.00	10.0	20.0
Dichlorofluoromethane	FB	Ave	103682 3477509	201291	471934	910063	1834172	1.00 40.0	2.00	5.00	10.0	20.0
Trichlorofluoromethane	FB	Ave	68568 3407867	157640	408598	855986	1703371	1.00 40.0	2.00	5.00	10.0	20.0
Ethyl ether	FB	Ave	44605 1653914	87432	212667	419544	869536	1.00 40.0	2.00	5.00	10.0	20.0
Acrolein	FB	Ave	32229 1148055	55282	144856	290454	603583	5.00 200	10.0	25.0	50.0	100
1,1-Dichloroethene	FB	Ave	57094 2182381	120279	277159	516876	1125076	1.00 40.0	2.00	5.00	10.0	20.0
1,1,2-Trichloro-1,2,2-trifluoroethane	FB	Ave	42550 1880984	91700	228498	462285	967763	1.00 40.0	2.00	5.00	10.0	20.0
Acetone	FB	Lin1	17181 332724	23252	50751	89168	190030	2.00 80.0	4.00	10.0	20.0	40.0
Iodomethane	FB	Ave	93588 3255913	156089	446819	909454	1760595	1.00 40.0	2.00	5.00	10.0	20.0
Carbon disulfide	FB	Ave	160687 5174431	283969	742305	1445567	2744114	1.00 40.0	2.00	5.00	10.0	20.0
3-Chloro-1-propene	FB	Ave	27003 1418141	58477	154763	317161	716236	1.00 40.0	2.00	5.00	10.0	20.0

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 364785

SDG No.: _____

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 01/21/2019 11:37 Calibration End Date: 01/21/2019 13:26 Calibration ID: 49063

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/L)				
			LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5
Methyl acetate	FB	Ave	79740 3067156	150240	380996	723073	1550490	2.00 80.0	4.00	10.0	20.0	40.0
Methylene Chloride	FB	Lin1	154915 2032773	198393	338417	580808	1097865	1.00 40.0	2.00	5.00	10.0	20.0
2-Methyl-2-propanol	FB	Ave	36240 1508945	82343	184510	328142	832907	10.0 400	20.0	50.0	100	200
Acrylonitrile	FB	Ave	175156 7643188	352291	868795	1699336	3768272	10.0 400	20.0	50.0	100	200
Methyl tert-butyl ether	FB	Ave	139955 6078083	285516	714788	1439491	3075308	1.00 40.0	2.00	5.00	10.0	20.0
trans-1,2-Dichloroethene	FB	Ave	54736 2072831	103155	251570	500088	1069131	1.00 40.0	2.00	5.00	10.0	20.0
Hexane	FB	Ave	10725 455555	22194	51427	104918	230317	1.00 40.0	2.00	5.00	10.0	20.0
1,1-Dichloroethane	FB	Ave	81126 3448011	166930	409730	832192	1773410	1.00 40.0	2.00	5.00	10.0	20.0
Vinyl acetate	FB	Ave	88027 4206187	177639	476538	992089	2098831	1.00 40.0	2.00	5.00	10.0	20.0
cis-1,2-Dichloroethene	FB	Ave	57035 2264494	116480	278759	565374	1189038	1.00 40.0	2.00	5.00	10.0	20.0
2,2-Dichloropropane	FB	Ave	11684 445323	21907	58586	120845	241048	1.00 40.0	2.00	5.00	10.0	20.0
2-Butanone (MEK)	FB	Ave	12505 511818	24387	59681	120010	256717	2.00 80.0	4.00	10.0	20.0	40.0
Chlorobromomethane	FB	Ave	28543 1150919	56716	140018	275468	595090	1.00 40.0	2.00	5.00	10.0	20.0
Tetrahydrofuran	FB	Ave	31985 1272086	62314	150812	279158	634219	2.00 80.0	4.00	10.0	20.0	40.0
Chloroform	FB	Ave	83929 3460990	169874	422474	847597	1793716	1.00 40.0	2.00	5.00	10.0	20.0
1,1,1-Trichloroethane	FB	Ave	63803 2533663	130349	334136	689229	1356966	1.00 40.0	2.00	5.00	10.0	20.0
Cyclohexane	FB	Ave	68915 3105972	144320	368473	742256	1592895	1.00 40.0	2.00	5.00	10.0	20.0
1,1-Dichloropropene	FB	Ave	66437 2816568	140972	343617	688186	1474449	1.00 40.0	2.00	5.00	10.0	20.0
Carbon tetrachloride	FB	Ave	56945 2706723	124489	318843	648774	1385341	1.00 40.0	2.00	5.00	10.0	20.0
Isobutyl alcohol	FB	Ave	31904 1496664	73133	177736	322275	821992	25.0 1000	50.0	125	250	500
Benzene	FB	Ave	207184 8390558	412777	1016161	2024145	4352726	1.00 40.0	2.00	5.00	10.0	20.0

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 364785

SDG No.: _____

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 01/21/2019 11:37 Calibration End Date: 01/21/2019 13:26 Calibration ID: 49063

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/L)				
			LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5
1,2-Dichloroethane	FB	Ave	68031 2877185	142414	350758	707702	1492475	1.00 40.0	2.00	5.00	10.0	20.0
n-Heptane	FB	Ave	8801 415529	19095	48491	101813	218160	1.00 40.0	2.00	5.00	10.0	20.0
Trichloroethene	FB	Ave	59121 2403531	120531	296310	593061	1274454	1.00 40.0	2.00	5.00	10.0	20.0
Methylcyclohexane	FB	Ave	60130 2723193	127322	326169	654672	1428283	1.00 40.0	2.00	5.00	10.0	20.0
1,2-Dichloropropane	FB	Ave	43707 1981074	97800	239856	483376	1020861	1.00 40.0	2.00	5.00	10.0	20.0
Dibromomethane	FB	Ave	31111 1256838	65366	157375	317664	664429	1.00 40.0	2.00	5.00	10.0	20.0
1,4-Dioxane	FB	Lin1	5185 326819	14227	39232	71764	183906	20.0 800	40.0	100	200	400
Dichlorobromomethane	FB	Ave	57382 2704273	121869	306036	637533	1409597	1.00 40.0	2.00	5.00	10.0	20.0
2-Chloroethyl vinyl ether	FB	Ave	61442 2796269	123378	319533	640886	1383920	2.00 80.0	4.00	10.0	20.0	40.0
cis-1,3-Dichloropropene	FB	Ave	62099 3274371	138202	356486	746312	1652896	1.00 40.0	2.00	5.00	10.0	20.0
4-Methyl-2-pentanone (MIBK)	FB	Ave	98299 4186633	187250	495239	970866	2114735	2.00 80.0	4.00	10.0	20.0	40.0
Toluene	CBNZ d5	Ave	209687 9034071	428381	1080197	2172974	4691174	1.00 40.0	2.00	5.00	10.0	20.0
trans-1,3-Dichloropropene	CBNZ d5	Ave	57707 3169409	125256	333930	706531	1578892	1.00 40.0	2.00	5.00	10.0	20.0
Ethyl methacrylate	CBNZ d5	Ave	61421 2869285	129112	316654	647349	1421429	1.00 40.0	2.00	5.00	10.0	20.0
1,1,2-Trichloroethane	CBNZ d5	Ave	41853 1840913	87521	218933	437245	940456	1.00 40.0	2.00	5.00	10.0	20.0
Tetrachloroethene	CBNZ d5	Ave	44889 1992562	95990	241473	482895	1049416	1.00 40.0	2.00	5.00	10.0	20.0
1,3-Dichloropropane	CBNZ d5	Ave	72217 3294456	153833	378264	779324	1663194	1.00 40.0	2.00	5.00	10.0	20.0
2-Hexanone	CBNZ d5	Ave	58616 2937075	135481	337296	666112	1487023	2.00 80.0	4.00	10.0	20.0	40.0
Chlorodibromomethane	CBNZ d5	Ave	36977 2200067	82891	222991	469828	1084488	1.00 40.0	2.00	5.00	10.0	20.0
Ethylene Dibromide	CBNZ d5	Ave	45295 1979685	91132	237763	469490	999728	1.00 40.0	2.00	5.00	10.0	20.0
Chlorobenzene	CBNZ d5	Ave	138107 6035718	286159	714138	1441533	3072375	1.00 40.0	2.00	5.00	10.0	20.0

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 364785

SDG No.: _____

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 01/21/2019 11:37 Calibration End Date: 01/21/2019 13:26 Calibration ID: 49063

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/L)				
			LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5
1,1,1,2-Tetrachloroethane	CBNZ d5	Ave	46702 2220799	95060	253355	522591	1113186	1.00 40.0	2.00	5.00	10.0	20.0
Ethylbenzene	CBNZ d5	Ave	73699 3208633	146696	374569	758362	1624425	1.00 40.0	2.00	5.00	10.0	20.0
m-Xylene & p-Xylene	CBNZ d5	Ave	168049 7830315	351787	872988	1818169	3890600	1.00 40.0	2.00	5.00	10.0	20.0
o-Xylene	CBNZ d5	Ave	88147 3825283	176986	449671	918350	1968397	1.00 40.0	2.00	5.00	10.0	20.0
Styrene	CBNZ d5	Ave	137740 6714209	289260	754758	1550959	3394736	1.00 40.0	2.00	5.00	10.0	20.0
Bromoform	CBNZ d5	Lin1	19867 1639155	51580	138461	306750	770158	1.00 40.0	2.00	5.00	10.0	20.0
Isopropylbenzene	CBNZ d5	Ave	198841 9246149	416807	1070139	2200366	4695119	1.00 40.0	2.00	5.00	10.0	20.0
1,1,2,2-Tetrachloroethane	DCBd 4	Ave	54865 2441270	112833	283791	577712	1238040	1.00 40.0	2.00	5.00	10.0	20.0
Bromobenzene	DCBd 4	Ave	63640 2849353	131876	330598	678570	1450127	1.00 40.0	2.00	5.00	10.0	20.0
1,2,3-Trichloropropane	DCBd 4	Ave	18846 848462	40526	99350	201010	423192	1.00 40.0	2.00	5.00	10.0	20.0
trans-1,4-Dichloro-2-butene	DCBd 4	Ave	3277 574750	16039	23872	66363	276245	1.00 40.0	2.00	5.00	10.0	20.0
N-Propylbenzene	DCBd 4	Ave	54874 2529942	117423	290505	601438	1274023	1.00 40.0	2.00	5.00	10.0	20.0
2-Chlorotoluene	DCBd 4	Ave	53702 2322347	107735	275264	555062	1188057	1.00 40.0	2.00	5.00	10.0	20.0
1,3,5-Trimethylbenzene	DCBd 4	Ave	160683 7300680	332591	840604	1766210	3784392	1.00 40.0	2.00	5.00	10.0	20.0
4-Chlorotoluene	DCBd 4	Ave	170473 7526400	356628	881245	1798133	3844885	1.00 40.0	2.00	5.00	10.0	20.0
tert-Butylbenzene	DCBd 4	Ave	141080 6973779	300630	749194	1551899	3285832	1.00 40.0	2.00	5.00	10.0	20.0
1,2,4-Trimethylbenzene	DCBd 4	Ave	169136 7660047	359918	876431	1848645	3917785	1.00 40.0	2.00	5.00	10.0	20.0
sec-Butylbenzene	DCBd 4	Ave	173094 7876504	373619	937713	1935865	4155710	1.00 40.0	2.00	5.00	10.0	20.0
1,3-Dichlorobenzene	DCBd 4	Ave	113176 4738373	229119	572540	1168047	2462678	1.00 40.0	2.00	5.00	10.0	20.0
4-Isopropyltoluene	DCBd 4	Ave	154318 7046098	334169	830781	1751242	3739399	1.00 40.0	2.00	5.00	10.0	20.0
1,4-Dichlorobenzene	DCBd 4	Ave	119243 4936571	241629	582208	1199185	2527274	1.00 40.0	2.00	5.00	10.0	20.0

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

Lab Name: TestAmerica Canton Job No.: 240-109102-1 Analy Batch No.: 364785

SDG No.: _____

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 01/21/2019 11:37 Calibration End Date: 01/21/2019 13:26 Calibration ID: 49063

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/L)				
			LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2	LVL 3	LVL 4	LVL 5
n-Butylbenzene	DCBd 4	Ave	127428 5435677	263153	659506	1384405	2918965	1.00 40.0	2.00	5.00	10.0	20.0
1,2-Dichlorobenzene	DCBd 4	Ave	109056 4571601	224899	556492	1159176	2381785	1.00 40.0	2.00	5.00	10.0	20.0
1,2-Dibromo-3-Chloropropane	DCBd 4	Ave	10572 585195	24334	65389	138143	306515	1.00 40.0	2.00	5.00	10.0	20.0
1,2,4-Trichlorobenzene	DCBd 4	Ave	73453 2553990	145803	356465	739836	1510157	1.00 40.0	2.00	5.00	10.0	20.0
Hexachlorobutadiene	DCBd 4	Qua	41867 1091039	78276	175613	359850	696597	1.00 40.0	2.00	5.00	10.0	20.0
Naphthalene	DCBd 4	Ave	172653 6820199	348942	926474	1891386	3879317	1.00 40.0	2.00	5.00	10.0	20.0
1,2,3-Trichlorobenzene	DCBd 4	Ave	65430 2286670	131338	324635	663704	1364016	1.00 40.0	2.00	5.00	10.0	20.0
Dibromofluoromethane (Surr)	FB	Ave	41840 1806247	80448	225163	447792	947952	1.00 40.0	2.00	5.00	10.0	20.0
1,2-Dichloroethane-d4 (Surr)	FB	Ave	55383 2257190	106482	288691	573781	1190875	1.00 40.0	2.00	5.00	10.0	20.0
Toluene-d8 (Surr)	CBNZ d5	Ave	174193 7359833	333303	912990	1816996	3883966	1.00 40.0	2.00	5.00	10.0	20.0
4-Bromofluorobenzene (Surr)	CBNZ d5	Ave	60446 2700470	120695	328418	649263	1374232	1.00 40.0	2.00	5.00	10.0	20.0

Curve Type Legend:

Ave = Average ISTD
Linl = Linear 1/conc ISTD
Qua = Quadratic ISTD

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8410.D
 Lims ID: STD8260 L6
 Client ID:
 Sample Type: IC Calib Level: 6
 Inject. Date: 21-Jan-2019 11:37:30 ALS Bottle#: 1 Worklist Smp#: 2
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0083824-002
 Misc. Info.: C90121A-IC,8260LLUX15,,43582
 Operator ID: Instrument ID: A3UX15
 Sublist: chrom-8260_15*sub83
 Method: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\8260_15.m
 Limit Group: MSV 8260B ICAL
 Last Update: 22-Jan-2019 09:12:44 Calib Date: 21-Jan-2019 19:15:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8431.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0303

First Level Reviewer: evansle

Date: 22-Jan-2019 09:12:32

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	5.208	5.196	0.012	99	1948244	10.0	10.0	
* 2 Chlorobenzene-d5	117	7.876	7.876	0.000	84	1580707	10.0	10.0	
* 3 1,4-Dichlorobenzene-d4	152	10.117	10.117	0.000	91	875018	10.0	10.0	
\$ 4 Dibromofluoromethane (Surr	113	4.639	4.639	0.000	95	1806247	40.0	39.2	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.923	4.923	0.000	100	2257190	40.0	38.1	
\$ 6 Toluene-d8 (Surr)	98	6.560	6.560	0.000	93	7359833	40.0	38.4	
\$ 7 4-Bromofluorobenzene (Surr	95	8.979	8.979	0.000	97	2700470	40.0	39.4	
9 Dichlorodifluoromethane	85	1.532	1.532	0.000	99	2392918	40.0	40.0	
10 Chloromethane	50	1.698	1.686	0.012	99	2319242	40.0	38.0	
11 Vinyl chloride	62	1.793	1.781	0.012	98	2025909	40.0	36.0	
12 Butadiene	54	1.804	1.793	0.011	98	2665435	40.0	36.1	
13 Bromomethane	94	2.065	2.054	0.011	90	1506749	40.0	33.5	
14 Chloroethane	64	2.160	2.137	0.023	100	1406432	40.0	35.4	
15 Dichlorofluoromethane	67	2.314	2.303	0.011	97	3477509	40.0	35.0	
16 Trichlorofluoromethane	101	2.362	2.362	0.000	98	3407867	40.0	40.3	
17 Ethyl ether	59	2.575	2.563	0.012	89	1653914	40.0	36.8	
18 Acrolein	56	2.694	2.694	0.000	100	1148055	200.0	185.9	
19 1,1-Dichloroethene	96	2.812	2.777	0.035	98	2182381	40.0	37.3	
20 1,1,2-Trichloro-1,2,2-trif	151	2.812	2.789	0.023	79	1880984	40.0	39.1	
22 Acetone	58	2.824	2.812	0.012	99	332724	80.0	76.8	
23 Iodomethane	142	2.955	2.931	0.024	98	3255913	40.0	35.7	
24 Carbon disulfide	76	3.002	2.990	0.012	100	5174431	40.0	34.3	
25 3-Chloro-1-propene	76	3.085	3.073	0.012	88	1418141	40.0	42.9	
27 Methyl acetate	43	3.097	3.085	0.012	98	3067156	80.0	76.8	
28 Methylene Chloride	84	3.192	3.180	0.012	92	2032773	40.0	39.6	
29 2-Methyl-2-propanol	59	3.263	3.263	0.000	99	1508945	400.0	381.9	
32 Acrylonitrile	53	3.393	3.394	-0.001	99	7643188	400.0	407.9	
31 Methyl tert-butyl ether	73	3.405	3.405	0.000	96	6078083	40.0	39.8	
30 trans-1,2-Dichloroethene	96	3.417	3.405	0.012	95	2072831	40.0	38.1	
33 Hexane	86	3.631	3.619	0.012	93	455555	40.0	39.9	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
34 1,1-Dichloroethane	63	3.761	3.761	0.000	96	3448011	40.0	39.2	
36 Vinyl acetate	43	3.797	3.785	0.012	97	4206187	40.0	41.5	
39 cis-1,2-Dichloroethene	96	4.247	4.235	0.012	81	2264494	40.0	37.8	
40 2,2-Dichloropropane	97	4.247	4.247	0.000	60	445323	40.0	36.6	
41 2-Butanone (MEK)	72	4.247	4.247	0.000	99	511818	80.0	79.0	
45 Chlorobromomethane	128	4.449	4.449	0.000	89	1150919	40.0	38.6	
46 Tetrahydrofuran	42	4.484	4.485	-0.001	87	1272086	80.0	78.9	
47 Chloroform	83	4.508	4.496	0.012	93	3460990	40.0	38.6	
48 1,1,1-Trichloroethane	97	4.662	4.662	0.000	98	2533663	40.0	36.7	
49 Cyclohexane	56	4.722	4.710	0.012	90	3105972	40.0	39.9	
51 1,1-Dichloropropene	75	4.793	4.793	0.000	95	2816568	40.0	38.6	
50 Carbon tetrachloride	117	4.805	4.805	0.000	97	2706723	40.0	40.4	
52 Isobutyl alcohol	41	4.864	4.864	0.000	95	1496664	1000.0	1000.0	
53 Benzene	78	4.982	4.971	0.011	96	8390558	40.0	38.6	
54 1,2-Dichloroethane	62	4.994	4.983	0.011	98	2877185	40.0	38.7	
56 n-Heptane	100	5.172	5.160	0.012	93	415529	40.0	40.1	
58 Trichloroethene	130	5.516	5.504	0.012	96	2403531	40.0	38.1	
60 Methylcyclohexane	83	5.682	5.670	0.012	88	2723193	40.0	39.7	
61 1,2-Dichloropropane	63	5.706	5.706	0.000	95	1981074	40.0	39.3	
64 1,4-Dioxane	88	5.813	5.813	0.000	59	326819	800.0	772.3	
63 Dibromomethane	93	5.813	5.813	0.000	88	1256838	40.0	37.6	
65 Dichlorobromomethane	83	5.931	5.931	0.000	99	2704273	40.0	40.7	
67 2-Chloroethyl vinyl ether	63	6.180	6.180	0.000	93	2796269	80.0	82.2	
68 cis-1,3-Dichloropropene	75	6.322	6.323	-0.001	95	3274371	40.0	42.6	
69 4-Methyl-2-pentanone (MIBK)	43	6.453	6.441	0.012	97	4186633	80.0	80.2	
70 Toluene	91	6.619	6.619	0.000	98	9034071	40.0	38.7	
71 trans-1,3-Dichloropropene	75	6.809	6.809	0.000	94	3169409	40.0	43.2	
72 Ethyl methacrylate	69	6.868	6.868	0.000	89	2869285	40.0	40.9	
73 1,1,2-Trichloroethane	97	6.975	6.975	0.000	91	1840913	40.0	39.1	
74 Tetrachloroethene	164	7.117	7.117	0.000	96	1992562	40.0	38.6	
75 1,3-Dichloropropane	76	7.129	7.129	0.000	93	3294456	40.0	39.7	
76 2-Hexanone	43	7.188	7.188	0.000	97	2937075	80.0	81.4	
79 Chlorodibromomethane	129	7.342	7.342	0.000	90	2200067	40.0	44.7	
80 Ethylene Dibromide	107	7.461	7.461	0.000	98	1979685	40.0	39.2	
82 Chlorobenzene	112	7.900	7.900	0.000	96	6035718	40.0	39.1	
83 1,1,1,2-Tetrachloroethane	131	7.971	7.971	0.000	94	2220799	40.0	40.7	
84 Ethylbenzene	106	7.995	7.995	0.000	98	3208633	40.0	39.5	
85 m-Xylene & p-Xylene	91	8.101	8.101	0.000	93	7830315	40.0	40.6	
86 o-Xylene	106	8.481	8.481	0.000	97	3825283	40.0	39.1	
87 Styrene	104	8.493	8.493	0.000	94	6714209	40.0	41.0	
88 Bromoform	173	8.682	8.682	0.000	98	1639155	40.0	41.4	
89 Isopropylbenzene	105	8.825	8.825	0.000	95	9246149	40.0	40.0	
93 1,1,2,2-Tetrachloroethane	83	9.121	9.121	0.000	97	2441270	40.0	39.4	
92 Bromobenzene	156	9.145	9.133	0.012	90	2849353	40.0	39.4	
94 1,2,3-Trichloropropane	110	9.169	9.169	0.000	85	848462	40.0	39.3	
95 trans-1,4-Dichloro-2-buten	53	9.169	9.169	0.000	80	574750	40.0	64.8	
96 N-Propylbenzene	120	9.228	9.228	0.000	98	2529942	40.0	39.7	
97 2-Chlorotoluene	126	9.323	9.323	0.000	98	2322347	40.0	38.9	
98 1,3,5-Trimethylbenzene	105	9.394	9.394	0.000	94	7300680	40.0	39.4	
99 4-Chlorotoluene	91	9.418	9.418	0.000	98	7526400	40.0	39.0	
101 tert-Butylbenzene	119	9.714	9.714	0.000	92	6973779	40.0	42.0	
103 1,2,4-Trimethylbenzene	105	9.761	9.762	-0.001	94	7660047	40.0	39.3	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
105 sec-Butylbenzene	105	9.927	9.928	-0.001	93	7876504	40.0	38.7	
106 1,3-Dichlorobenzene	146	10.046	10.046	0.000	98	4738373	40.0	38.0	
107 4-Isopropyltoluene	119	10.070	10.070	0.000	97	7046098	40.0	38.7	
108 1,4-Dichlorobenzene	146	10.141	10.141	0.000	97	4936571	40.0	38.2	
111 n-Butylbenzene	91	10.473	10.473	0.000	97	5435677	40.0	37.7	
112 1,2-Dichlorobenzene	146	10.509	10.509	0.000	99	4571601	40.0	37.7	
113 1,2-Dibromo-3-Chloropropan	157	11.291	11.291	0.000	93	585195	40.0	41.5	
115 1,2,4-Trichlorobenzene	180	12.109	12.110	-0.001	94	2553990	40.0	33.2	
116 Hexachlorobutadiene	225	12.275	12.276	-0.001	93	1091039	40.0	39.8	
117 Naphthalene	128	12.358	12.359	-0.001	96	6820199	40.0	35.3	
118 1,2,3-Trichlorobenzene	180	12.608	12.608	0.000	95	2286670	40.0	33.1	
S 127 1,2-Dichloroethene, Total	96				0			75.8	
S 128 1,3-Dichloropropene, Total	75				0			85.8	
S 129 Xylenes, Total	106				0		80.0	79.7	
S 130 Trihalomethanes, Total	1				0		160.0	165.3	
S 157 Total BTEX	1		0.000				200.0	ND	

Reagents:

vm50ss_stk_00079	Amount Added: 32.00	Units: uL
VM50IS_00072	Amount Added: 1.00	Units: uL
VMRPRIMW_00318	Amount Added: 32.00	Units: uL
VMAROLISTDW_00281	Amount Added: 32.00	Units: uL
VMRGAS_00278	Amount Added: 32.00	Units: uL

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8410.D

Injection Date: 21-Jan-2019 11:37:30

Instrument ID: A3UX15

Operator ID:

Lims ID: STD8260 L6

Worklist Smp#: 2

Client ID:

Purge Vol: 5.000 mL

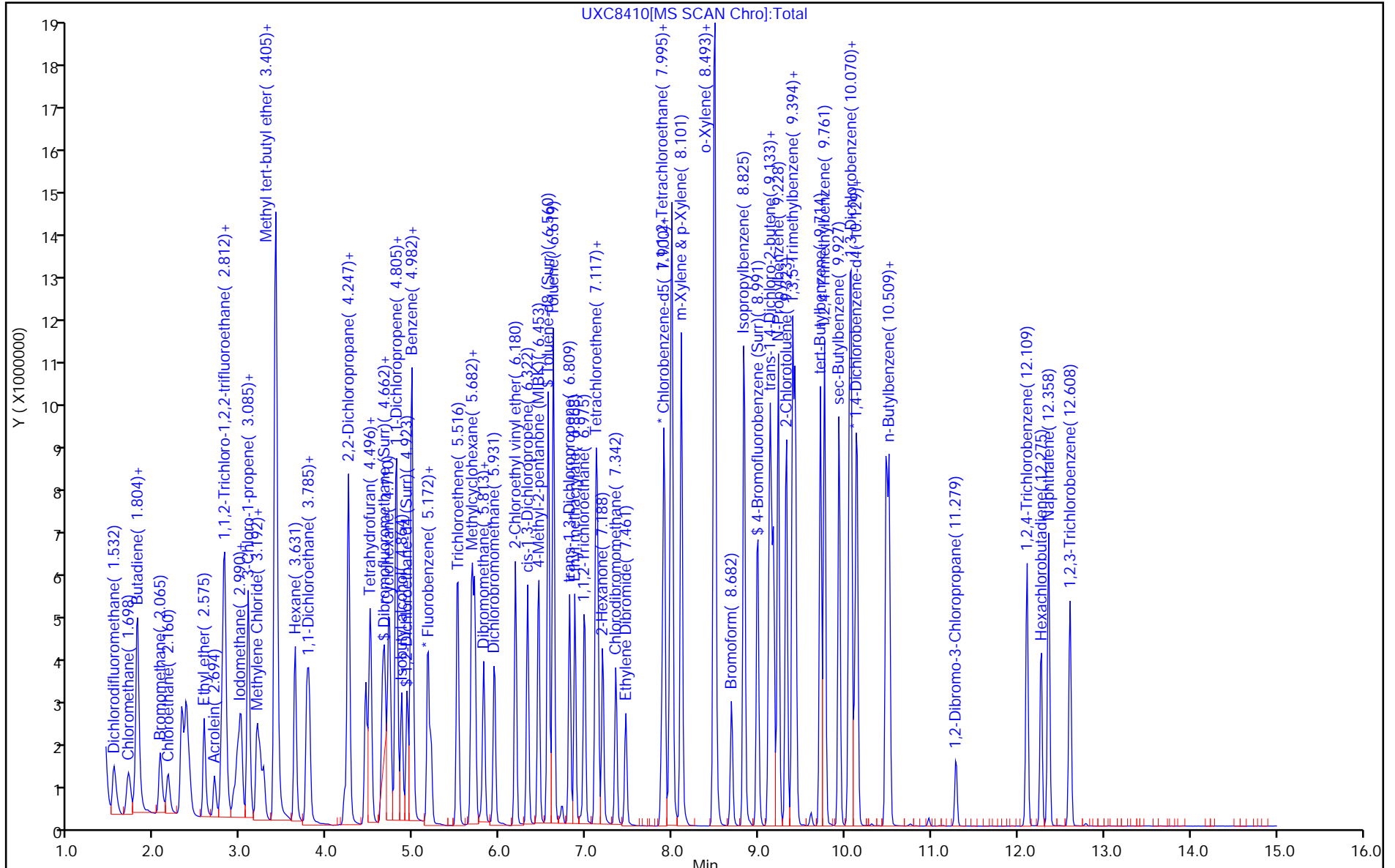
Dil. Factor: 1.0000

ALS Bottle#: 1

Method: 8260_15

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8411.D
 Lims ID: STD8260 L5
 Client ID:
 Sample Type: IC Calib Level: 5
 Inject. Date: 21-Jan-2019 11:59:30 ALS Bottle#: 2 Worklist Smp#: 3
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0083824-003
 Misc. Info.: C90121A-IC,8260LLUX15,,43582
 Operator ID: Instrument ID: A3UX15
 Sublist: chrom-8260_15*sub83
 Method: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\8260_15.m
 Limit Group: MSV 8260B ICAL
 Last Update: 22-Jan-2019 09:09:37 Calib Date: 21-Jan-2019 19:15:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8431.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0303

First Level Reviewer: evansle

Date: 21-Jan-2019 12:41:50

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	5.196	5.196	0.000	99	1920375	10.0	10.0	
* 2 Chlorobenzene-d5	117	7.876	7.876	0.000	84	1534484	10.0	10.0	
* 3 1,4-Dichlorobenzene-d4	152	10.117	10.117	0.000	90	858675	10.0	10.0	
\$ 4 Dibromofluoromethane (Surr	113	4.639	4.639	0.000	95	947952	20.0	20.9	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.923	4.923	0.000	100	1190875	20.0	20.4	
\$ 6 Toluene-d8 (Surr)	98	6.560	6.560	0.000	93	3883966	20.0	20.9	
\$ 7 4-Bromofluorobenzene (Surr	95	8.979	8.979	0.000	96	1374232	20.0	20.6	
9 Dichlorodifluoromethane	85	1.532	1.532	0.000	99	1222220	20.0	20.7	
10 Chloromethane	50	1.698	1.686	0.012	99	1207742	20.0	20.1	
11 Vinyl chloride	62	1.793	1.781	0.012	98	1027051	20.0	18.5	
12 Butadiene	54	1.793	1.793	0.000	98	1378009	20.0	18.9	
13 Bromomethane	94	2.065	2.054	0.011	90	791041	20.0	17.8	
14 Chloroethane	64	2.160	2.137	0.023	100	733356	20.0	18.7	
15 Dichlorofluoromethane	67	2.314	2.303	0.011	98	1834172	20.0	18.7	
16 Trichlorofluoromethane	101	2.362	2.362	0.000	99	1703371	20.0	20.5	
17 Ethyl ether	59	2.575	2.563	0.012	90	869536	20.0	19.6	
18 Acrolein	56	2.694	2.694	0.000	99	603583	100.0	99.2	
19 1,1-Dichloroethene	96	2.801	2.777	0.024	98	1125076	20.0	19.5	
20 1,1,2-Trichloro-1,2,2-trif	151	2.801	2.789	0.012	78	967763	20.0	20.4	
22 Acetone	58	2.824	2.812	0.012	100	190030	40.0	43.5	
23 Iodomethane	142	2.955	2.931	0.024	98	1760595	20.0	19.6	
24 Carbon disulfide	76	3.002	2.990	0.012	100	2744114	20.0	18.5	
25 3-Chloro-1-propene	76	3.073	3.073	0.000	88	716236	20.0	22.0	
27 Methyl acetate	43	3.085	3.085	0.000	98	1550490	40.0	39.4	
28 Methylene Chloride	84	3.180	3.180	0.000	92	1097865	20.0	20.6	
29 2-Methyl-2-propanol	59	3.263	3.263	0.000	99	832907	200.0	213.8	
32 Acrylonitrile	53	3.394	3.394	0.000	98	3768272	200.0	204.0	
31 Methyl tert-butyl ether	73	3.405	3.405	0.000	96	3075308	20.0	20.4	
30 trans-1,2-Dichloroethene	96	3.405	3.405	0.000	95	1069131	20.0	19.9	
33 Hexane	86	3.631	3.619	0.012	94	230317	20.0	20.5	

Compound	Sig	RT (min.)	Exp RT (min.)	Diff RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
34 1,1-Dichloroethane	63	3.761	3.761	0.000	96	1773410	20.0	20.4	
36 Vinyl acetate	43	3.785	3.785	0.000	97	2098831	20.0	21.0	
39 cis-1,2-Dichloroethene	96	4.247	4.235	0.012	81	1189038	20.0	20.1	
40 2,2-Dichloropropane	97	4.247	4.247	0.000	65	241048	20.0	20.1	
41 2-Butanone (MEK)	72	4.247	4.247	0.000	99	256717	40.0	40.2	
45 Chlorobromomethane	128	4.449	4.449	0.000	87	595090	20.0	20.2	
46 Tetrahydrofuran	42	4.485	4.485	0.000	87	634219	40.0	39.9	
47 Chloroform	83	4.496	4.496	0.000	93	1793716	20.0	20.3	
48 1,1,1-Trichloroethane	97	4.662	4.662	0.000	98	1356966	20.0	19.9	
49 Cyclohexane	56	4.710	4.710	0.000	90	1592895	20.0	20.8	
51 1,1-Dichloropropene	75	4.793	4.793	0.000	94	1474449	20.0	20.5	
50 Carbon tetrachloride	117	4.805	4.805	0.000	96	1385341	20.0	21.0	
52 Isobutyl alcohol	41	4.864	4.864	0.000	94	821992	500.0	557.2	
53 Benzene	78	4.971	4.971	0.000	95	4352726	20.0	20.3	
54 1,2-Dichloroethane	62	4.983	4.983	0.000	97	1492475	20.0	20.3	
56 n-Heptane	100	5.160	5.160	0.000	92	218160	20.0	21.3	
58 Trichloroethene	130	5.504	5.504	0.000	96	1274454	20.0	20.5	
60 Methylcyclohexane	83	5.670	5.670	0.000	90	1428283	20.0	21.1	
61 1,2-Dichloropropane	63	5.706	5.706	0.000	95	1020861	20.0	20.5	
64 1,4-Dioxane	88	5.813	5.813	0.000	37	183906	400.0	443.3	
63 Dibromomethane	93	5.813	5.813	0.000	88	664429	20.0	20.2	
65 Dichlorobromomethane	83	5.931	5.931	0.000	99	1409597	20.0	21.5	
67 2-Chloroethyl vinyl ether	63	6.180	6.180	0.000	93	1383920	40.0	41.3	
68 cis-1,3-Dichloropropene	75	6.323	6.323	0.000	94	1652896	20.0	21.8	
69 4-Methyl-2-pentanone (MIBK)	43	6.453	6.441	0.012	97	2114735	40.0	41.1	
70 Toluene	91	6.619	6.619	0.000	98	4691174	20.0	20.7	
71 trans-1,3-Dichloropropene	75	6.809	6.809	0.000	94	1578892	20.0	22.1	
72 Ethyl methacrylate	69	6.868	6.868	0.000	90	1421429	20.0	20.9	
73 1,1,2-Trichloroethane	97	6.975	6.975	0.000	91	940456	20.0	20.6	
74 Tetrachloroethene	164	7.117	7.117	0.000	96	1049416	20.0	20.9	
75 1,3-Dichloropropane	76	7.129	7.129	0.000	93	1663194	20.0	20.7	
76 2-Hexanone	43	7.188	7.188	0.000	96	1487023	40.0	42.4	
79 Chlorodibromomethane	129	7.342	7.342	0.000	90	1084488	20.0	22.7	
80 Ethylene Dibromide	107	7.461	7.461	0.000	99	999728	20.0	20.4	
82 Chlorobenzene	112	7.900	7.900	0.000	95	3072375	20.0	20.5	
83 1,1,1,2-Tetrachloroethane	131	7.971	7.971	0.000	95	1113186	20.0	21.0	
84 Ethylbenzene	106	7.995	7.995	0.000	98	1624425	20.0	20.6	
85 m-Xylene & p-Xylene	91	8.101	8.101	0.000	93	3890600	20.0	20.8	
86 o-Xylene	106	8.481	8.481	0.000	96	1968397	20.0	20.7	
87 Styrene	104	8.493	8.493	0.000	93	3394736	20.0	21.4	
88 Bromoform	173	8.682	8.682	0.000	98	770158	20.0	20.3	
89 Isopropylbenzene	105	8.825	8.825	0.000	96	4695119	20.0	20.9	
93 1,1,2,2-Tetrachloroethane	83	9.121	9.121	0.000	97	1238040	20.0	20.3	
92 Bromobenzene	156	9.145	9.133	0.012	89	1450127	20.0	20.4	
94 1,2,3-Trichloropropane	110	9.169	9.169	0.000	84	423192	20.0	20.0	
95 trans-1,4-Dichloro-2-buten	53	9.169	9.169	0.000	75	276245	20.0	31.7	
96 N-Propylbenzene	120	9.228	9.228	0.000	98	1274023	20.0	20.4	
97 2-Chlorotoluene	126	9.323	9.323	0.000	98	1188057	20.0	20.3	
98 1,3,5-Trimethylbenzene	105	9.394	9.394	0.000	94	3784392	20.0	20.8	
99 4-Chlorotoluene	91	9.418	9.418	0.000	98	3844885	20.0	20.3	
101 tert-Butylbenzene	119	9.714	9.714	0.000	92	3285832	20.0	20.2	
103 1,2,4-Trimethylbenzene	105	9.762	9.762	0.000	94	3917785	20.0	20.5	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
105 sec-Butylbenzene	105	9.928	9.928	0.000	93	4155710	20.0	20.8	
106 1,3-Dichlorobenzene	146	10.046	10.046	0.000	98	2462678	20.0	20.1	
107 4-Isopropyltoluene	119	10.070	10.070	0.000	97	3739399	20.0	20.9	
108 1,4-Dichlorobenzene	146	10.141	10.141	0.000	97	2527274	20.0	19.9	
111 n-Butylbenzene	91	10.473	10.473	0.000	97	2918965	20.0	20.6	
112 1,2-Dichlorobenzene	146	10.509	10.509	0.000	99	2381785	20.0	20.0	
113 1,2-Dibromo-3-Chloropropan	157	11.291	11.291	0.000	93	306515	20.0	22.2	
115 1,2,4-Trichlorobenzene	180	12.110	12.110	0.000	94	1510157	20.0	20.0	
116 Hexachlorobutadiene	225	12.276	12.276	0.000	93	696597	20.0	20.4	
117 Naphthalene	128	12.359	12.359	0.000	96	3879317	20.0	20.5	
118 1,2,3-Trichlorobenzene	180	12.608	12.608	0.000	96	1364016	20.0	20.1	
S 127 1,2-Dichloroethene, Total	96				0			40.0	
S 128 1,3-Dichloropropene, Total	75				0			44.0	
S 129 Xylenes, Total	106				0		40.0	41.5	
S 130 Trihalomethanes, Total	1				0		80.0	84.8	
S 157 Total BTEX	1		0.000				100.0	ND	

Reagents:

vm50ss_stk_00079	Amount Added: 16.00	Units: uL
VM50IS_00072	Amount Added: 1.00	Units: uL
VMRPRIMW_00318	Amount Added: 16.00	Units: uL
VMAROLISTDW_00281	Amount Added: 16.00	Units: uL
VMRGAS_00278	Amount Added: 16.00	Units: uL

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8411.D

Injection Date: 21-Jan-2019 11:59:30

Instrument ID: A3UX15

Operator ID:

Lims ID: STD8260 L5

Worklist Smp#: 3

Client ID:

Purge Vol: 5.000 mL

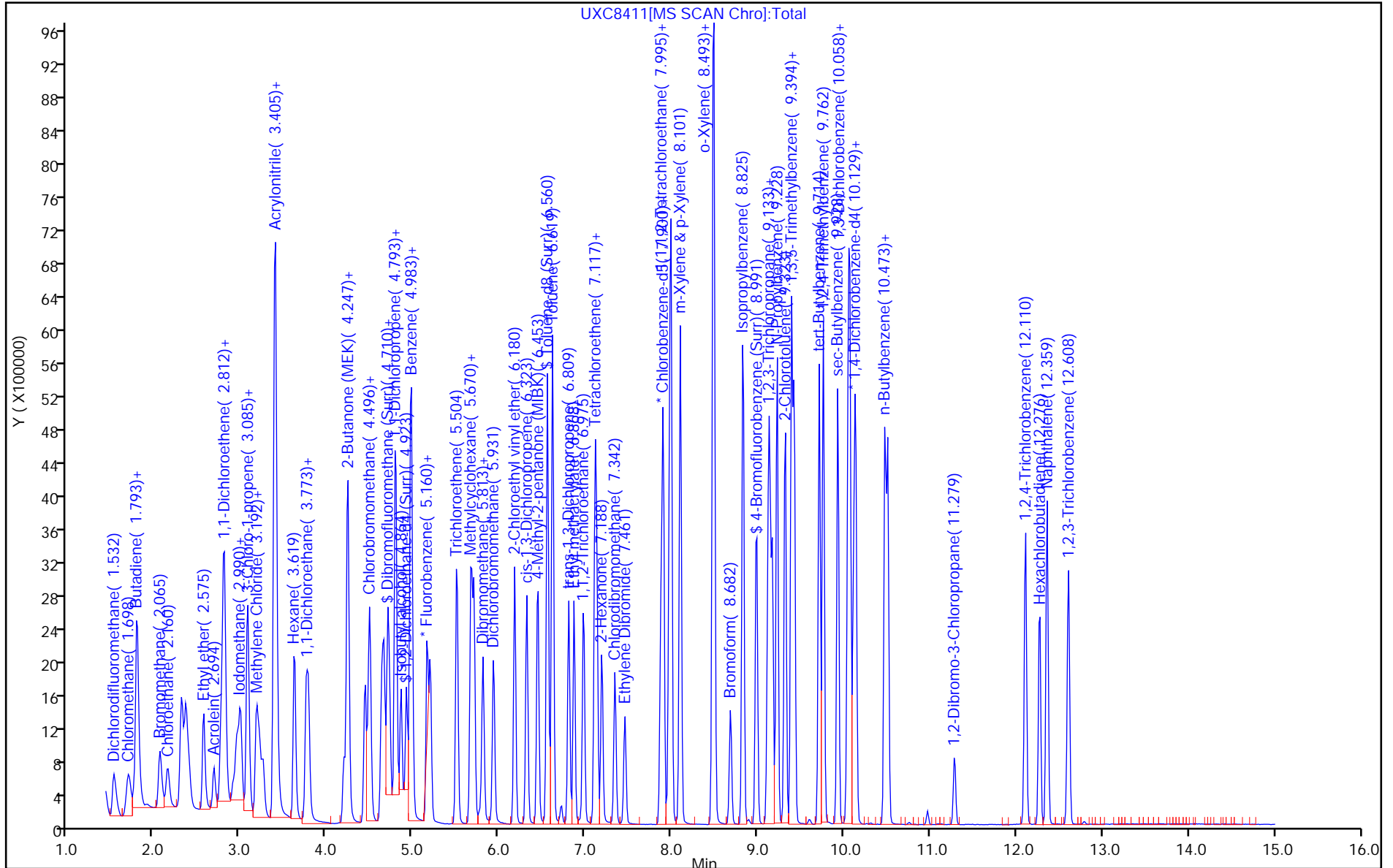
Dil. Factor: 1.0000

ALS Bottle#: 2

Method: 8260_15

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8412.D
 Lims ID: STD8260 L4
 Client ID:
 Sample Type: ICIS Calib Level: 4
 Inject. Date: 21-Jan-2019 12:21:30 ALS Bottle#: 3 Worklist Smp#: 4
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0083824-004
 Misc. Info.: C90121A-IC,8260LLUX15,,43582
 Operator ID: Instrument ID: A3UX15
 Sublist: chrom-8260_15*sub83
 Method: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\8260_15.m
 Limit Group: MSV 8260B ICAL
 Last Update: 22-Jan-2019 09:09:43 Calib Date: 21-Jan-2019 19:15:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8431.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0303

First Level Reviewer: evansle

Date: 21-Jan-2019 12:44:31

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	5.196	5.196	0.000	99	1873821	10.0	10.0	
* 2 Chlorobenzene-d5	117	7.876	7.876	0.000	85	1485792	10.0	10.0	
* 3 1,4-Dichlorobenzene-d4	152	10.117	10.117	0.000	93	831741	10.0	10.0	
\$ 4 Dibromofluoromethane (Surr	113	4.639	4.639	0.000	95	447792	10.0	10.1	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.923	4.923	0.000	100	573781	10.0	10.1	
\$ 6 Toluene-d8 (Surr)	98	6.560	6.560	0.000	93	1816996	10.0	10.1	
\$ 7 4-Bromofluorobenzene (Surr	95	8.979	8.979	0.000	96	649263	10.0	10.1	
9 Dichlorodifluoromethane	85	1.532	1.532	0.000	99	593761	10.0	10.3	
10 Chloromethane	50	1.686	1.686	0.000	99	585358	10.0	9.96	
11 Vinyl chloride	62	1.781	1.781	0.000	98	574312	10.0	10.6	
12 Butadiene	54	1.793	1.793	0.000	98	698736	10.0	9.83	
13 Bromomethane	94	2.054	2.054	0.000	91	423055	10.0	9.78	
14 Chloroethane	64	2.137	2.137	0.000	100	372694	10.0	9.74	
15 Dichlorofluoromethane	67	2.303	2.303	0.000	97	910063	10.0	9.52	
16 Trichlorofluoromethane	101	2.362	2.362	0.000	99	855986	10.0	10.5	
17 Ethyl ether	59	2.563	2.563	0.000	91	419544	10.0	9.69	
18 Acrolein	56	2.694	2.694	0.000	99	290454	50.0	48.9	
19 1,1-Dichloroethene	96	2.777	2.777	0.000	98	516876	10.0	9.18	
20 1,1,2-Trichloro-1,2,2-trif	151	2.789	2.789	0.000	87	462285	10.0	10.0	
22 Acetone	58	2.812	2.812	0.000	99	89168	20.0	19.7	
23 Iodomethane	142	2.931	2.931	0.000	98	909454	10.0	10.4	
24 Carbon disulfide	76	2.990	2.990	0.000	100	1445567	10.0	9.97	
25 3-Chloro-1-propene	76	3.073	3.073	0.000	88	317161	10.0	9.97	
27 Methyl acetate	43	3.085	3.085	0.000	97	723073	20.0	18.8	
28 Methylene Chloride	84	3.180	3.180	0.000	92	580808	10.0	10.0	
29 2-Methyl-2-propanol	59	3.263	3.263	0.000	99	328142	100.0	86.3	
32 Acrylonitrile	53	3.394	3.394	0.000	98	1699336	100.0	94.3	
30 trans-1,2-Dichloroethene	96	3.405	3.405	0.000	97	500088	10.0	9.55	
31 Methyl tert-butyl ether	73	3.405	3.405	0.000	96	1439491	10.0	9.81	
33 Hexane	86	3.619	3.619	0.000	95	104918	10.0	9.55	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
34 1,1-Dichloroethane	63	3.761	3.761	0.000	96	832192	10.0	9.83	
36 Vinyl acetate	43	3.785	3.785	0.000	97	992089	10.0	10.2	
39 cis-1,2-Dichloroethene	96	4.235	4.235	0.000	81	565374	10.0	9.80	
41 2-Butanone (MEK)	72	4.247	4.247	0.000	98	120010	20.0	19.3	
40 2,2-Dichloropropane	97	4.247	4.247	0.000	70	120845	10.0	10.3	
45 Chlorobromomethane	128	4.449	4.449	0.000	88	275468	10.0	9.60	
46 Tetrahydrofuran	42	4.485	4.485	0.000	87	279158	20.0	18.0	
47 Chloroform	83	4.496	4.496	0.000	93	847597	10.0	9.82	
48 1,1,1-Trichloroethane	97	4.662	4.662	0.000	98	689229	10.0	10.4	
49 Cyclohexane	56	4.710	4.710	0.000	91	742256	10.0	9.92	
51 1,1-Dichloropropene	75	4.793	4.793	0.000	94	688186	10.0	9.80	
50 Carbon tetrachloride	117	4.805	4.805	0.000	96	648774	10.0	10.1	
52 Isobutyl alcohol	41	4.864	4.864	0.000	95	322275	250.0	223.9	
53 Benzene	78	4.971	4.971	0.000	96	2024145	10.0	9.68	
54 1,2-Dichloroethane	62	4.983	4.983	0.000	98	707702	10.0	9.89	
56 n-Heptane	100	5.160	5.160	0.000	93	101813	10.0	10.2	
58 Trichloroethene	130	5.504	5.504	0.000	96	593061	10.0	9.77	
60 Methylcyclohexane	83	5.670	5.670	0.000	90	654672	10.0	9.91	
61 1,2-Dichloropropane	63	5.706	5.706	0.000	94	483376	10.0	9.96	
63 Dibromomethane	93	5.813	5.813	0.000	89	317664	10.0	9.87	
64 1,4-Dioxane	88	5.813	5.813	0.000	33	71764	200.0	180.7	
65 Dichlorobromomethane	83	5.931	5.931	0.000	99	637533	10.0	9.98	
67 2-Chloroethyl vinyl ether	63	6.180	6.180	0.000	93	640886	20.0	19.6	
68 cis-1,3-Dichloropropene	75	6.323	6.323	0.000	95	746312	10.0	10.1	
69 4-Methyl-2-pentanone (MIBK)	43	6.441	6.441	0.000	97	970866	20.0	19.3	
70 Toluene	91	6.619	6.619	0.000	98	2172974	10.0	9.91	
71 trans-1,3-Dichloropropene	75	6.809	6.809	0.000	94	706531	10.0	10.2	
72 Ethyl methacrylate	69	6.868	6.868	0.000	89	647349	10.0	9.81	
73 1,1,2-Trichloroethane	97	6.975	6.975	0.000	91	437245	10.0	9.87	
74 Tetrachloroethene	164	7.117	7.117	0.000	96	482895	10.0	9.95	
75 1,3-Dichloropropane	76	7.129	7.129	0.000	93	779324	10.0	10.0	
76 2-Hexanone	43	7.188	7.188	0.000	97	666112	20.0	19.6	
79 Chlorodibromomethane	129	7.342	7.342	0.000	90	469828	10.0	10.1	
80 Ethylene Dibromide	107	7.461	7.461	0.000	98	469490	10.0	9.90	
82 Chlorobenzene	112	7.900	7.900	0.000	95	1441533	10.0	9.92	
83 1,1,1,2-Tetrachloroethane	131	7.971	7.971	0.000	94	522591	10.0	10.2	
84 Ethylbenzene	106	7.995	7.995	0.000	98	758362	10.0	9.92	
85 m-Xylene & p-Xylene	91	8.101	8.101	0.000	93	1818169	10.0	10.0	
86 o-Xylene	106	8.481	8.481	0.000	96	918350	10.0	9.99	
87 Styrene	104	8.493	8.493	0.000	93	1550959	10.0	10.1	
88 Bromoform	173	8.682	8.682	0.000	98	306750	10.0	8.74	
89 Isopropylbenzene	105	8.825	8.825	0.000	95	2200366	10.0	10.1	
93 1,1,2,2-Tetrachloroethane	83	9.121	9.121	0.000	97	577712	10.0	9.80	
92 Bromobenzene	156	9.133	9.133	0.000	89	678570	10.0	9.86	
95 trans-1,4-Dichloro-2-buten	53	9.169	9.169	0.000	75	66363	10.0	7.87	
94 1,2,3-Trichloropropane	110	9.169	9.169	0.000	83	201010	10.0	9.79	
96 N-Propylbenzene	120	9.228	9.228	0.000	98	601438	10.0	9.92	
97 2-Chlorotoluene	126	9.323	9.323	0.000	98	555062	10.0	9.78	
98 1,3,5-Trimethylbenzene	105	9.394	9.394	0.000	94	1766210	10.0	10.0	
99 4-Chlorotoluene	91	9.418	9.418	0.000	98	1798133	10.0	9.80	
101 tert-Butylbenzene	119	9.714	9.714	0.000	91	1551899	10.0	9.83	
103 1,2,4-Trimethylbenzene	105	9.762	9.762	0.000	94	1848645	10.0	9.98	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
105 sec-Butylbenzene	105	9.928	9.928	0.000	94	1935865	10.0	10.0	
106 1,3-Dichlorobenzene	146	10.046	10.046	0.000	99	1168047	10.0	9.86	
107 4-Isopropyltoluene	119	10.070	10.070	0.000	97	1751242	10.0	10.1	
108 1,4-Dichlorobenzene	146	10.141	10.141	0.000	97	1199185	10.0	9.76	
111 n-Butylbenzene	91	10.473	10.473	0.000	97	1384405	10.0	10.1	
112 1,2-Dichlorobenzene	146	10.509	10.509	0.000	99	1159176	10.0	10.0	
113 1,2-Dibromo-3-Chloropropan	157	11.291	11.291	0.000	91	138143	10.0	10.3	
115 1,2,4-Trichlorobenzene	180	12.110	12.110	0.000	93	739836	10.0	10.1	
116 Hexachlorobutadiene	225	12.276	12.276	0.000	93	359850	10.0	9.75	
117 Naphthalene	128	12.359	12.359	0.000	96	1891386	10.0	10.3	
118 1,2,3-Trichlorobenzene	180	12.608	12.608	0.000	95	663704	10.0	10.1	
S 129 Xylenes, Total	106				0		20.0	20.0	
S 157 Total BTEX	1		0.000				50.0	ND	
S 130 Trihalomethanes, Total	1				0		40.0	38.7	

Reagents:

VM50IS_00072	Amount Added: 1.00	Units: uL
vm50ss_stk_00079	Amount Added: 8.00	Units: uL
VMRPRIMW_00318	Amount Added: 8.00	Units: uL
VMAROLISTDW_00281	Amount Added: 8.00	Units: uL
VMRGAS_00278	Amount Added: 8.00	Units: uL



Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8412.D

Injection Date: 21-Jan-2019 12:21:30

Instrument ID: A3UX15

Operator ID:

Lims ID: STD8260 L4

Worklist Smp#: 4

Client ID:

Purge Vol: 5.000 mL

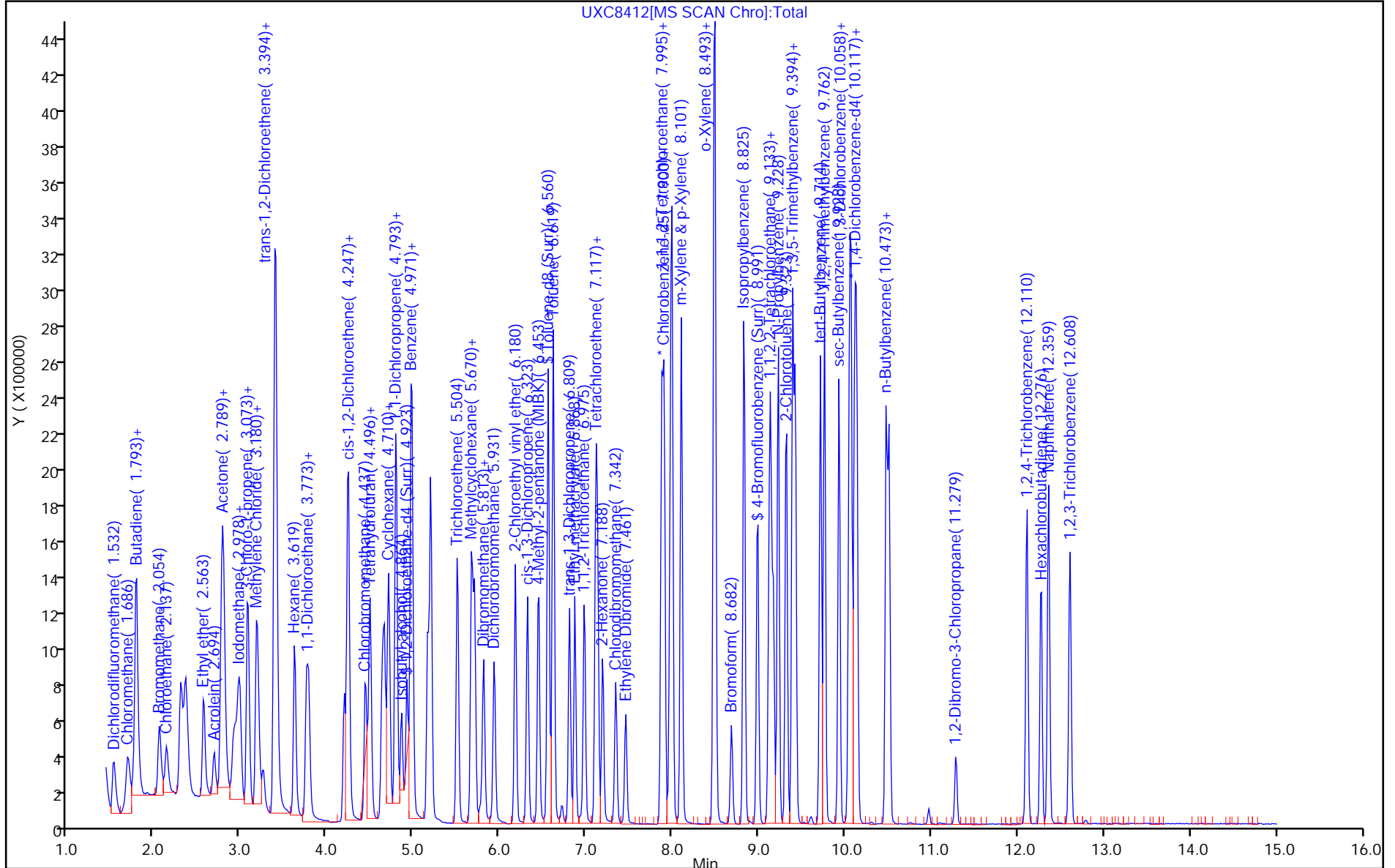
Dil. Factor: 1.0000

ALS Bottle#: 3

Method: 8260_15

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8413.D
 Lims ID: STD8260 L3
 Client ID:
 Sample Type: IC Calib Level: 3
 Inject. Date: 21-Jan-2019 12:43:30 ALS Bottle#: 4 Worklist Smp#: 5
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0083824-005
 Misc. Info.: C90121A-IC,8260LLUX15,,43582
 Operator ID: Instrument ID: A3UX15
 Sublist: chrom-8260_15*sub83
 Method: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\8260_15.m
 Limit Group: MSV 8260B ICAL
 Last Update: 22-Jan-2019 09:09:48 Calib Date: 21-Jan-2019 19:15:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8431.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0303

First Level Reviewer: evansle

Date: 21-Jan-2019 13:20:00

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	5.196	5.196	0.000	99	1838760	10.0	10.0	
* 2 Chlorobenzene-d5	117	7.876	7.876	0.000	85	1476200	10.0	10.0	
* 3 1,4-Dichlorobenzene-d4	152	10.117	10.117	0.000	93	809369	10.0	10.0	
\$ 4 Dibromofluoromethane (Surr	113	4.639	4.639	0.000	95	225163	5.00	5.18	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.923	4.923	0.000	100	288691	5.00	5.16	
\$ 6 Toluene-d8 (Surr)	98	6.560	6.560	0.000	93	912990	5.00	5.10	
\$ 7 4-Bromofluorobenzene (Surr	95	8.979	8.979	0.000	96	328418	5.00	5.13	
9 Dichlorodifluoromethane	85	1.520	1.532	-0.012	99	296202	5.00	5.24	
10 Chloromethane	50	1.686	1.686	0.000	99	276149	5.00	4.79	
11 Vinyl chloride	62	1.781	1.781	0.000	67	261470	5.00	4.93	
12 Butadiene	54	1.793	1.793	0.000	95	361713	5.00	5.18	
13 Bromomethane	94	2.053	2.054	-0.001	91	198900	5.00	4.69	
14 Chloroethane	64	2.136	2.137	-0.001	100	182320	5.00	4.86	
15 Dichlorofluoromethane	67	2.302	2.303	-0.001	98	471934	5.00	5.03	
16 Trichlorofluoromethane	101	2.362	2.362	0.000	98	408598	5.00	5.12	
17 Ethyl ether	59	2.563	2.563	0.000	90	212667	5.00	5.01	
18 Acrolein	56	2.694	2.694	0.000	99	144856	25.0	24.9	
19 1,1-Dichloroethene	96	2.789	2.777	0.012	98	277159	5.00	5.02	
20 1,1,2-Trichloro-1,2,2-trif	151	2.789	2.789	0.000	87	228498	5.00	5.04	
22 Acetone	58	2.812	2.812	0.000	99	50751	10.0	10.4	
23 Iodomethane	142	2.943	2.931	0.012	98	446819	5.00	5.20	
24 Carbon disulfide	76	2.990	2.990	0.000	100	742305	5.00	5.21	
25 3-Chloro-1-propene	76	3.073	3.073	0.000	88	154763	5.00	4.96	
27 Methyl acetate	43	3.085	3.085	0.000	98	380996	10.0	10.1	
28 Methylene Chloride	84	3.180	3.180	0.000	93	338417	5.00	4.93	
29 2-Methyl-2-propanol	59	3.263	3.263	0.000	99	184510	50.0	49.5	
32 Acrylonitrile	53	3.393	3.394	-0.001	98	868795	50.0	49.1	
31 Methyl tert-butyl ether	73	3.405	3.405	0.000	96	714788	5.00	4.96	
30 trans-1,2-Dichloroethene	96	3.405	3.405	0.000	98	251570	5.00	4.89	
33 Hexane	86	3.631	3.619	0.012	94	51427	5.00	4.77	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
34 1,1-Dichloroethane	63	3.761	3.761	0.000	96	409730	5.00	4.93	
36 Vinyl acetate	43	3.785	3.785	0.000	97	476538	5.00	4.99	
39 cis-1,2-Dichloroethene	96	4.235	4.235	0.000	80	278759	5.00	4.92	
40 2,2-Dichloropropane	97	4.247	4.247	0.000	68	58586	5.00	5.11	
41 2-Butanone (MEK)	72	4.247	4.247	0.000	99	59681	10.0	9.76	
45 Chlorobromomethane	128	4.449	4.449	0.000	89	140018	5.00	4.97	
46 Tetrahydrofuran	42	4.484	4.485	-0.001	88	150812	10.0	9.91	
47 Chloroform	83	4.496	4.496	0.000	93	422474	5.00	4.99	
48 1,1,1-Trichloroethane	97	4.662	4.662	0.000	97	334136	5.00	5.12	
49 Cyclohexane	56	4.710	4.710	0.000	90	368473	5.00	5.02	
51 1,1-Dichloropropene	75	4.793	4.793	0.000	93	343617	5.00	4.99	
50 Carbon tetrachloride	117	4.805	4.805	0.000	96	318843	5.00	5.04	
52 Isobutyl alcohol	41	4.864	4.864	0.000	94	177736	125.0	125.8	
53 Benzene	78	4.971	4.971	0.000	96	1016161	5.00	4.95	
54 1,2-Dichloroethane	62	4.982	4.983	-0.001	97	350758	5.00	4.99	
56 n-Heptane	100	5.160	5.160	0.000	93	48491	5.00	4.96	
58 Trichloroethene	130	5.504	5.504	0.000	96	296310	5.00	4.97	
60 Methylcyclohexane	83	5.670	5.670	0.000	89	326169	5.00	5.03	
61 1,2-Dichloropropane	63	5.706	5.706	0.000	94	239856	5.00	5.04	
64 1,4-Dioxane	88	5.813	5.813	0.000	33	39232	100.0	103.2	M
63 Dibromomethane	93	5.813	5.813	0.000	89	157375	5.00	4.99	
65 Dichlorobromomethane	83	5.931	5.931	0.000	99	306036	5.00	4.88	
67 2-Chloroethyl vinyl ether	63	6.180	6.180	0.000	93	319533	10.0	9.96	
68 cis-1,3-Dichloropropene	75	6.322	6.323	-0.001	94	356486	5.00	4.92	
69 4-Methyl-2-pentanone (MIBK)	43	6.441	6.441	0.000	97	495239	10.0	10.1	
70 Toluene	91	6.619	6.619	0.000	98	1080197	5.00	4.96	
71 trans-1,3-Dichloropropene	75	6.809	6.809	0.000	94	333930	5.00	4.87	
72 Ethyl methacrylate	69	6.868	6.868	0.000	89	316654	5.00	4.83	
73 1,1,2-Trichloroethane	97	6.975	6.975	0.000	91	218933	5.00	4.97	
74 Tetrachloroethene	164	7.117	7.117	0.000	96	241473	5.00	5.01	
75 1,3-Dichloropropane	76	7.129	7.129	0.000	92	378264	5.00	4.89	
76 2-Hexanone	43	7.188	7.188	0.000	97	337296	10.0	10.0	
79 Chlorodibromomethane	129	7.342	7.342	0.000	89	222991	5.00	4.85	
80 Ethylene Dibromide	107	7.461	7.461	0.000	97	237763	5.00	5.05	
82 Chlorobenzene	112	7.900	7.900	0.000	95	714138	5.00	4.95	
83 1,1,1,2-Tetrachloroethane	131	7.971	7.971	0.000	93	253355	5.00	4.97	
84 Ethylbenzene	106	7.995	7.995	0.000	98	374569	5.00	4.93	
85 m-Xylene & p-Xylene	91	8.101	8.101	0.000	93	872988	5.00	4.84	
86 o-Xylene	106	8.481	8.481	0.000	96	449671	5.00	4.92	
87 Styrene	104	8.493	8.493	0.000	94	754758	5.00	4.94	
88 Bromoform	173	8.682	8.682	0.000	98	138461	5.00	4.32	
89 Isopropylbenzene	105	8.825	8.825	0.000	96	1070139	5.00	4.95	
93 1,1,2,2-Tetrachloroethane	83	9.121	9.121	0.000	97	283791	5.00	4.95	
92 Bromobenzene	156	9.133	9.133	0.000	89	330598	5.00	4.94	
94 1,2,3-Trichloropropane	110	9.169	9.169	0.000	83	99350	5.00	4.97	
95 trans-1,4-Dichloro-2-buten	53	9.169	9.169	0.000	62	23872	5.00	2.91	
96 N-Propylbenzene	120	9.228	9.228	0.000	98	290505	5.00	4.93	
97 2-Chlorotoluene	126	9.323	9.323	0.000	97	275264	5.00	4.99	
98 1,3,5-Trimethylbenzene	105	9.394	9.394	0.000	94	840604	5.00	4.90	
99 4-Chlorotoluene	91	9.418	9.418	0.000	98	881245	5.00	4.94	
101 tert-Butylbenzene	119	9.714	9.714	0.000	92	749194	5.00	4.88	
103 1,2,4-Trimethylbenzene	105	9.761	9.762	-0.001	94	876431	5.00	4.86	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
105 sec-Butylbenzene	105	9.927	9.928	-0.001	94	937713	5.00	4.98	
106 1,3-Dichlorobenzene	146	10.046	10.046	0.000	98	572540	5.00	4.96	
107 4-Isopropyltoluene	119	10.070	10.070	0.000	97	830781	5.00	4.93	
108 1,4-Dichlorobenzene	146	10.141	10.141	0.000	97	582208	5.00	4.87	
111 n-Butylbenzene	91	10.473	10.473	0.000	97	659506	5.00	4.94	
112 1,2-Dichlorobenzene	146	10.509	10.509	0.000	99	556492	5.00	4.96	
113 1,2-Dibromo-3-Chloropropan	157	11.291	11.291	0.000	90	65389	5.00	5.01	
115 1,2,4-Trichlorobenzene	180	12.109	12.110	-0.001	94	356465	5.00	5.01	
116 Hexachlorobutadiene	225	12.275	12.276	-0.001	94	175613	5.00	4.68	
117 Naphthalene	128	12.358	12.359	-0.001	96	926474	5.00	5.18	
118 1,2,3-Trichlorobenzene	180	12.607	12.608	-0.001	96	324635	5.00	5.08	
S 127 1,2-Dichloroethene, Total	96				0			9.82	
S 128 1,3-Dichloropropene, Total	75				0			9.78	
S 129 Xylenes, Total	106				0		10.0	9.77	
S 130 Trihalomethanes, Total	1				0		20.0	19.0	
S 157 Total BTEX	1		0.000				25.0	ND	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

VM50IS_00072	Amount Added: 1.00	Units: uL
vm50ss_stk_00079	Amount Added: 4.00	Units: uL
VMRPRIMW_00318	Amount Added: 4.00	Units: uL
VMAROLISTDW_00281	Amount Added: 4.00	Units: uL
VMRGAS_00278	Amount Added: 4.00	Units: uL

Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8413.D

Injection Date: 21-Jan-2019 12:43:30

Instrument ID: A3UX15

Operator ID:

Lims ID: STD8260 L3

Worklist Smp#: 5

Client ID:

Purge Vol: 5.000 mL

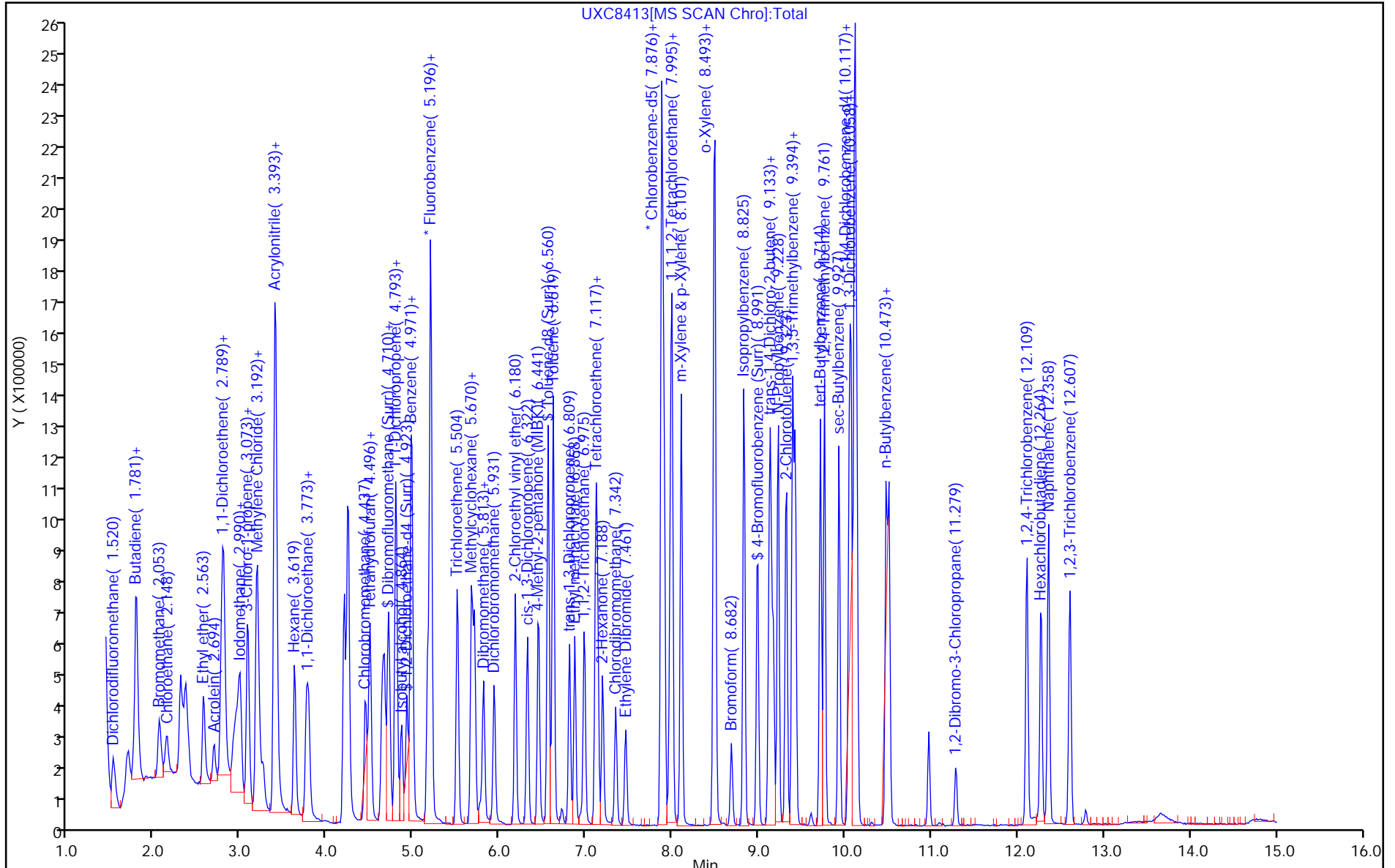
Dil. Factor: 1.0000

ALS Bottle#: 4

Method: 8260_15

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8413.D
Injection Date: 21-Jan-2019 12:43:30 Instrument ID: A3UX15
Lims ID: STD8260 L3
Client ID:
Operator ID:
Purge Vol: 5.000 mL
Method: 8260_15
Column: DB-624 (0.18 mm)

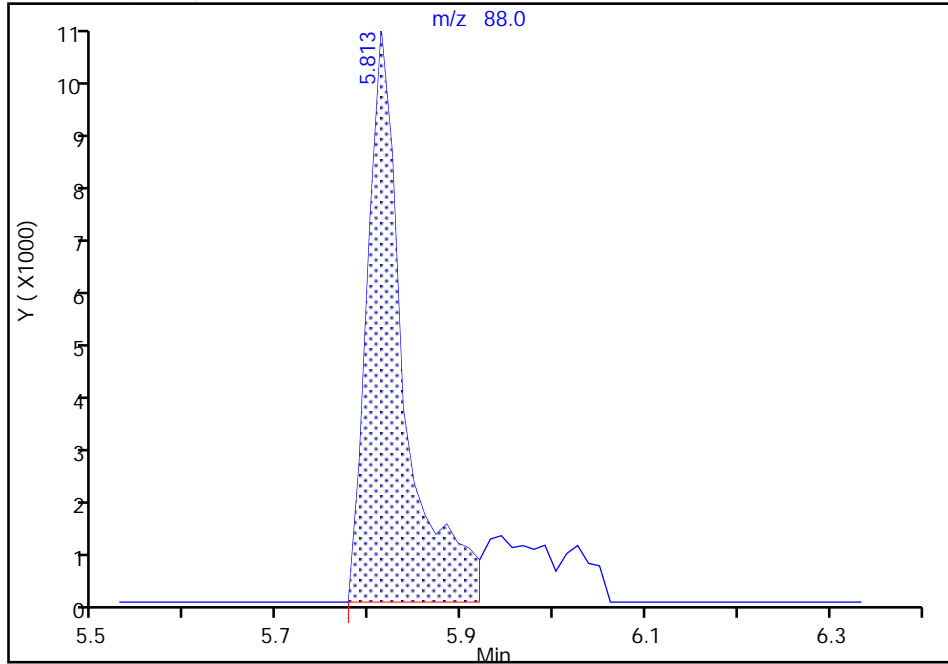
ALS Bottle#: 4 Worklist Smp#: 5
Dil. Factor: 1.0000
Limit Group: MSV 8260B ICAL
Detector MS SCAN

64 1,4-Dioxane, CAS: 123-91-1

Signal: 1

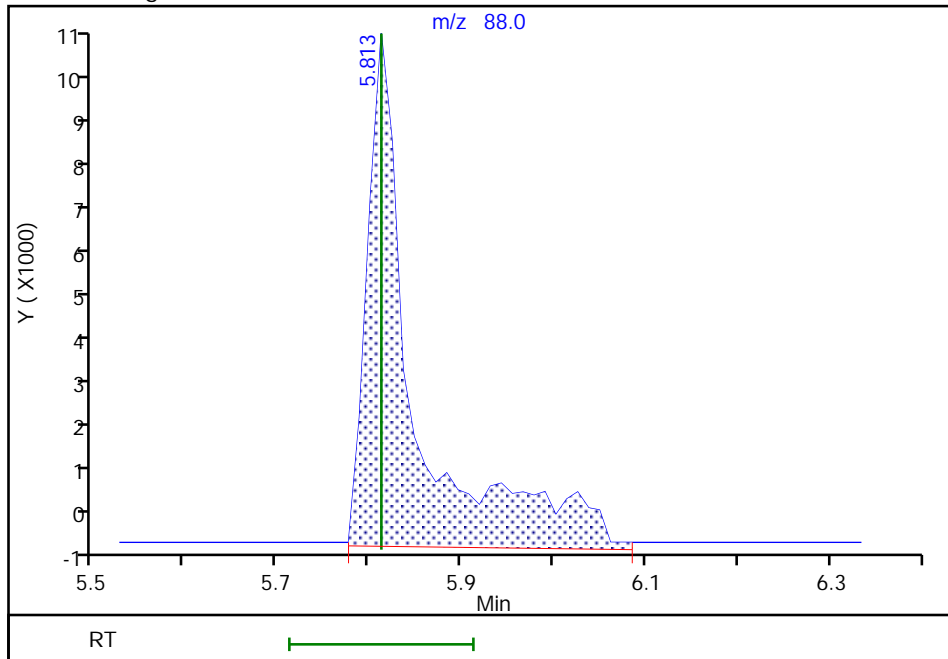
RT: 5.81
Area: 29665
Amount: 80.472140
Amount Units: ug/l

Processing Integration Results



RT: 5.81
Area: 39232
Amount: 103.2242
Amount Units: ug/l

Manual Integration Results



Reviewer: evansle, 21-Jan-2019 13:18:53
Audit Action: Manually Integrated

Audit Reason: Baseline



TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8414.D
 Lims ID: STD8260 L2
 Client ID:
 Sample Type: IC Calib Level: 2
 Inject. Date: 21-Jan-2019 13:05:30 ALS Bottle#: 5 Worklist Smp#: 6
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0083824-006
 Misc. Info.: C90121A-IC,8260LLUX15,,43582
 Operator ID: Instrument ID: A3UX15
 Sublist: chrom-8260_15*sub83
 Method: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\8260_15.m
 Limit Group: MSV 8260B ICAL
 Last Update: 22-Jan-2019 09:09:53 Calib Date: 21-Jan-2019 19:15:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8431.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0303

First Level Reviewer: evansle

Date: 21-Jan-2019 13:21:36

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	5.196	5.196	0.000	99	1827101	10.0	10.0	
* 2 Chlorobenzene-d5	117	7.876	7.876	0.000	84	1454104	10.0	10.0	
* 3 1,4-Dichlorobenzene-d4	152	10.117	10.117	0.000	93	793660	10.0	10.0	
\$ 4 Dibromofluoromethane (Surr	113	4.639	4.639	0.000	94	80448	2.00	1.86	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.923	4.923	0.000	100	106482	2.00	1.92	
\$ 6 Toluene-d8 (Surr)	98	6.560	6.560	0.000	93	333303	2.00	1.89	
\$ 7 4-Bromofluorobenzene (Surr	95	8.979	8.979	0.000	95	120695	2.00	1.91	
9 Dichlorodifluoromethane	85	1.532	1.532	0.000	99	112693	2.00	2.01	
10 Chloromethane	50	1.698	1.686	0.012	99	119628	2.00	2.09	
11 Vinyl chloride	62	1.793	1.781	0.012	62	104647	2.00	1.99	
12 Butadiene	54	1.793	1.793	0.000	96	148568	2.00	2.14	
13 Bromomethane	94	2.065	2.054	0.011	91	100403	2.00	2.38	
14 Chloroethane	64	2.148	2.137	0.011	100	82718	2.00	2.22	
15 Dichlorofluoromethane	67	2.314	2.303	0.011	98	201291	2.00	2.16	
16 Trichlorofluoromethane	101	2.362	2.362	0.000	97	157640	2.00	1.99	
17 Ethyl ether	59	2.575	2.563	0.012	90	87432	2.00	2.07	
18 Acrolein	56	2.694	2.694	0.000	99	55282	10.0	9.55	
19 1,1-Dichloroethene	96	2.801	2.777	0.024	98	120279	2.00	2.19	
20 1,1,2-Trichloro-1,2,2-trif	151	2.801	2.789	0.012	77	91700	2.00	2.03	
22 Acetone	58	2.812	2.812	0.000	98	23252	4.00	3.50	
23 Iodomethane	142	2.955	2.931	0.024	98	156089	2.00	1.83	
24 Carbon disulfide	76	3.002	2.990	0.012	100	283969	2.00	2.01	
25 3-Chloro-1-propene	76	3.073	3.073	0.000	86	58477	2.00	1.88	
27 Methyl acetate	43	3.085	3.085	0.000	97	150240	4.00	4.01	
28 Methylene Chloride	84	3.180	3.180	0.000	96	198393	2.00	1.89	
29 2-Methyl-2-propanol	59	3.263	3.263	0.000	99	82343	20.0	22.2	
32 Acrylonitrile	53	3.393	3.394	-0.001	100	352291	20.0	20.0	
30 trans-1,2-Dichloroethene	96	3.405	3.405	0.000	96	103155	2.00	2.02	
31 Methyl tert-butyl ether	73	3.405	3.405	0.000	96	285516	2.00	1.99	
33 Hexane	86	3.631	3.619	0.012	93	22194	2.00	2.07	

Compound	Sig	RT (min.)	Exp RT (min.)	Diff RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
34 1,1-Dichloroethane	63	3.761	3.761	0.000	96	166930	2.00	2.02	
36 Vinyl acetate	43	3.785	3.785	0.000	97	177639	2.00	1.87	
39 cis-1,2-Dichloroethene	96	4.247	4.235	0.012	81	116480	2.00	2.07	
41 2-Butanone (MEK)	72	4.247	4.247	0.000	72	24387	4.00	4.01	
40 2,2-Dichloropropane	97	4.247	4.247	0.000	60	21907	2.00	1.92	
45 Chlorobromomethane	128	4.449	4.449	0.000	90	56716	2.00	2.03	
46 Tetrahydrofuran	42	4.484	4.485	-0.001	86	62314	4.00	4.12	
47 Chloroform	83	4.496	4.496	0.000	94	169874	2.00	2.02	
48 1,1,1-Trichloroethane	97	4.662	4.662	0.000	97	130349	2.00	2.01	
49 Cyclohexane	56	4.710	4.710	0.000	90	144320	2.00	1.98	
51 1,1-Dichloropropene	75	4.793	4.793	0.000	94	140972	2.00	2.06	
50 Carbon tetrachloride	117	4.805	4.805	0.000	97	124489	2.00	1.98	
52 Isobutyl alcohol	41	4.864	4.864	0.000	96	73133	50.0	52.1	
53 Benzene	78	4.971	4.971	0.000	95	412777	2.00	2.02	
54 1,2-Dichloroethane	62	4.983	4.983	0.000	98	142414	2.00	2.04	
56 n-Heptane	100	5.160	5.160	0.000	92	19095	2.00	1.96	
58 Trichloroethene	130	5.504	5.504	0.000	96	120531	2.00	2.04	
60 Methylcyclohexane	83	5.670	5.670	0.000	90	127322	2.00	1.98	
61 1,2-Dichloropropane	63	5.706	5.706	0.000	94	97800	2.00	2.07	
63 Dibromomethane	93	5.813	5.813	0.000	89	65366	2.00	2.08	
64 1,4-Dioxane	88	5.813	5.813	0.000	37	14227	40.0	41.3	
65 Dichlorobromomethane	83	5.931	5.931	0.000	98	121869	2.00	1.96	
67 2-Chloroethyl vinyl ether	63	6.180	6.180	0.000	93	123378	4.00	3.87	
68 cis-1,3-Dichloropropene	75	6.323	6.323	0.000	94	138202	2.00	1.92	
69 4-Methyl-2-pentanone (MIBK)	43	6.453	6.441	0.012	96	187250	4.00	3.83	
70 Toluene	91	6.619	6.619	0.000	98	428381	2.00	2.00	
71 trans-1,3-Dichloropropene	75	6.809	6.809	0.000	95	125256	2.00	1.85	
72 Ethyl methacrylate	69	6.868	6.868	0.000	89	129112	2.00	2.00	
73 1,1,2-Trichloroethane	97	6.975	6.975	0.000	91	87521	2.00	2.02	
74 Tetrachloroethene	164	7.117	7.117	0.000	96	95990	2.00	2.02	
75 1,3-Dichloropropane	76	7.129	7.129	0.000	92	153833	2.00	2.02	
76 2-Hexanone	43	7.188	7.188	0.000	97	135481	4.00	4.08	
79 Chlorodibromomethane	129	7.342	7.342	0.000	90	82891	2.00	1.83	
80 Ethylene Dibromide	107	7.461	7.461	0.000	98	91132	2.00	1.96	
82 Chlorobenzene	112	7.900	7.900	0.000	97	286159	2.00	2.01	
83 1,1,1,2-Tetrachloroethane	131	7.971	7.971	0.000	93	95060	2.00	1.89	
84 Ethylbenzene	106	7.995	7.995	0.000	98	146696	2.00	1.96	
85 m-Xylene & p-Xylene	91	8.101	8.101	0.000	93	351787	2.00	1.98	
86 o-Xylene	106	8.481	8.481	0.000	96	176986	2.00	1.97	
87 Styrene	104	8.493	8.493	0.000	91	289260	2.00	1.92	
88 Bromoform	173	8.682	8.682	0.000	97	51580	2.00	2.03	
89 Isopropylbenzene	105	8.825	8.825	0.000	95	416807	2.00	1.96	
93 1,1,2,2-Tetrachloroethane	83	9.121	9.121	0.000	97	112833	2.00	2.01	
92 Bromobenzene	156	9.133	9.133	0.000	89	131876	2.00	2.01	
95 trans-1,4-Dichloro-2-buten	53	9.169	9.169	0.000	67	16039	2.00	1.99	
94 1,2,3-Trichloropropane	110	9.169	9.169	0.000	86	40526	2.00	2.07	
96 N-Propylbenzene	120	9.228	9.228	0.000	98	117423	2.00	2.03	
97 2-Chlorotoluene	126	9.323	9.323	0.000	98	107735	2.00	1.99	
98 1,3,5-Trimethylbenzene	105	9.394	9.394	0.000	94	332591	2.00	1.98	
99 4-Chlorotoluene	91	9.418	9.418	0.000	98	356628	2.00	2.04	
101 tert-Butylbenzene	119	9.714	9.714	0.000	92	300630	2.00	2.00	
103 1,2,4-Trimethylbenzene	105	9.762	9.762	0.000	94	359918	2.00	2.04	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
105 sec-Butylbenzene	105	9.928	9.928	0.000	94	373619	2.00	2.02	
106 1,3-Dichlorobenzene	146	10.046	10.046	0.000	98	229119	2.00	2.03	
107 4-Isopropyltoluene	119	10.070	10.070	0.000	96	334169	2.00	2.02	
108 1,4-Dichlorobenzene	146	10.141	10.141	0.000	97	241629	2.00	2.06	
111 n-Butylbenzene	91	10.473	10.473	0.000	97	263153	2.00	2.01	
112 1,2-Dichlorobenzene	146	10.509	10.509	0.000	98	224899	2.00	2.04	
113 1,2-Dibromo-3-Chloropropan	157	11.291	11.291	0.000	88	24334	2.00	1.90	
115 1,2,4-Trichlorobenzene	180	12.109	12.110	-0.001	93	145803	2.00	2.09	
116 Hexachlorobutadiene	225	12.276	12.276	0.000	94	78276	2.00	2.11	
117 Naphthalene	128	12.359	12.359	0.000	96	348942	2.00	1.99	
118 1,2,3-Trichlorobenzene	180	12.608	12.608	0.000	95	131338	2.00	2.09	
S 127 1,2-Dichloroethene, Total	96				0			4.09	
S 128 1,3-Dichloropropene, Total	75				0			3.77	
S 129 Xylenes, Total	106				0		4.00	3.95	
S 157 Total BTEX	1		0.000				10.0	ND	
S 130 Trihalomethanes, Total	1				0		8.00	7.83	

Reagents:

VM50IS_00072	Amount Added: 1.00	Units: uL
vm50ss_stk_00079	Amount Added: 1.60	Units: uL
VMRPRIMW_00318	Amount Added: 1.60	Units: uL
VMAROLISTDW_00281	Amount Added: 1.60	Units: uL
VMRGAS_00278	Amount Added: 1.60	Units: uL

Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8414.D

Injection Date: 21-Jan-2019 13:05:30

Instrument ID: A3UX15

Operator ID:

Lims ID: STD8260 L2

Worklist Smp#: 6

Client ID:

Purge Vol: 5.000 mL

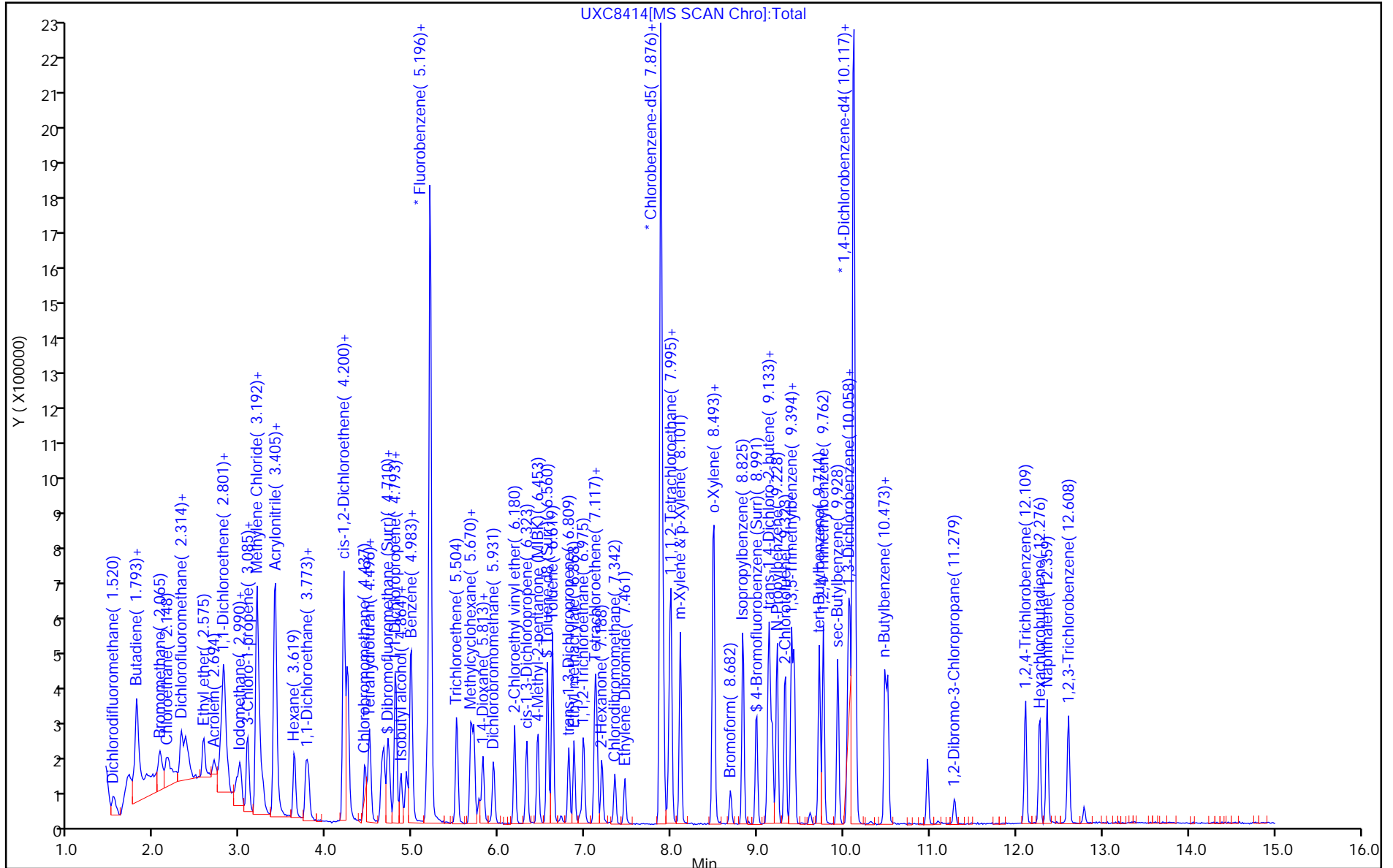
Dil. Factor: 1.0000

ALS Bottle#: 5

Method: 8260_15

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8415.D
 Lims ID: STD8260 L1
 Client ID:
 Sample Type: IC Calib Level: 1
 Inject. Date: 21-Jan-2019 13:26:30 ALS Bottle#: 6 Worklist Smp#: 7
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0083824-007
 Misc. Info.: C90121A-IC,8260LLUX15,,43582
 Operator ID: Instrument ID: A3UX15
 Sublist: chrom-8260_15*sub83
 Method: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\8260_15.m
 Limit Group: MSV 8260B ICAL
 Last Update: 22-Jan-2019 09:09:57 Calib Date: 21-Jan-2019 19:15:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8431.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0303

First Level Reviewer: evansle

Date: 21-Jan-2019 13:51:17

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	5.196	5.196	0.000	99	1764972	10.0	10.0	
* 2 Chlorobenzene-d5	117	7.876	7.876	0.000	84	1396761	10.0	10.0	
* 3 1,4-Dichlorobenzene-d4	152	10.117	10.117	0.000	94	755367	10.0	10.0	
\$ 4 Dibromofluoromethane (Surr	113	4.639	4.639	0.000	94	41840	1.00	1.00	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.923	4.923	0.000	100	55383	1.00	1.03	
\$ 6 Toluene-d8 (Surr)	98	6.560	6.560	0.000	93	174193	1.00	1.03	
\$ 7 4-Bromofluorobenzene (Surr	95	8.979	8.979	0.000	95	60446	1.00	1.00	
9 Dichlorodifluoromethane	85	1.532	1.532	0.000	99	47729	1.00	0.8803	
10 Chloromethane	50	1.698	1.686	0.012	96	58046	1.00	1.05	
11 Vinyl chloride	62	1.793	1.781	0.012	72	57558	1.00	1.13	
12 Butadiene	54	1.793	1.793	0.000	95	71088	1.00	1.06	
13 Bromomethane	94	2.054	2.054	0.000	88	47411	1.00	1.16	
14 Chloroethane	64	2.137	2.137	0.000	99	40558	1.00	1.13	
15 Dichlorofluoromethane	67	2.303	2.303	0.000	97	103682	1.00	1.15	
16 Trichlorofluoromethane	101	2.350	2.362	-0.012	98	68568	1.00	0.8959	
17 Ethyl ether	59	2.575	2.563	0.012	91	44605	1.00	1.09	
18 Acrolein	56	2.694	2.694	0.000	97	32229	5.00	5.76	
19 1,1-Dichloroethene	96	2.777	2.777	0.000	98	57094	1.00	1.08	
20 1,1,2-Trichloro-1,2,2-trif	151	2.789	2.789	0.000	88	42550	1.00	0.9770	
22 Acetone	58	2.813	2.812	0.001	98	17181	2.00	2.11	
23 Iodomethane	142	2.931	2.931	0.000	99	93588	1.00	1.13	
24 Carbon disulfide	76	2.990	2.990	0.000	97	160687	1.00	1.18	
25 3-Chloro-1-propene	76	3.073	3.073	0.000	89	27003	1.00	0.9008	
27 Methyl acetate	43	3.085	3.085	0.000	98	79740	2.00	2.20	
28 Methylene Chloride	84	3.180	3.180	0.000	93	154915	1.00	1.05	
29 2-Methyl-2-propanol	59	3.263	3.263	0.000	99	36240	10.0	10.1	
32 Acrylonitrile	53	3.394	3.394	0.000	97	175156	10.0	10.3	
31 Methyl tert-butyl ether	73	3.405	3.405	0.000	96	139955	1.00	1.01	
30 trans-1,2-Dichloroethene	96	3.405	3.405	0.000	97	54736	1.00	1.11	
33 Hexane	86	3.619	3.619	0.000	94	10725	1.00	1.04	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
34 1,1-Dichloroethane	63	3.761	3.761	0.000	97	81126	1.00	1.02	
36 Vinyl acetate	43	3.785	3.785	0.000	97	88027	1.00	0.9595	
39 cis-1,2-Dichloroethene	96	4.236	4.235	0.001	80	57035	1.00	1.05	
40 2,2-Dichloropropane	97	4.247	4.247	0.000	42	11684	1.00	1.06	
41 2-Butanone (MEK)	72	4.247	4.247	0.000	57	12505	2.00	2.13	
45 Chlorobromomethane	128	4.437	4.449	-0.012	88	28543	1.00	1.06	
46 Tetrahydrofuran	42	4.485	4.485	0.000	86	31985	2.00	2.19	
47 Chloroform	83	4.496	4.496	0.000	93	83929	1.00	1.03	
48 1,1,1-Trichloroethane	97	4.662	4.662	0.000	98	63803	1.00	1.02	
49 Cyclohexane	56	4.710	4.710	0.000	91	68915	1.00	0.9779	
51 1,1-Dichloropropene	75	4.793	4.793	0.000	93	66437	1.00	1.00	
50 Carbon tetrachloride	117	4.805	4.805	0.000	97	56945	1.00	0.9377	
52 Isobutyl alcohol	41	4.864	4.864	0.000	95	31904	25.0	23.5	
53 Benzene	78	4.971	4.971	0.000	96	207184	1.00	1.05	
54 1,2-Dichloroethane	62	4.983	4.983	0.000	97	68031	1.00	1.01	
56 n-Heptane	100	5.161	5.160	0.000	80	8801	1.00	0.9370	a
58 Trichloroethene	130	5.504	5.504	0.000	95	59121	1.00	1.03	
60 Methylcyclohexane	83	5.670	5.670	0.000	88	60130	1.00	0.9667	
61 1,2-Dichloropropane	63	5.706	5.706	0.000	90	43707	1.00	0.9560	
64 1,4-Dioxane	88	5.825	5.813	0.012	33	5185	20.0	19.2	
63 Dibromomethane	93	5.813	5.813	0.000	86	31111	1.00	1.03	
65 Dichlorobromomethane	83	5.931	5.931	0.000	98	57382	1.00	0.9535	
67 2-Chloroethyl vinyl ether	63	6.180	6.180	0.000	93	61442	2.00	1.99	
68 cis-1,3-Dichloropropene	75	6.323	6.323	0.000	93	62099	1.00	0.8920	
69 4-Methyl-2-pentanone (MIBK)	43	6.453	6.441	0.012	98	98299	2.00	2.08	
70 Toluene	91	6.619	6.619	0.000	98	209687	1.00	1.02	
71 trans-1,3-Dichloropropene	75	6.809	6.809	0.000	94	57707	1.00	0.8893	
72 Ethyl methacrylate	69	6.868	6.868	0.000	90	61421	1.00	0.9899	
73 1,1,2-Trichloroethane	97	6.975	6.975	0.000	91	41853	1.00	1.00	
74 Tetrachloroethene	164	7.117	7.117	0.000	97	44889	1.00	0.9835	
75 1,3-Dichloropropane	76	7.129	7.129	0.000	92	72217	1.00	0.9861	
76 2-Hexanone	43	7.188	7.188	0.000	96	58616	2.00	1.84	
79 Chlorodibromomethane	129	7.342	7.342	0.000	89	36977	1.00	0.8497	
80 Ethylene Dibromide	107	7.461	7.461	0.000	97	45295	1.00	1.02	
82 Chlorobenzene	112	7.900	7.900	0.000	96	138107	1.00	1.01	
83 1,1,1,2-Tetrachloroethane	131	7.971	7.971	0.000	93	46702	1.00	0.9691	
84 Ethylbenzene	106	7.995	7.995	0.000	98	73699	1.00	1.03	
85 m-Xylene & p-Xylene	91	8.101	8.101	0.000	93	168049	1.00	0.9855	
86 o-Xylene	106	8.481	8.481	0.000	97	88147	1.00	1.02	
87 Styrene	104	8.493	8.493	0.000	93	137740	1.00	0.9520	
88 Bromoform	173	8.682	8.682	0.000	96	19867	1.00	1.20	
89 Isopropylbenzene	105	8.825	8.825	0.000	95	198841	1.00	0.9729	
93 1,1,2,2-Tetrachloroethane	83	9.121	9.121	0.000	97	54865	1.00	1.03	
92 Bromobenzene	156	9.133	9.133	0.000	90	63640	1.00	1.02	
94 1,2,3-Trichloropropane	110	9.169	9.169	0.000	85	18846	1.00	1.01	
95 trans-1,4-Dichloro-2-buten	53	9.169	9.169	0.000	62	3277	1.00	0.4279	
96 N-Propylbenzene	120	9.228	9.228	0.000	99	54874	1.00	1.00	
97 2-Chlorotoluene	126	9.323	9.323	0.000	97	53702	1.00	1.04	
98 1,3,5-Trimethylbenzene	105	9.394	9.394	0.000	94	160683	1.00	1.00	
99 4-Chlorotoluene	91	9.418	9.418	0.000	98	170473	1.00	1.02	
101 tert-Butylbenzene	119	9.714	9.714	0.000	92	141080	1.00	0.9843	
103 1,2,4-Trimethylbenzene	105	9.762	9.762	0.000	94	169136	1.00	1.01	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/l	OnCol Amt ug/l	Flags
105 sec-Butylbenzene	105	9.928	9.928	0.000	94	173094	1.00	0.9849	
106 1,3-Dichlorobenzene	146	10.046	10.046	0.000	97	113176	1.00	1.05	
107 4-Isopropyltoluene	119	10.070	10.070	0.000	95	154318	1.00	0.9808	
108 1,4-Dichlorobenzene	146	10.141	10.141	0.000	96	119243	1.00	1.07	
111 n-Butylbenzene	91	10.473	10.473	0.000	97	127428	1.00	1.02	
112 1,2-Dichlorobenzene	146	10.509	10.509	0.000	98	109056	1.00	1.04	
113 1,2-Dibromo-3-Chloropropan	157	11.291	11.291	0.000	84	10572	1.00	0.8688	
115 1,2,4-Trichlorobenzene	180	12.110	12.110	0.000	94	73453	1.00	1.11	
116 Hexachlorobutadiene	225	12.276	12.276	0.000	93	41867	1.00	1.20	
117 Naphthalene	128	12.359	12.359	0.000	96	172653	1.00	1.03	
118 1,2,3-Trichlorobenzene	180	12.608	12.608	0.000	95	65430	1.00	1.10	
S 127 1,2-Dichloroethene, Total	96				0			2.16	
S 128 1,3-Dichloropropene, Total	75				0			1.78	
S 129 Xylenes, Total	106				0		2.00	2.01	
S 130 Trihalomethanes, Total	1				0		4.00	4.03	
S 157 Total BTEX	1		0.000				5.00	ND	

QC Flag Legend

Review Flags

a - User Assigned ID

Reagents:

VM50IS_00072	Amount Added: 1.00	Units: uL
vm50ss_stk_00079	Amount Added: 0.80	Units: uL
VMRPRIMW_00318	Amount Added: 0.80	Units: uL
VMAROLISTDW_00281	Amount Added: 0.80	Units: uL
VMRGAS_00278	Amount Added: 0.80	Units: uL

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8415.D

Injection Date: 21-Jan-2019 13:26:30

Instrument ID: A3UX15

Operator ID:

Lims ID: STD8260 L1

Worklist Smp#: 7

Client ID:

Purge Vol: 5.000 mL

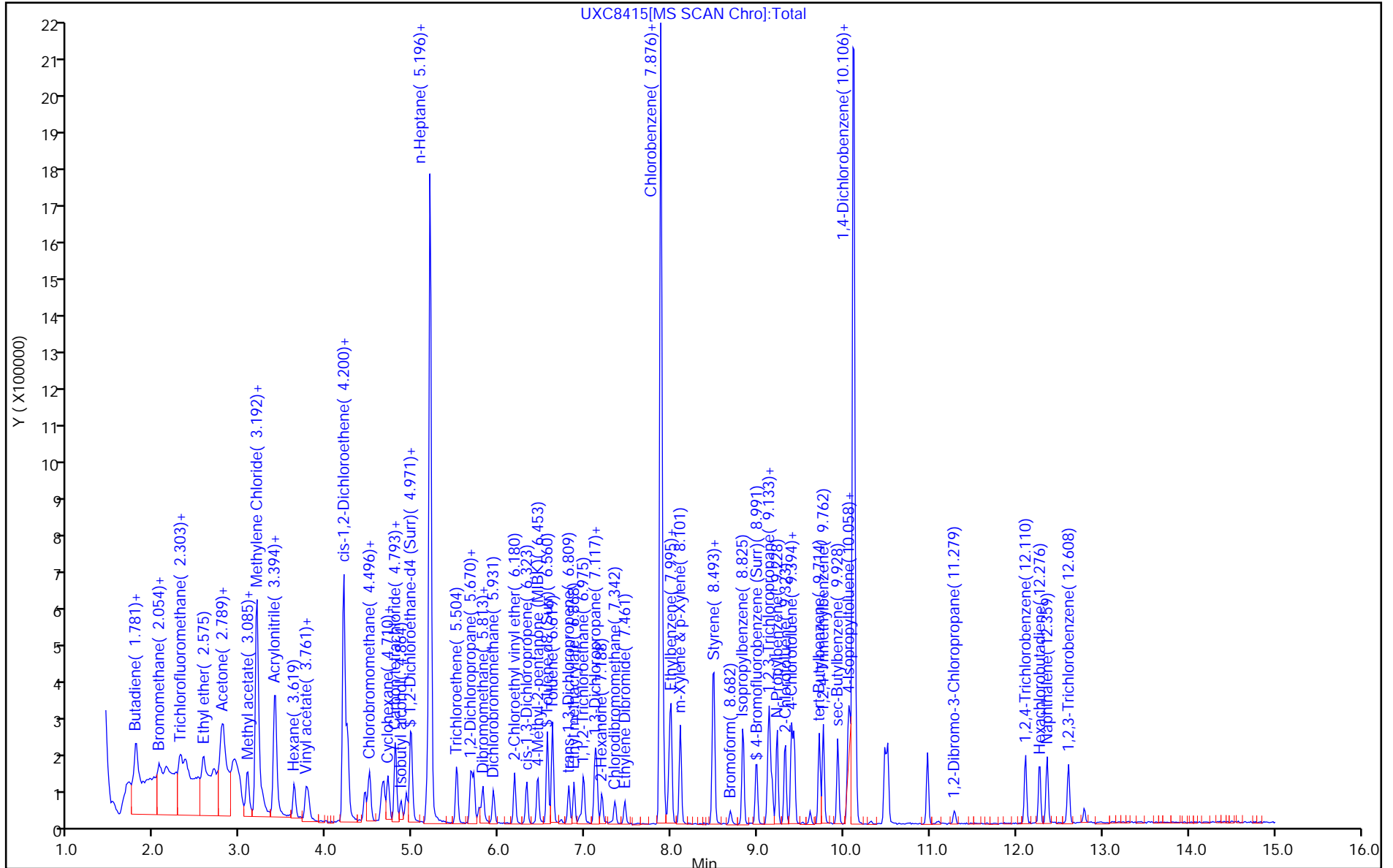
Dil. Factor: 1.0000

ALS Bottle#: 6

Method: 8260_15

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton

Data File:	\\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8415.D		
Injection Date:	21-Jan-2019 13:26:30	Instrument ID:	A3UX15
Lims ID:	STD8260 L1		
Client ID:			
Operator ID:		ALS Bottle#:	6
Purge Vol:	5.000 mL	Dil. Factor:	1.0000
Method:	8260_15	Limit Group:	MSV 8260B ICAL
Column:	DB-624 (0.18 mm)	Detector:	MS SCAN

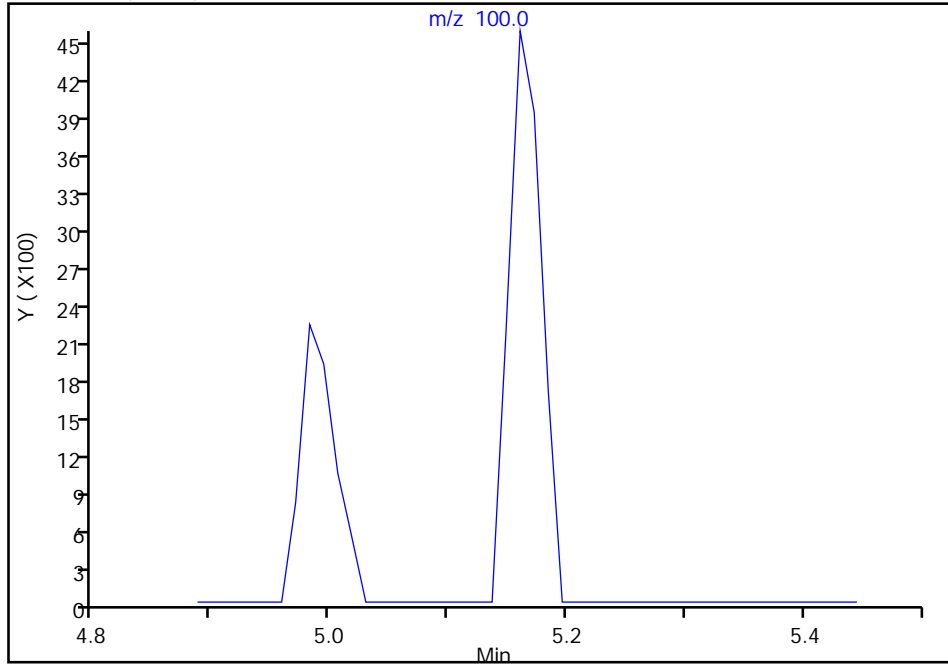
Worklist Smp#: 7

56 n-Heptane, CAS: 142-82-5

Signal: 1

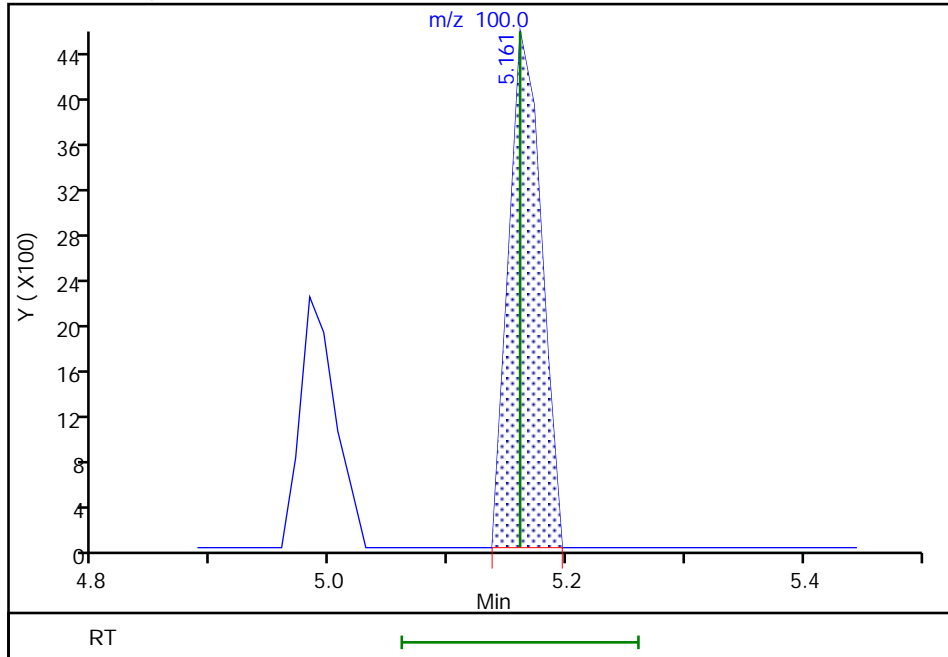
Not Detected
Expected RT: 5.16

Processing Integration Results



Manual Integration Results

RT: 5.16
 Area: 8801
 Amount: 0.936957
 Amount Units: ug/l



Reviewer: evansle, 21-Jan-2019 13:47:12

Audit Action: Assigned Compound ID

Audit Reason: Peak assignment corrected

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Client Sample ID: MW-72_030519 Lab Sample ID: 240-109102-1
 Matrix: Water Lab File ID: UXC9436.D
 Analysis Method: 8260B Date Collected: 03/05/2019 14:20
 Sample wt/vol: 5(mL) Date Analyzed: 03/18/2019 16:27
 Soil Aliquot Vol: _____ Dilution Factor: 1
 Soil Extract Vol.: _____ GC Column: DB-624 ID: 0.18 (mm)
 % Moisture: _____ Level: (low/med) Low
 Analysis Batch No.: 372032 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-35-4	1,1-Dichloroethene	1.0	U	1.0	0.19
156-59-2	cis-1,2-Dichloroethene	1.0	U	1.0	0.16
127-18-4	Tetrachloroethene	1.0	U	1.0	0.15
156-60-5	trans-1,2-Dichloroethene	1.0	U	1.0	0.19
79-01-6	Trichloroethene	1.0	U	1.0	0.10
75-01-4	Vinyl chloride	1.9		1.0	0.20

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	89		70-121
460-00-4	4-Bromofluorobenzene (Surr)	68		59-120
2037-26-5	Toluene-d8 (Surr)	79		70-123
1868-53-7	Dibromofluoromethane (Surr)	91		75-128

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX15\20190318-85248.b\UXC9436.D
 Lims ID: 240-109102-B-1
 Client ID: MW-72_030519
 Sample Type: Client
 Inject. Date: 18-Mar-2019 16:27:30 ALS Bottle#: 16 Worklist Smp#: 41
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0085248-041
 Misc. Info.: C90318A,8260LLUX15,,43582
 Operator ID: Instrument ID: A3UX15
 Method: \\chromna\Canton\ChromData\A3UX15\20190318-85248.b\8260_15.m
 Limit Group: MSV 8260B ICAL
 Last Update: 19-Mar-2019 08:40:05 Calib Date: 21-Jan-2019 19:15:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8431.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0304

First Level Reviewer: evansle

Date: 19-Mar-2019 08:30:58

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	5.196	5.196	0.000	99	1808468	10.0	
* 2 Chlorobenzene-d5	117	7.876	7.876	0.000	84	1429308	10.0	
* 3 1,4-Dichlorobenzene-d4	152	10.117	10.117	0.000	93	724641	10.0	
\$ 4 Dibromofluoromethane (Surr	113	4.639	4.639	0.000	95	309482	7.23	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.923	4.923	0.000	99	389940	7.09	
\$ 6 Toluene-d8 (Surr)	98	6.560	6.560	0.000	93	1086195	6.26	
\$ 7 4-Bromofluorobenzene (Surr	95	8.991	8.991	0.000	96	338549	5.46	
11 Vinyl chloride	62	1.781	1.769	0.012	98	98713	1.89	
19 1,1-Dichloroethene	96		2.777				ND	
30 trans-1,2-Dichloroethene	96		3.405				ND	
39 cis-1,2-Dichloroethene	96	4.247	4.236	0.011	77	7745	0.1391	
58 Trichloroethene	130		5.504				ND	
74 Tetrachloroethene	164	7.117	7.117	0.000	95	5889	0.1261	

Reagents:

vm40ml_vials_00009	Amount Added: 0.00	Units:	Run Reagent
VM50IS_00075	Amount Added: 1.00	Units: uL	Run Reagent
vm50ss_stk_00080	Amount Added: 0.80	Units: uL	Run Reagent
vmDist_H2o_00140	Amount Added: 0.00	Units:	Run Reagent

Data File: \\chromna\Canton\ChromData\A3UX15\20190318-85248.b\UXC9436.D

Injection Date: 18-Mar-2019 16:27:30

Instrument ID: A3UX15

Operator ID:

Lims ID: 240-109102-B-1

Lab Sample ID: 240-109102-1

Worklist Smp#: 41

Client ID: MW-72_030519

Purge Vol: 5.000 mL

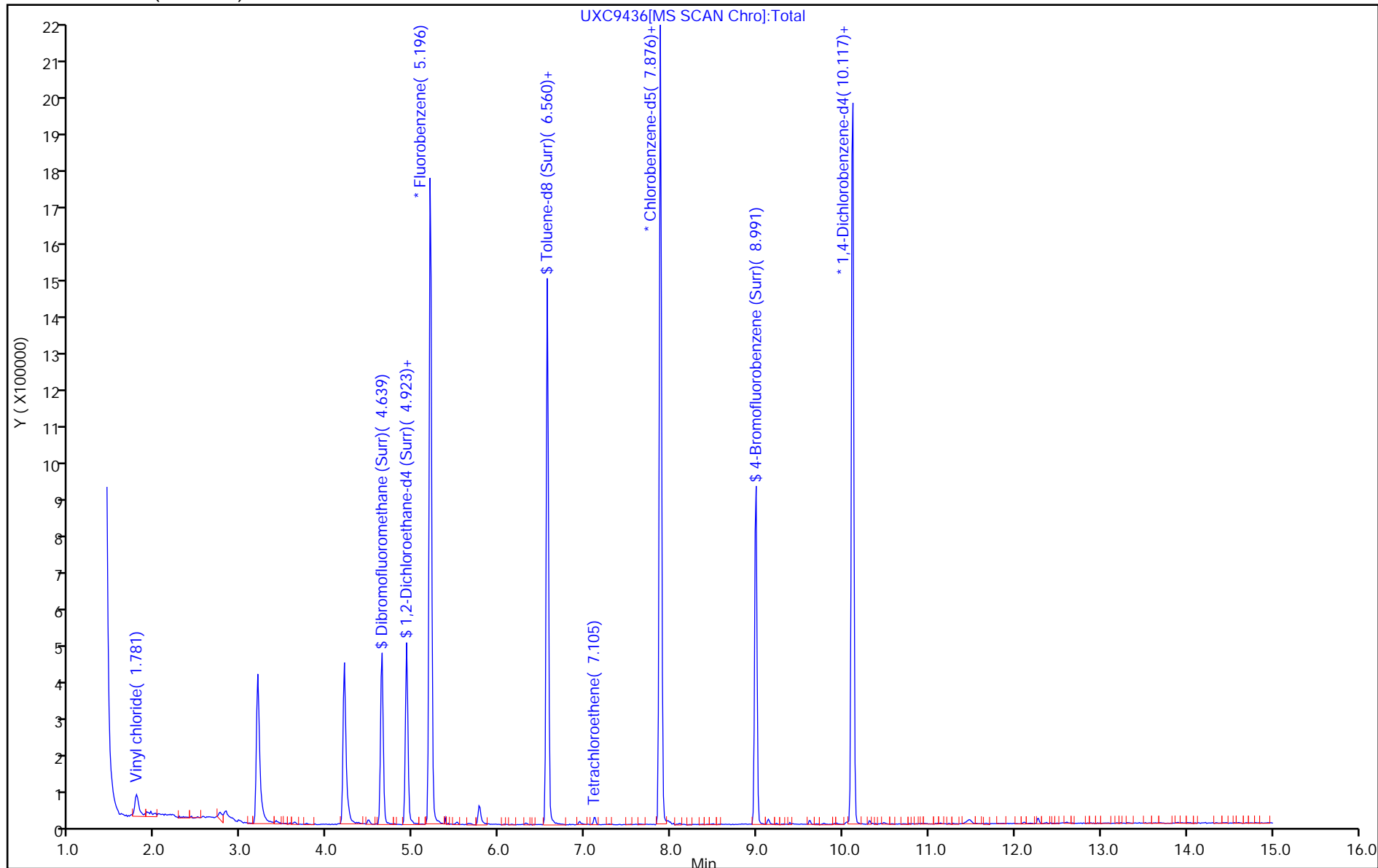
Dil. Factor: 1.0000

ALS Bottle#: 16

Method: 8260_15

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Recovery Report

Data File: \\chromna\Canton\ChromData\A3UX15\20190318-85248.b\UXC9436.D
 Lims ID: 240-109102-B-1
 Client ID: MW-72_030519
 Sample Type: Client
 Inject. Date: 18-Mar-2019 16:27:30 ALS Bottle#: 16 Worklist Smp#: 41
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0085248-041
 Misc. Info.: C90318A,8260LLUX15,,43582
 Operator ID: Instrument ID: A3UX15
 Method: \\chromna\Canton\ChromData\A3UX15\20190318-85248.b\8260_15.m
 Limit Group: MSV 8260B ICAL
 Last Update: 19-Mar-2019 08:40:05 Calib Date: 21-Jan-2019 19:15:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX15\20190121-83824.b\UXC8431.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0304

First Level Reviewer: evansle

Date: 19-Mar-2019 08:30:58

Compound	Amount Added	Amount Recovered	% Rec.
\$ 4 Dibromofluoromethane (Surr)	7.97	7.23	90.75
\$ 5 1,2-Dichloroethane-d4 (Surr)	7.97	7.09	88.94
\$ 6 Toluene-d8 (Surr)	7.97	6.26	78.56
\$ 7 4-Bromofluorobenzene (Surr)	7.97	5.46	68.46

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX15\20190318-85248.b\UXC9436.D

Injection Date: 18-Mar-2019 16:27:30

Instrument ID: A3UX15

Lims ID: 240-109102-B-1

Lab Sample ID: 240-109102-1

Client ID: MW-72_030519

Operator ID:

ALS Bottle#: 16

Worklist Smp#: 41

Purge Vol: 5.000 mL

Dil. Factor: 1.0000

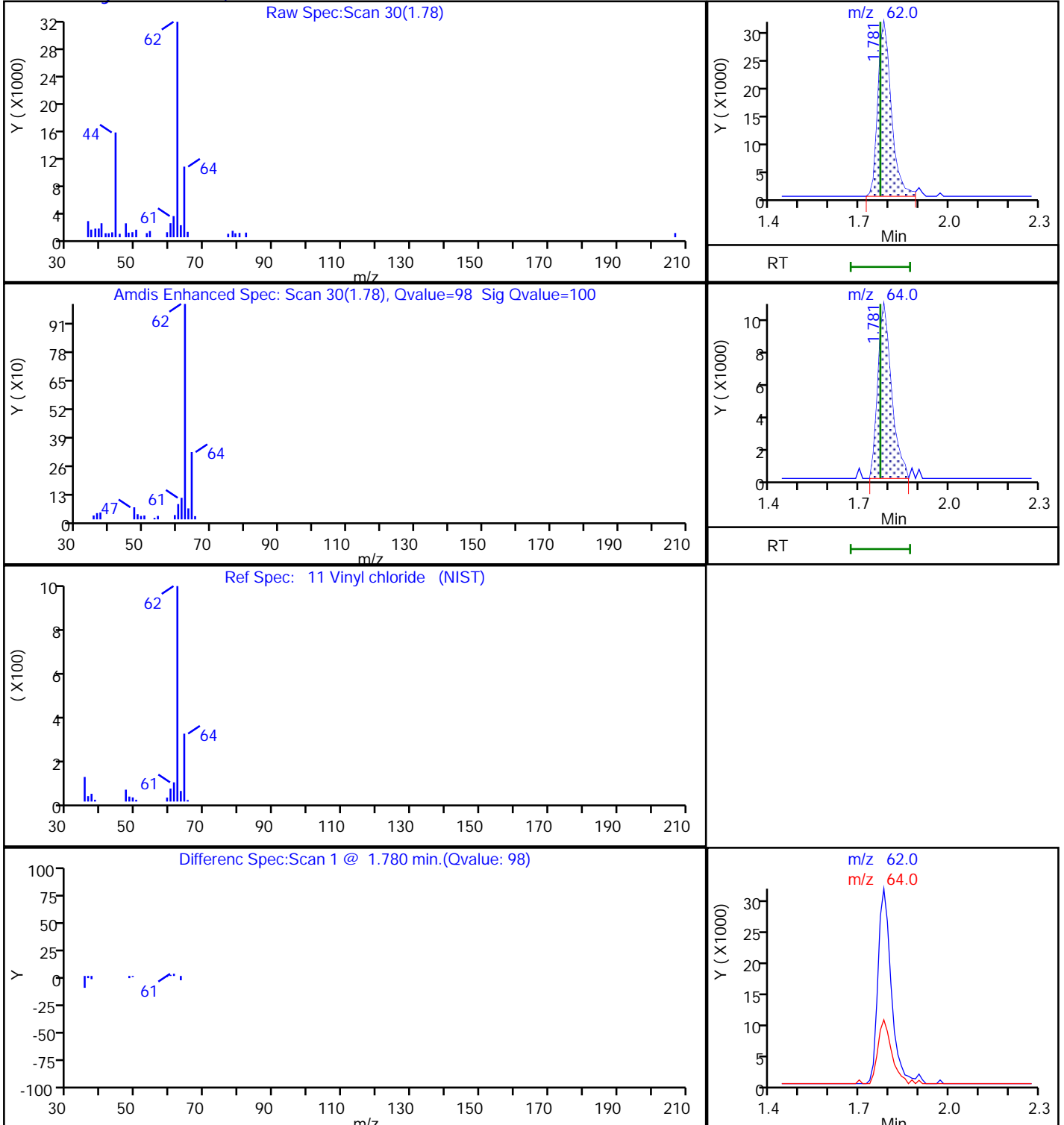
Method: 8260_15

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)

Detector: MS SCAN

11 Vinyl chloride, CAS: 75-01-4



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Client Sample ID: MW-72S_030519 Lab Sample ID: 240-109102-2
 Matrix: Water Lab File ID: UXJ8803.D
 Analysis Method: 8260B Date Collected: 03/05/2019 15:50
 Sample wt/vol: 5(mL) Date Analyzed: 03/14/2019 18:59
 Soil Aliquot Vol: _____ Dilution Factor: 1
 Soil Extract Vol.: _____ GC Column: DB-624 ID: 0.18 (mm)
 % Moisture: _____ Level: (low/med) Low
 Analysis Batch No.: 371553 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-35-4	1,1-Dichloroethene	1.0	U	1.0	0.19
156-59-2	cis-1,2-Dichloroethene	1.0	U	1.0	0.16
127-18-4	Tetrachloroethene	1.0	U	1.0	0.15
156-60-5	trans-1,2-Dichloroethene	1.0	U	1.0	0.19
79-01-6	Trichloroethene	0.20	J B	1.0	0.10
75-01-4	Vinyl chloride	0.24	J *	1.0	0.20

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	113		70-121
460-00-4	4-Bromofluorobenzene (Surr)	100		59-120
2037-26-5	Toluene-d8 (Surr)	114		70-123
1868-53-7	Dibromofluoromethane (Surr)	98		75-128

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8803.D
 Lims ID: 240-109102-E-2
 Client ID: MW-72S_030519
 Sample Type: Client
 Inject. Date: 14-Mar-2019 18:59:30 ALS Bottle#: 25 Worklist Smp#: 53
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0085154-053
 Misc. Info.: J90314A,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Method: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 15-Mar-2019 07:49:44 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0312

First Level Reviewer: evansle Date: 15-Mar-2019 07:46:27

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	4.659	4.659	0.000	98	1530908	20.0	
* 2 Chlorobenzene-d5	117	7.309	7.297	0.012	88	931259	20.0	
* 3 1,4-Dichlorobenzene-d4	152	9.522	9.522	0.000	97	346683	20.0	
\$ 4 Dibromofluoromethane (Surr	113	4.091	4.091	0.000	99	351498	19.7	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.375	4.375	0.000	98	487583	22.6	
\$ 6 Toluene-d8 (Surr)	98	6.008	6.007	0.001	94	1262173	22.8	
\$ 7 4-Bromofluorobenzene (Surr	95	8.410	8.410	0.000	98	354798	20.1	
12 Vinyl chloride	62	1.274	1.369	-0.095	40	4789	0.2363	
21 1,1-Dichloroethene	96		2.233				ND	
34 trans-1,2-Dichloroethene	96		2.872				ND	
41 cis-1,2-Dichloroethene	96		3.688				ND	
60 Trichloroethene	130	4.966	4.966	0.000	88	4599	0.2018	
76 Tetrachloroethene	164		6.552				ND	

Reagents:

vm40ml_vials_00009 Amount Added: 0.00 Units: Run Reagent
 vmDist_H2o_00139 Amount Added: 0.00 Units: Run Reagent
 vm50ss_stk_00080 Amount Added: 2.00 Units: uL Run Reagent
 VM50IS_00075 Amount Added: 2.00 Units: uL Run Reagent

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8803.D

Injection Date: 14-Mar-2019 18:59:30

Instrument ID: A3UX11

Operator ID: 43582

Lims ID: 240-109102-E-2

Lab Sample ID: 240-109102-2

Worklist Smp#: 53

Client ID: MW-72S_030519

Purge Vol: 5.000 mL

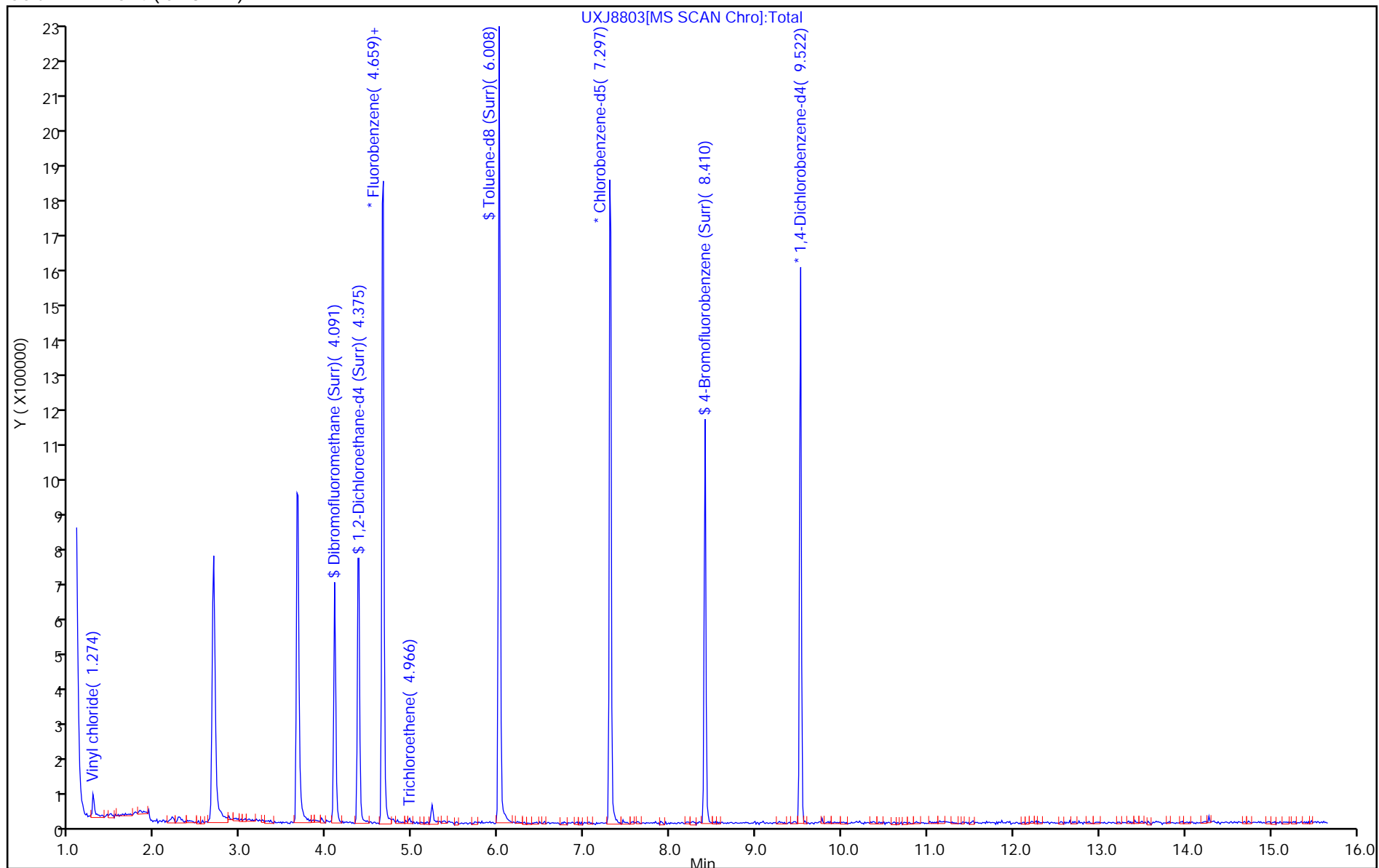
Dil. Factor: 1.0000

ALS Bottle#: 25

Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Recovery Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8803.D
 Lims ID: 240-109102-E-2
 Client ID: MW-72S_030519
 Sample Type: Client
 Inject. Date: 14-Mar-2019 18:59:30 ALS Bottle#: 25 Worklist Smp#: 53
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0085154-053
 Misc. Info.: J90314A,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Method: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 15-Mar-2019 07:49:44 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0312

First Level Reviewer: evansle Date: 15-Mar-2019 07:46:27

Compound	Amount Added	Amount Recovered	% Rec.
\$ 4 Dibromofluoromethane (Surr)	20.0	19.7	98.40
\$ 5 1,2-Dichloroethane-d4 (Surr)	20.0	22.6	112.79
\$ 6 Toluene-d8 (Surr)	20.0	22.8	114.21
\$ 7 4-Bromofluorobenzene (Surr)	20.0	20.1	100.38

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8803.D

Injection Date: 14-Mar-2019 18:59:30

Instrument ID: A3UX11

Lims ID: 240-109102-E-2

Lab Sample ID: 240-109102-2

Client ID: MW-72S_030519

Operator ID: 43582

ALS Bottle#: 25

Worklist Smp#: 53

Purge Vol: 5.000 mL

Dil. Factor: 1.0000

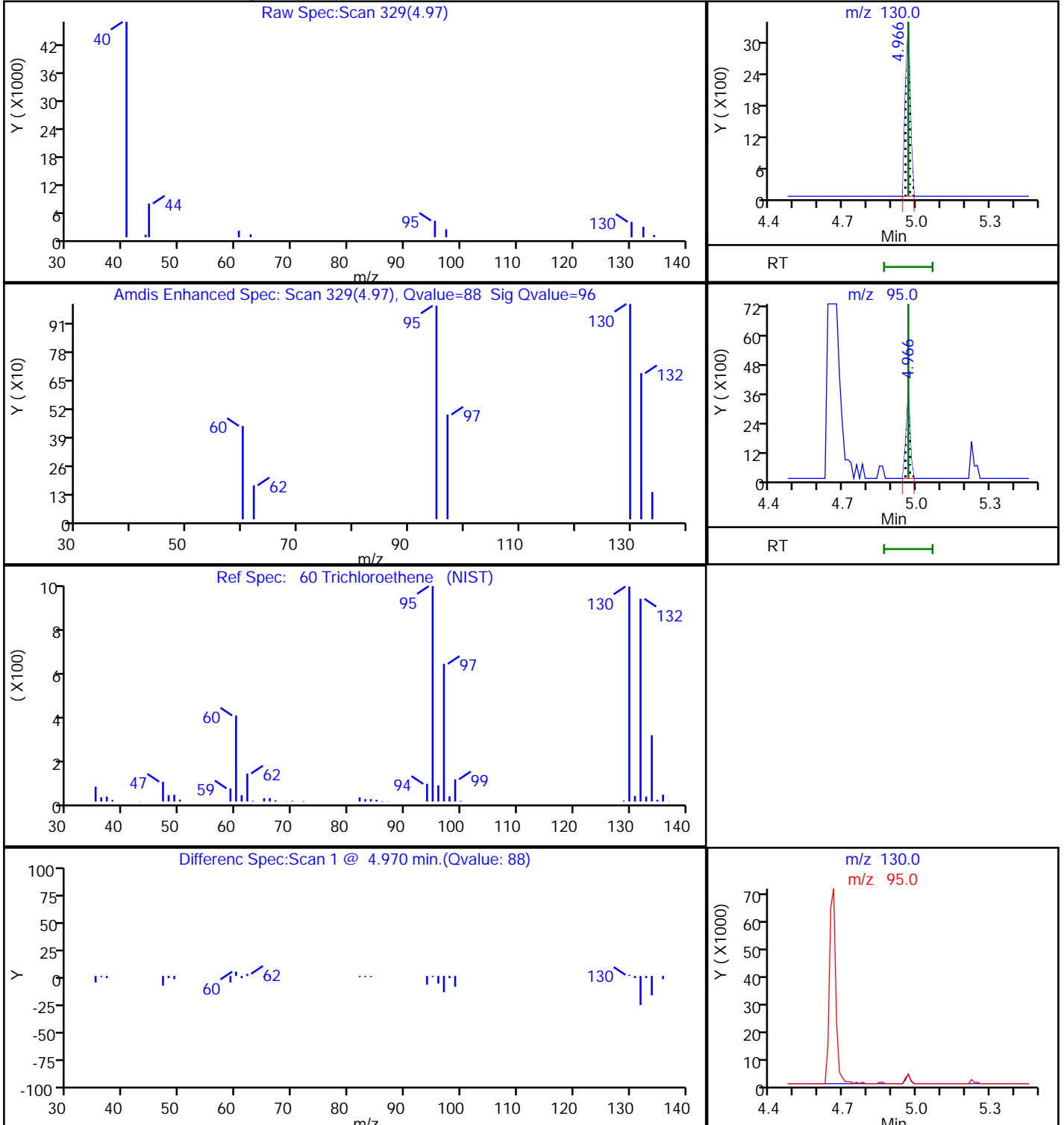
Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)

Detector: MS SCAN

60 Trichloroethene, CAS: 79-01-6



TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8803.D

Injection Date: 14-Mar-2019 18:59:30

Instrument ID: A3UX11

Lims ID: 240-109102-E-2

Lab Sample ID: 240-109102-2

Client ID: MW-72S_030519

Operator ID: 43582

ALS Bottle#: 25

Worklist Smp#: 53

Purge Vol: 5.000 mL

Dil. Factor: 1.0000

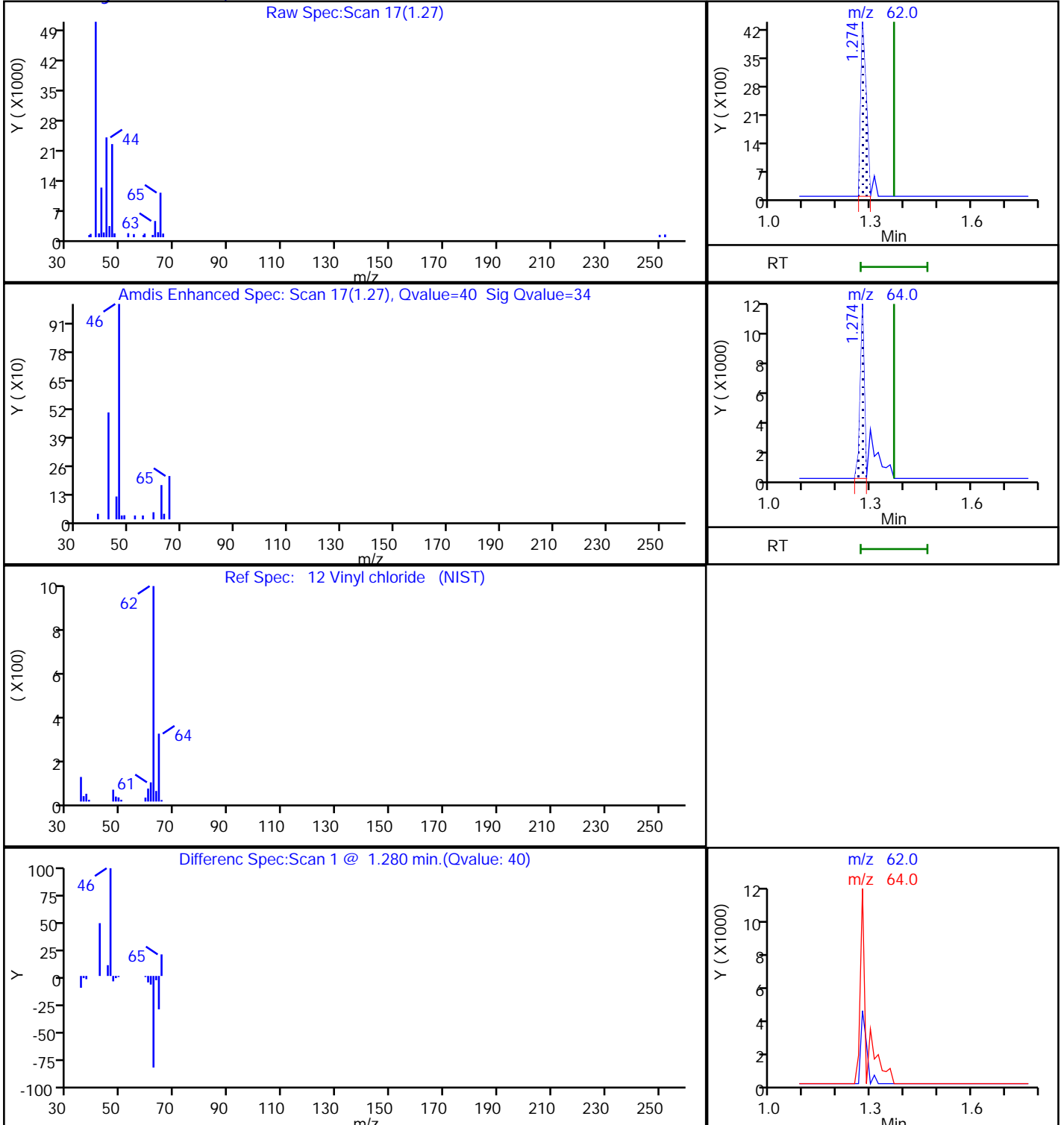
Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)

Detector: MS SCAN

12 Vinyl chloride, CAS: 75-01-4



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Client Sample ID: MW-83_030519 Lab Sample ID: 240-109102-3
 Matrix: Water Lab File ID: UXJ8804.D
 Analysis Method: 8260B Date Collected: 03/05/2019 12:13
 Sample wt/vol: 5(mL) Date Analyzed: 03/14/2019 19:22
 Soil Aliquot Vol: _____ Dilution Factor: 1
 Soil Extract Vol.: _____ GC Column: DB-624 ID: 0.18 (mm)
 % Moisture: _____ Level: (low/med) Low
 Analysis Batch No.: 371553 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-35-4	1,1-Dichloroethene	1.0	U	1.0	0.19
156-59-2	cis-1,2-Dichloroethene	1.0	U	1.0	0.16
127-18-4	Tetrachloroethene	1.0	U	1.0	0.15
156-60-5	trans-1,2-Dichloroethene	1.0	U	1.0	0.19
79-01-6	Trichloroethene	0.17	J B	1.0	0.10
75-01-4	Vinyl chloride	1.0	U *	1.0	0.20

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	109		70-121
460-00-4	4-Bromofluorobenzene (Surr)	94		59-120
2037-26-5	Toluene-d8 (Surr)	113		70-123
1868-53-7	Dibromofluoromethane (Surr)	98		75-128

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8804.D
 Lims ID: 240-109102-E-3
 Client ID: MW-83_030519
 Sample Type: Client
 Inject. Date: 14-Mar-2019 19:22:30 ALS Bottle#: 26 Worklist Smp#: 54
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0085154-054
 Misc. Info.: J90314A,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Method: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 15-Mar-2019 07:49:44 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0312

First Level Reviewer: evansle

Date: 15-Mar-2019 07:46:41

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	4.659	4.659	0.000	98	1512205	20.0	
* 2 Chlorobenzene-d5	117	7.309	7.297	0.012	88	880557	20.0	
* 3 1,4-Dichlorobenzene-d4	152	9.522	9.522	0.000	97	335209	20.0	
\$ 4 Dibromofluoromethane (Surr	113	4.091	4.091	0.000	97	344686	19.5	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.375	4.375	0.000	98	464263	21.7	
\$ 6 Toluene-d8 (Surr)	98	6.007	6.007	0.000	94	1176196	22.5	
\$ 7 4-Bromofluorobenzene (Surr	95	8.410	8.410	0.000	98	313151	18.7	
12 Vinyl chloride	62		1.369				ND	
21 1,1-Dichloroethene	96		2.233				ND	
34 trans-1,2-Dichloroethene	96		2.872				ND	
41 cis-1,2-Dichloroethene	96	3.700	3.688	0.012	78	3140	0.1460	
60 Trichloroethene	130	4.966	4.966	0.000	86	3755	0.1668	
76 Tetrachloroethene	164		6.552				ND	

Reagents:

vm40ml_vials_00009	Amount Added: 0.00	Units:	Run Reagent
vmDist_H2o_00139	Amount Added: 0.00	Units:	Run Reagent
vm50ss_stk_00080	Amount Added: 2.00	Units: uL	Run Reagent
VM50IS_00075	Amount Added: 2.00	Units: uL	Run Reagent

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8804.D

Injection Date: 14-Mar-2019 19:22:30

Instrument ID: A3UX11

Operator ID: 43582

Lims ID: 240-109102-E-3

Lab Sample ID: 240-109102-3

Worklist Smp#: 54

Client ID: MW-83_030519

Purge Vol: 5.000 mL

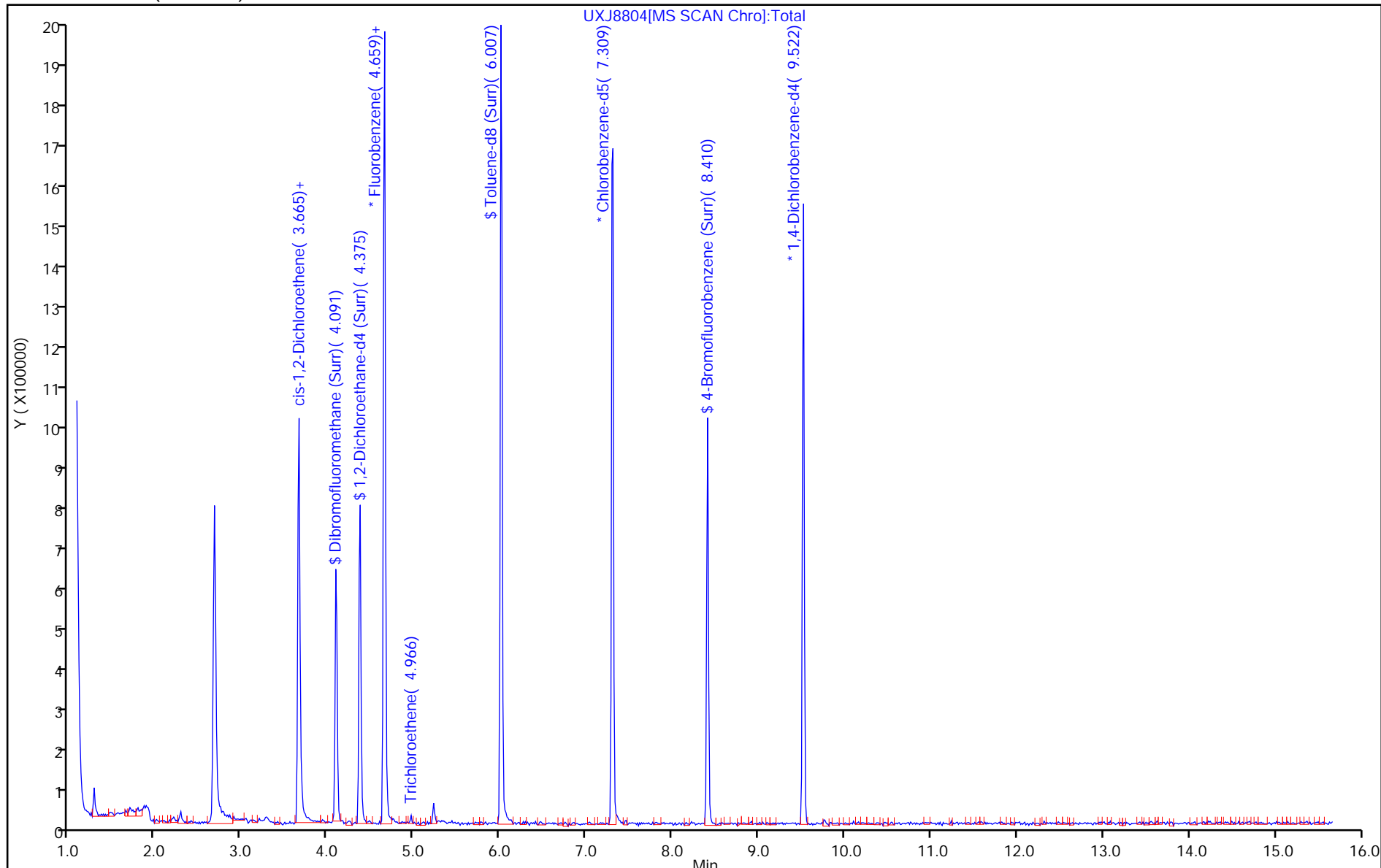
Dil. Factor: 1.0000

ALS Bottle#: 26

Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Recovery Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8804.D
 Lims ID: 240-109102-E-3
 Client ID: MW-83_030519
 Sample Type: Client
 Inject. Date: 14-Mar-2019 19:22:30 ALS Bottle#: 26 Worklist Smp#: 54
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0085154-054
 Misc. Info.: J90314A,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Method: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 15-Mar-2019 07:49:44 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0312

First Level Reviewer: evansle

Date: 15-Mar-2019 07:46:41

Compound	Amount Added	Amount Recovered	% Rec.
\$ 4 Dibromofluoromethane (Surr)	20.0	19.5	97.69
\$ 5 1,2-Dichloroethane-d4 (Surr)	20.0	21.7	108.73
\$ 6 Toluene-d8 (Surr)	20.0	22.5	112.56
\$ 7 4-Bromofluorobenzene (Surr)	20.0	18.7	93.70

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8804.D

Injection Date: 14-Mar-2019 19:22:30

Instrument ID: A3UX11

Lims ID: 240-109102-E-3

Lab Sample ID: 240-109102-3

Client ID: MW-83_030519

Operator ID: 43582

ALS Bottle#: 26

Worklist Smp#: 54

Purge Vol: 5.000 mL

Dil. Factor: 1.0000

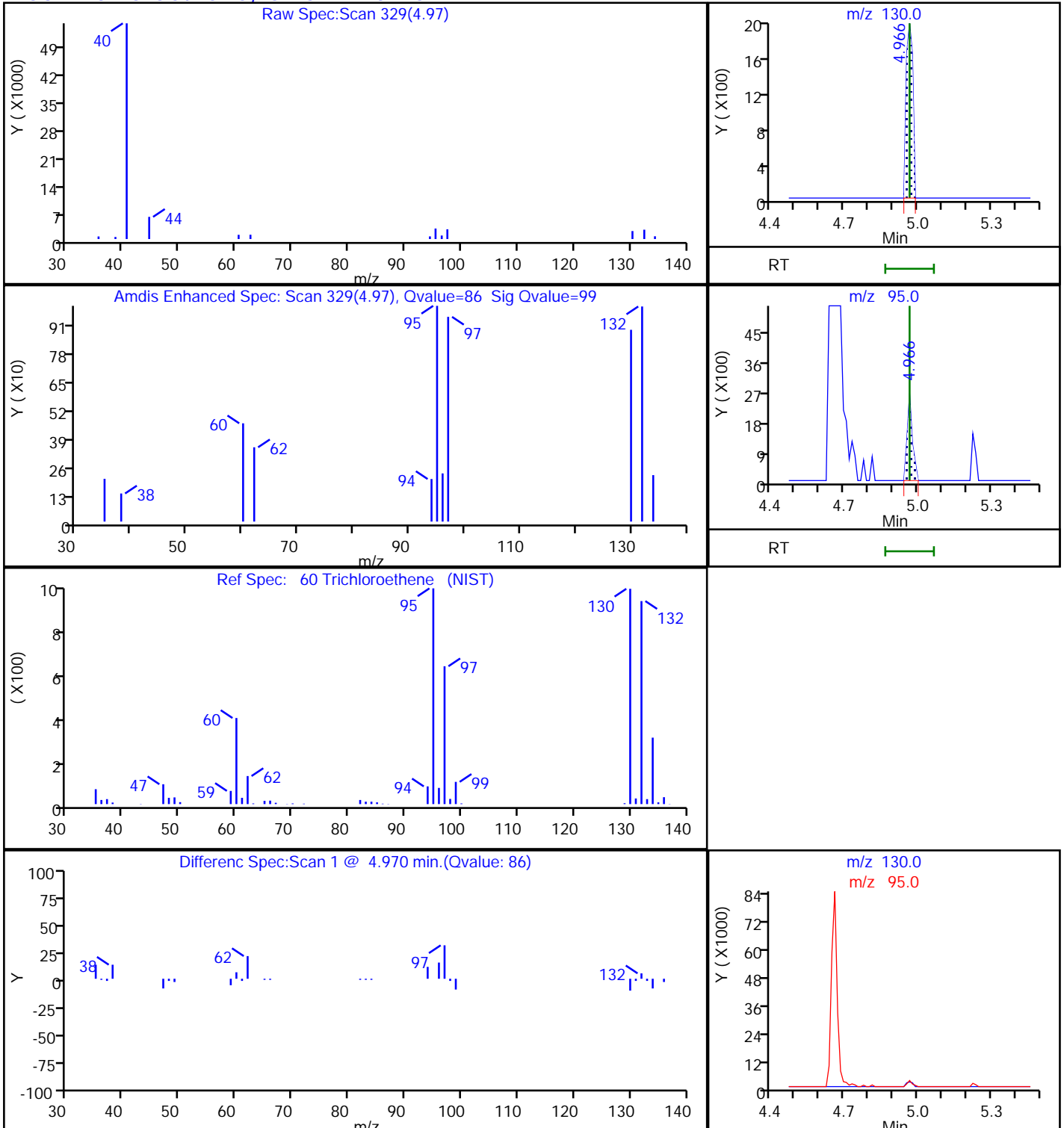
Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)

Detector: MS SCAN

60 Trichloroethene, CAS: 79-01-6



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Client Sample ID: MW-83S_030519 Lab Sample ID: 240-109102-4
 Matrix: Water Lab File ID: UXJ8805.D
 Analysis Method: 8260B Date Collected: 03/05/2019 14:07
 Sample wt/vol: 5(mL) Date Analyzed: 03/14/2019 19:44
 Soil Aliquot Vol: _____ Dilution Factor: 1
 Soil Extract Vol.: _____ GC Column: DB-624 ID: 0.18 (mm)
 % Moisture: _____ Level: (low/med) Low
 Analysis Batch No.: 371553 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-35-4	1,1-Dichloroethene	1.0	U	1.0	0.19
156-59-2	cis-1,2-Dichloroethene	1.0	U	1.0	0.16
127-18-4	Tetrachloroethene	1.0	U	1.0	0.15
156-60-5	trans-1,2-Dichloroethene	1.0	U	1.0	0.19
79-01-6	Trichloroethene	0.20	J B	1.0	0.10
75-01-4	Vinyl chloride	0.27	J *	1.0	0.20

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	105		70-121
460-00-4	4-Bromofluorobenzene (Surr)	88		59-120
2037-26-5	Toluene-d8 (Surr)	104		70-123
1868-53-7	Dibromofluoromethane (Surr)	95		75-128

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8805.D
 Lims ID: 240-109102-E-4
 Client ID: MW-83S_030519
 Sample Type: Client
 Inject. Date: 14-Mar-2019 19:44:30 ALS Bottle#: 27 Worklist Smp#: 55
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0085154-055
 Misc. Info.: J90314A,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Method: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 15-Mar-2019 07:50:44 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0312

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	4.658	4.659	-0.001	98	1482743	20.0	
* 2 Chlorobenzene-d5	117	7.309	7.297	0.012	88	914019	20.0	
* 3 1,4-Dichlorobenzene-d4	152	9.522	9.522	0.000	96	334563	20.0	
\$ 4 Dibromofluoromethane (Surr	113	4.090	4.091	-0.001	96	327103	18.9	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.374	4.375	-0.001	97	439410	21.0	
\$ 6 Toluene-d8 (Surr)	98	6.007	6.007	0.000	95	1130212	20.8	
\$ 7 4-Bromofluorobenzene (Surr	95	8.410	8.410	0.000	98	305470	17.6	
12 Vinyl chloride	62	1.286	1.369	-0.083	44	5383	0.2742	
21 1,1-Dichloroethene	96		2.233				ND	
34 trans-1,2-Dichloroethene	96		2.872				ND	
41 cis-1,2-Dichloroethene	96	3.700	3.688	0.012	81	2486	0.1179	
60 Trichloroethene	130	4.966	4.966	0.000	79	4386	0.1987	
76 Tetrachloroethene	164		6.552				ND	

Reagents:

vm40ml_vials_00009 Amount Added: 0.00 Units: Run Reagent
 vmDist_H2o_00139 Amount Added: 0.00 Units: Run Reagent
 vm50ss_stk_00080 Amount Added: 2.00 Units: uL Run Reagent
 VM50IS_00075 Amount Added: 2.00 Units: uL Run Reagent

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8805.D

Injection Date: 14-Mar-2019 19:44:30

Instrument ID: A3UX11

Operator ID: 43582

Lims ID: 240-109102-E-4

Lab Sample ID: 240-109102-4

Worklist Smp#: 55

Client ID: MW-83S_030519

Purge Vol: 5.000 mL

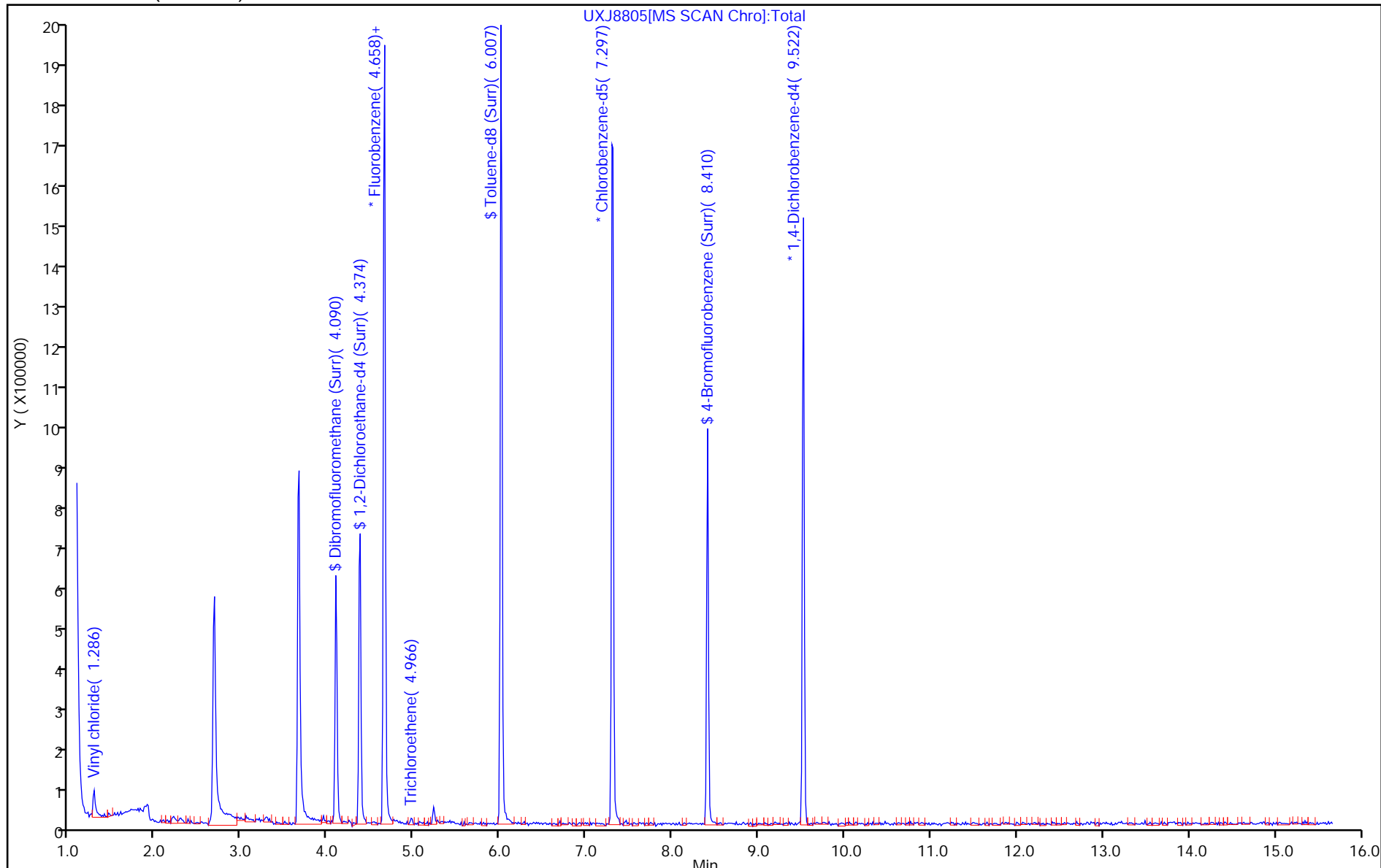
Dil. Factor: 1.0000

ALS Bottle#: 27

Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Recovery Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8805.D
 Lims ID: 240-109102-E-4
 Client ID: MW-83S_030519
 Sample Type: Client
 Inject. Date: 14-Mar-2019 19:44:30 ALS Bottle#: 27 Worklist Smp#: 55
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0085154-055
 Misc. Info.: J90314A,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Method: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 15-Mar-2019 07:50:44 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: RT Order ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0312

Compound	Amount Added	Amount Recovered	% Rec.
\$ 4 Dibromofluoromethane (Surr)	20.0	18.9	94.55
\$ 5 1,2-Dichloroethane-d4 (Surr)	20.0	21.0	104.95
\$ 6 Toluene-d8 (Surr)	20.0	20.8	104.20
\$ 7 4-Bromofluorobenzene (Surr)	20.0	17.6	88.05

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8805.D

Injection Date: 14-Mar-2019 19:44:30

Instrument ID: A3UX11

Lims ID: 240-109102-E-4

Lab Sample ID: 240-109102-4

Client ID: MW-83S_030519

Operator ID: 43582

ALS Bottle#: 27

Worklist Smp#: 55

Purge Vol: 5.000 mL

Dil. Factor: 1.0000

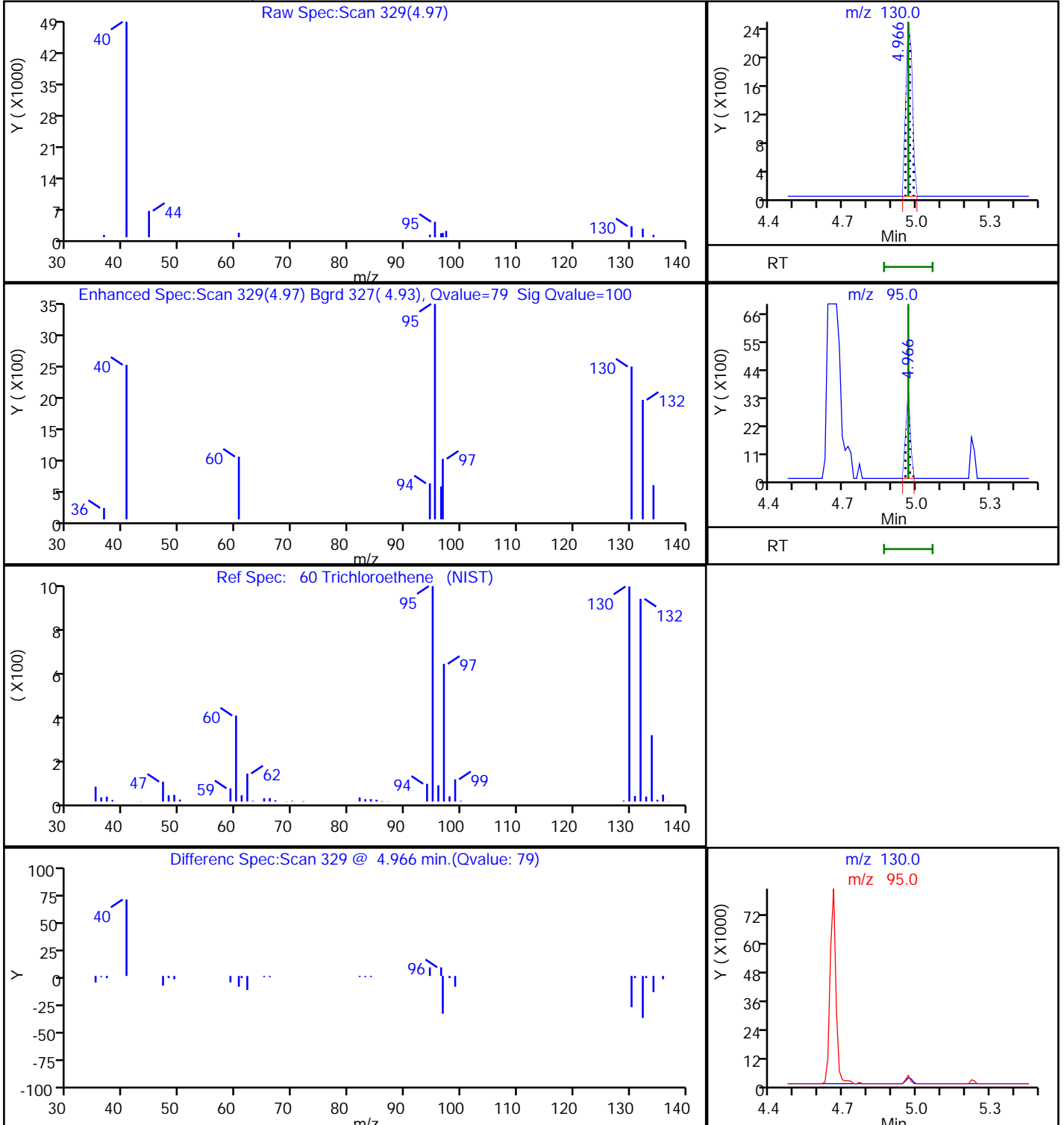
Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)

Detector: MS SCAN

60 Trichloroethene, CAS: 79-01-6



TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8805.D

Injection Date: 14-Mar-2019 19:44:30

Instrument ID: A3UX11

Lims ID: 240-109102-E-4

Lab Sample ID: 240-109102-4

Client ID: MW-83S_030519

Operator ID: 43582

ALS Bottle#: 27

Worklist Smp#: 55

Purge Vol: 5.000 mL

Dil. Factor: 1.0000

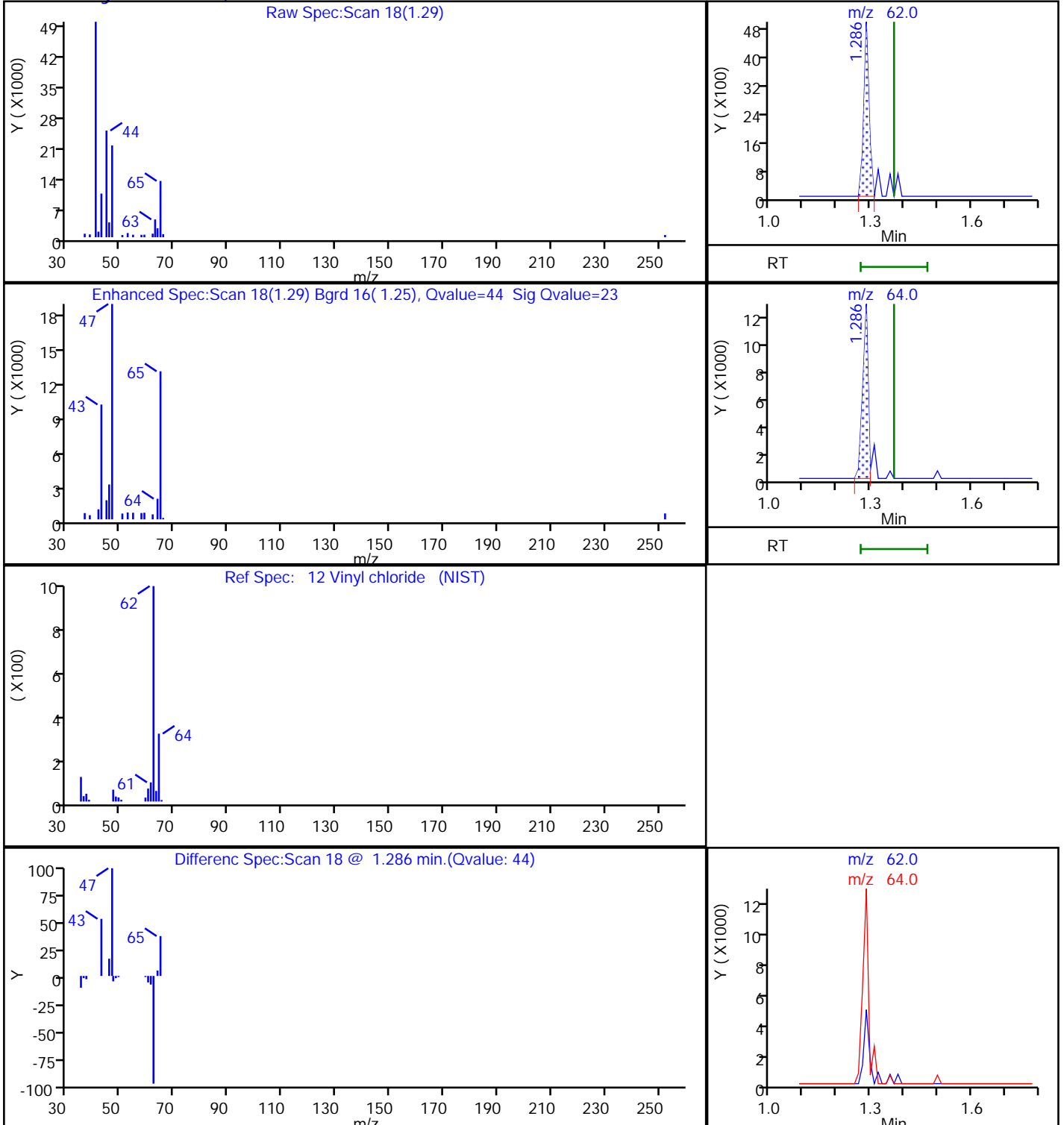
Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)

Detector: MS SCAN

12 Vinyl chloride, CAS: 75-01-4



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Client Sample ID: MW-82SR_030519 Lab Sample ID: 240-109102-5
 Matrix: Water Lab File ID: UXJ8806.D
 Analysis Method: 8260B Date Collected: 03/05/2019 17:16
 Sample wt/vol: 5(mL) Date Analyzed: 03/14/2019 20:06
 Soil Aliquot Vol: _____ Dilution Factor: 1
 Soil Extract Vol.: _____ GC Column: DB-624 ID: 0.18 (mm)
 % Moisture: _____ Level: (low/med) Low
 Analysis Batch No.: 371553 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-35-4	1,1-Dichloroethene	1.0	U	1.0	0.19
156-59-2	cis-1,2-Dichloroethene	1.0	U	1.0	0.16
127-18-4	Tetrachloroethene	1.0	U	1.0	0.15
156-60-5	trans-1,2-Dichloroethene	1.0	U	1.0	0.19
79-01-6	Trichloroethene	0.16	J B	1.0	0.10
75-01-4	Vinyl chloride	0.27	J *	1.0	0.20

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	110		70-121
460-00-4	4-Bromofluorobenzene (Surr)	90		59-120
2037-26-5	Toluene-d8 (Surr)	109		70-123
1868-53-7	Dibromofluoromethane (Surr)	96		75-128

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8806.D
 Lims ID: 240-109102-E-5
 Client ID: MW-82SR_030519
 Sample Type: Client
 Inject. Date: 14-Mar-2019 20:06:30 ALS Bottle#: 28 Worklist Smp#: 56
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0085154-056
 Misc. Info.: J90314A,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Method: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 15-Mar-2019 07:49:44 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0312

First Level Reviewer: evansle

Date: 15-Mar-2019 07:50:06

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	4.658	4.659	-0.001	98	1506099	20.0	
* 2 Chlorobenzene-d5	117	7.309	7.297	0.012	89	886816	20.0	
* 3 1,4-Dichlorobenzene-d4	152	9.522	9.522	0.000	96	324903	20.0	
\$ 4 Dibromofluoromethane (Surr	113	4.090	4.091	-0.001	98	338483	19.3	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.374	4.375	-0.001	98	469621	22.1	
\$ 6 Toluene-d8 (Surr)	98	6.007	6.007	0.000	94	1151201	21.9	
\$ 7 4-Bromofluorobenzene (Surr	95	8.410	8.410	0.000	98	301927	17.9	
12 Vinyl chloride	62	1.286	1.369	-0.083	44	5456	0.2736	
21 1,1-Dichloroethene	96		2.233				ND	
34 trans-1,2-Dichloroethene	96		2.872				ND	
41 cis-1,2-Dichloroethene	96		3.688				ND	
60 Trichloroethene	130	4.966	4.966	0.000	84	3614	0.1612	
76 Tetrachloroethene	164		6.552				ND	

Reagents:

vm40ml_vials_00009	Amount Added: 0.00	Units:	Run Reagent
vmDist_H2o_00139	Amount Added: 0.00	Units:	Run Reagent
vm50ss_stk_00080	Amount Added: 2.00	Units: uL	Run Reagent
VM50IS_00075	Amount Added: 2.00	Units: uL	Run Reagent

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8806.D

Injection Date: 14-Mar-2019 20:06:30

Instrument ID: A3UX11

Operator ID: 43582

Lims ID: 240-109102-E-5

Lab Sample ID: 240-109102-5

Worklist Smp#: 56

Client ID: MW-82SR_030519

Purge Vol: 5.000 mL

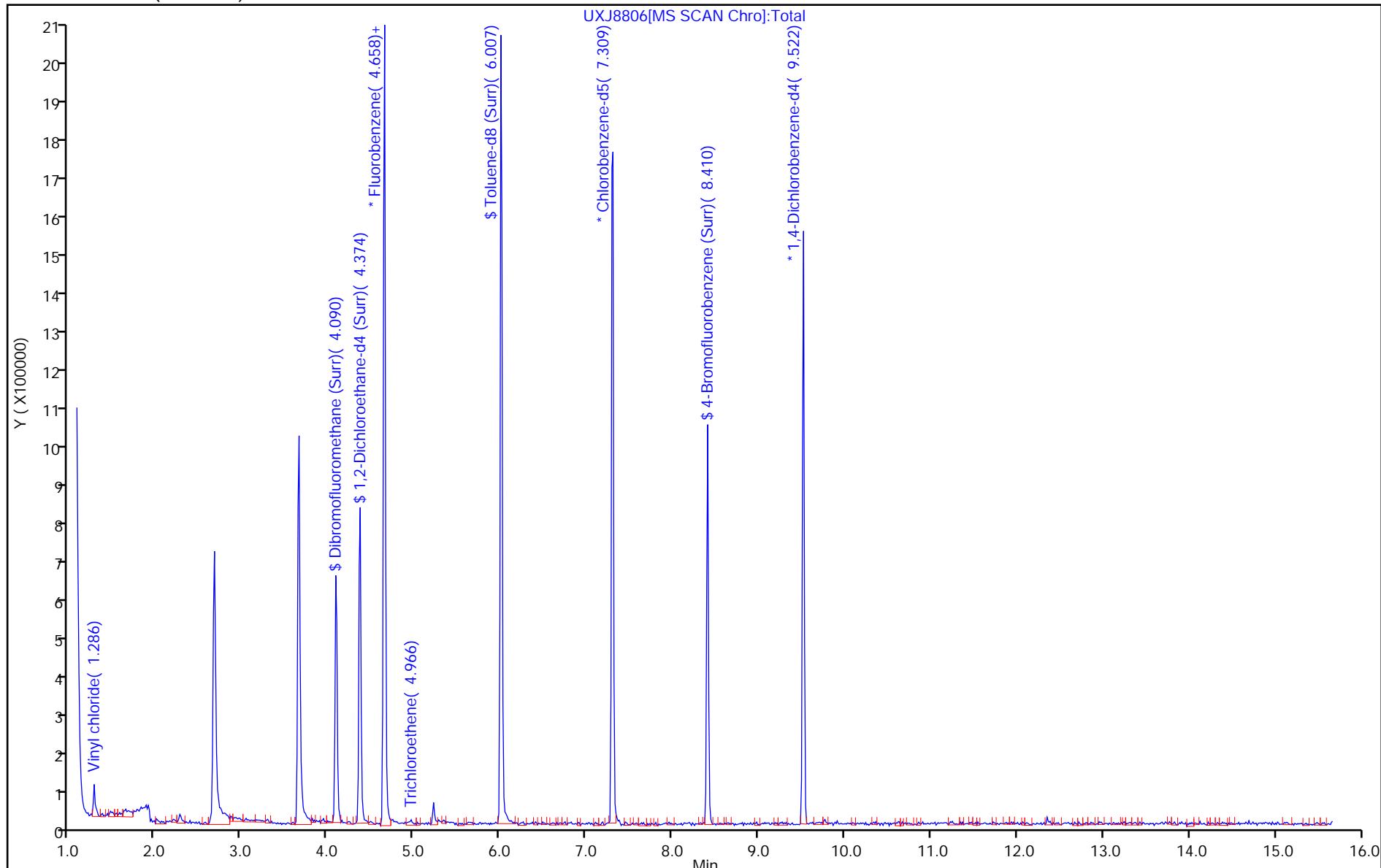
Dil. Factor: 1.0000

ALS Bottle#: 28

Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Recovery Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8806.D
 Lims ID: 240-109102-E-5
 Client ID: MW-82SR_030519
 Sample Type: Client
 Inject. Date: 14-Mar-2019 20:06:30 ALS Bottle#: 28 Worklist Smp#: 56
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0085154-056
 Misc. Info.: J90314A,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Method: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 15-Mar-2019 07:49:44 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0312

First Level Reviewer: evansle Date: 15-Mar-2019 07:50:06

Compound	Amount Added	Amount Recovered	% Rec.
\$ 4 Dibromofluoromethane (Surr)	20.0	19.3	96.32
\$ 5 1,2-Dichloroethane-d4 (Surr)	20.0	22.1	110.43
\$ 6 Toluene-d8 (Surr)	20.0	21.9	109.39
\$ 7 4-Bromofluorobenzene (Surr)	20.0	17.9	89.70

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8806.D

Injection Date: 14-Mar-2019 20:06:30

Instrument ID: A3UX11

Lims ID: 240-109102-E-5

Lab Sample ID: 240-109102-5

Client ID: MW-82SR_030519

Operator ID: 43582

ALS Bottle#: 28 Worklist Smp#: 56

Purge Vol: 5.000 mL

Dil. Factor: 1.0000

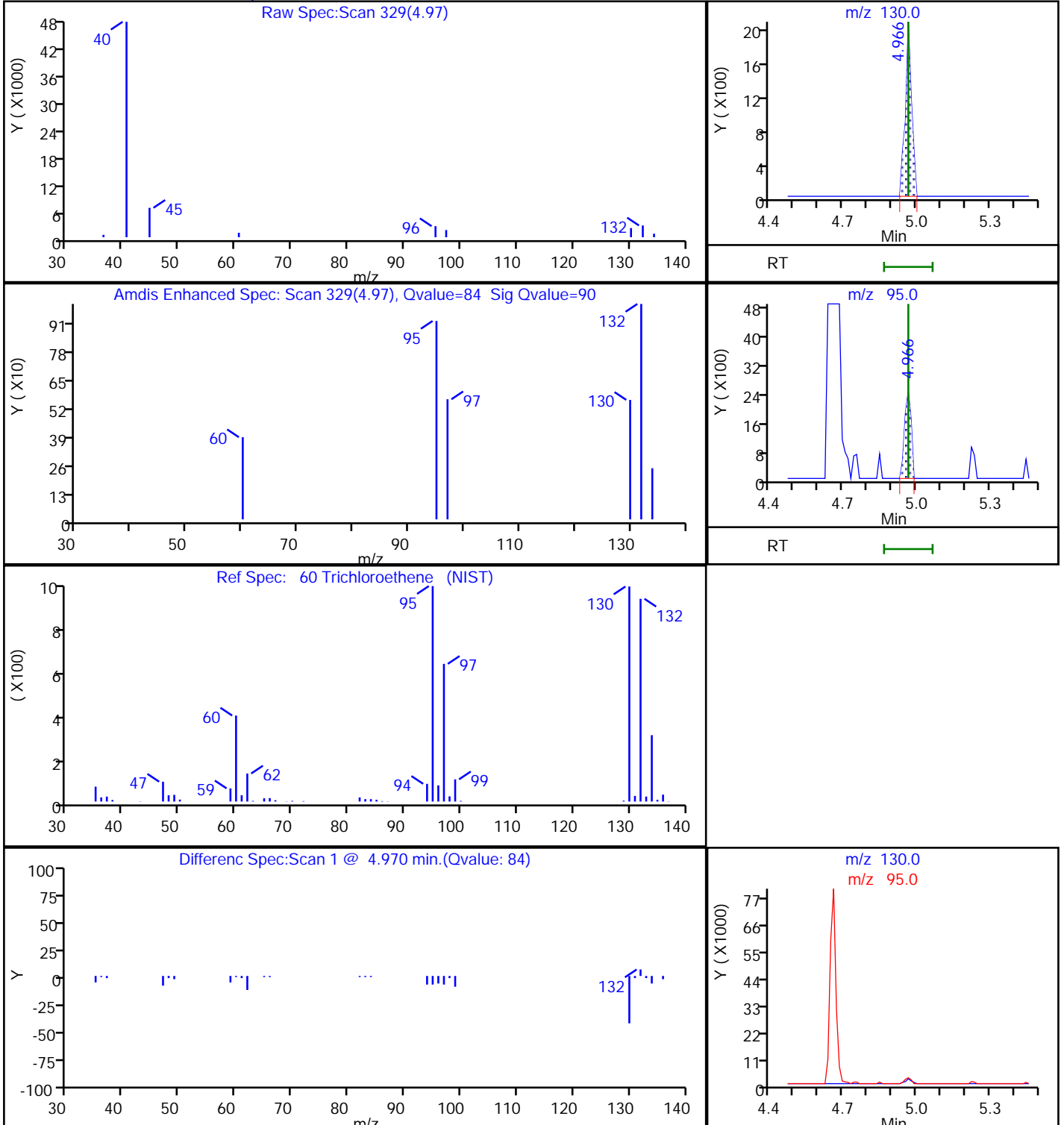
Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)

Detector: MS SCAN

60 Trichloroethene, CAS: 79-01-6



TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8806.D

Injection Date: 14-Mar-2019 20:06:30

Instrument ID: A3UX11

Lims ID: 240-109102-E-5

Lab Sample ID: 240-109102-5

Client ID: MW-82SR_030519

Operator ID: 43582

ALS Bottle#: 28

Worklist Smp#: 56

Purge Vol: 5.000 mL

Dil. Factor: 1.0000

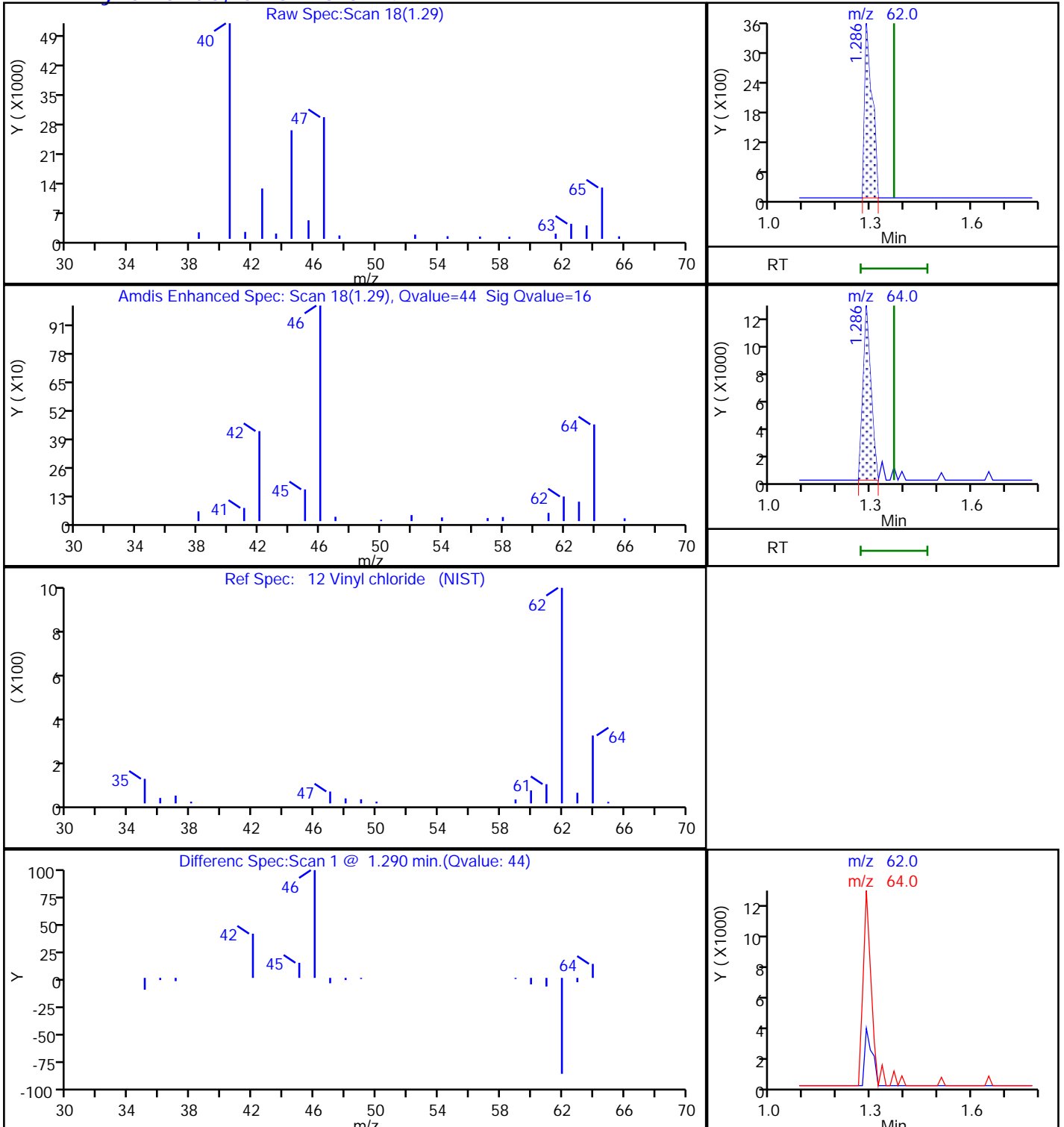
Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)

Detector: MS SCAN

12 Vinyl chloride, CAS: 75-01-4



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton Job No.: 240-109102-1
 SDG No.: _____
 Client Sample ID: TRIP BLANK Lab Sample ID: 240-109102-6
 Matrix: Water Lab File ID: UXJ8807.D
 Analysis Method: 8260B Date Collected: 03/05/2019 00:00
 Sample wt/vol: 5 (mL) Date Analyzed: 03/14/2019 20:28
 Soil Aliquot Vol: _____ Dilution Factor: 1
 Soil Extract Vol.: _____ GC Column: DB-624 ID: 0.18 (mm)
 % Moisture: _____ Level: (low/med) Low
 Analysis Batch No.: 371553 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
75-35-4	1,1-Dichloroethene	1.0	U	1.0	0.19
156-59-2	cis-1,2-Dichloroethene	1.0	U	1.0	0.16
127-18-4	Tetrachloroethene	1.0	U	1.0	0.15
156-60-5	trans-1,2-Dichloroethene	1.0	U	1.0	0.19
79-01-6	Trichloroethene	0.16	J B	1.0	0.10
75-01-4	Vinyl chloride	1.0	U *	1.0	0.20

CAS NO.	SURROGATE	%REC	Q	LIMITS
17060-07-0	1,2-Dichloroethane-d4 (Surr)	113		70-121
460-00-4	4-Bromofluorobenzene (Surr)	91		59-120
2037-26-5	Toluene-d8 (Surr)	110		70-123
1868-53-7	Dibromofluoromethane (Surr)	97		75-128

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8807.D
 Lims ID: 240-109102-A-6
 Client ID: TRIP BLANK
 Sample Type: Client
 Inject. Date: 14-Mar-2019 20:28:30 ALS Bottle#: 29 Worklist Smp#: 57
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0085154-057
 Misc. Info.: J90314A,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Method: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 15-Mar-2019 07:49:44 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0312

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/l	Flags
* 1 Fluorobenzene	96	4.659	4.659	0.000	98	1477808	20.0	
* 2 Chlorobenzene-d5	117	7.309	7.297	0.012	88	923845	20.0	
* 3 1,4-Dichlorobenzene-d4	152	9.522	9.522	0.000	97	340630	20.0	
\$ 4 Dibromofluoromethane (Surr	113	4.091	4.091	0.000	98	332842	19.3	
\$ 5 1,2-Dichloroethane-d4 (Sur	65	4.375	4.375	0.000	98	472086	22.6	
\$ 6 Toluene-d8 (Surr)	98	6.008	6.007	0.001	95	1210872	22.1	
\$ 7 4-Bromofluorobenzene (Surr	95	8.410	8.410	0.000	98	319965	18.2	
12 Vinyl chloride	62		1.369				ND	
21 1,1-Dichloroethene	96		2.233				ND	
34 trans-1,2-Dichloroethene	96		2.872				ND	
41 cis-1,2-Dichloroethene	96		3.688				ND	
60 Trichloroethene	130	4.966	4.966	0.000	83	3576	0.1625	
76 Tetrachloroethene	164		6.552				ND	

Reagents:

vm40ml_vials_00009 Amount Added: 0.00 Units: Run Reagent
 vmDist_H2o_00139 Amount Added: 0.00 Units: Run Reagent
 vm50ss_stk_00080 Amount Added: 2.00 Units: uL Run Reagent
 VM50IS_00075 Amount Added: 2.00 Units: uL Run Reagent

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8807.D

Injection Date: 14-Mar-2019 20:28:30

Instrument ID: A3UX11

Operator ID: 43582

Lims ID: 240-109102-A-6

Lab Sample ID: 240-109102-6

Worklist Smp#: 57

Client ID: TRIP BLANK

Purge Vol: 5.000 mL

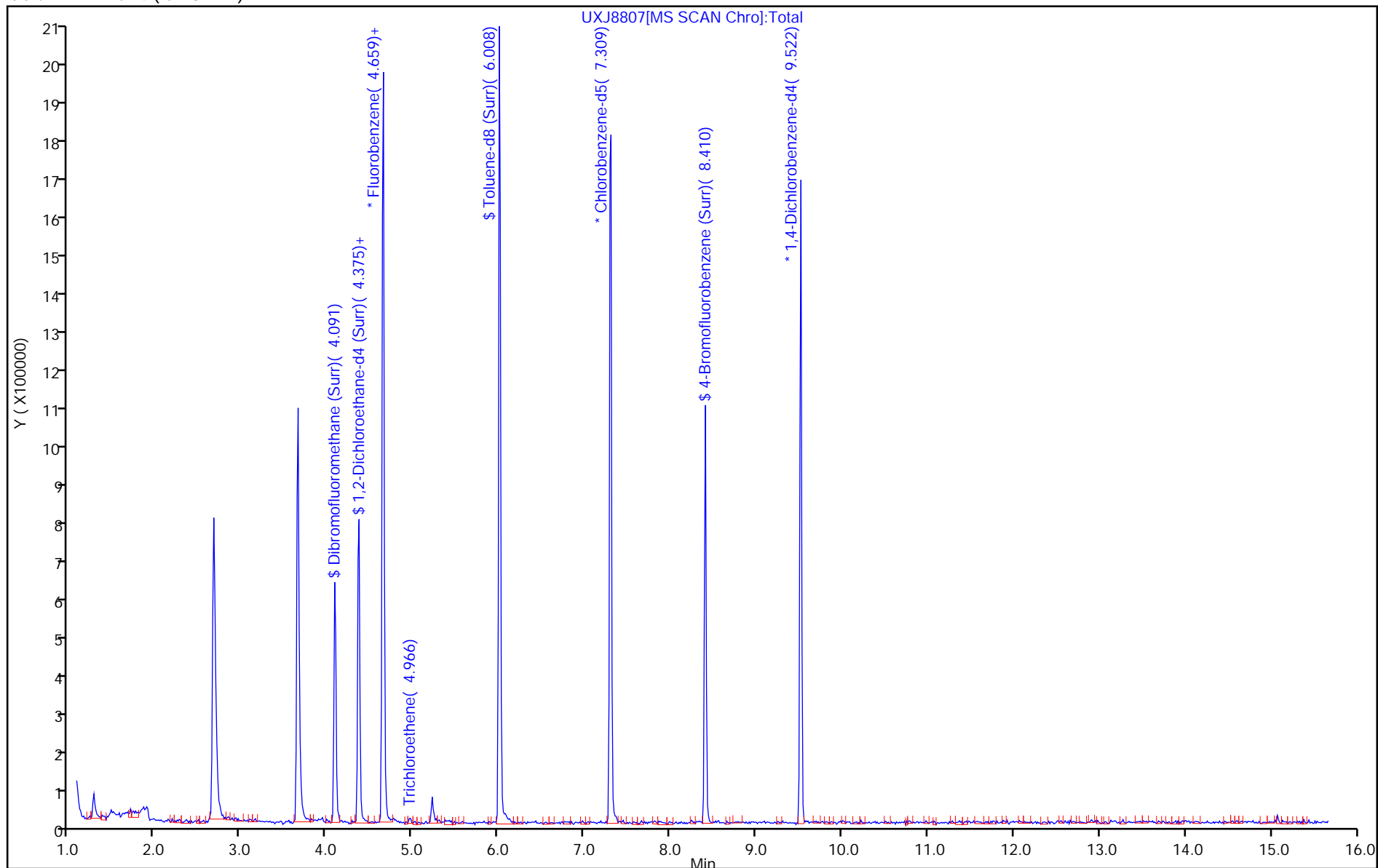
Dil. Factor: 1.0000

ALS Bottle#: 29

Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)



TestAmerica Canton
Recovery Report

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8807.D
 Lims ID: 240-109102-A-6
 Client ID: TRIP BLANK
 Sample Type: Client
 Inject. Date: 14-Mar-2019 20:28:30 ALS Bottle#: 29 Worklist Smp#: 57
 Purge Vol: 5.000 mL Dil. Factor: 1.0000
 Sample Info: 240-0085154-057
 Misc. Info.: J90314A,8260LLUX11,,43582
 Operator ID: 43582 Instrument ID: A3UX11
 Method: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\8260_11.m
 Limit Group: MSV 8260B ICAL
 Last Update: 15-Mar-2019 07:49:44 Calib Date: 07-Feb-2019 20:09:30
 Integrator: RTE ID Type: Deconvolution ID
 Quant Method: Internal Standard Quant By: Initial Calibration
 Last ICal File: \\chromna\Canton\ChromData\A3UX11\20190207-84305.b\UXJ8040.D
 Column 1 : DB-624 (0.18 mm) Det: MS SCAN
 Process Host: CTX0312

Compound	Amount Added	Amount Recovered	% Rec.
\$ 4 Dibromofluoromethane (Surr)	20.0	19.3	96.53
\$ 5 1,2-Dichloroethane-d4 (Surr)	20.0	22.6	113.13
\$ 6 Toluene-d8 (Surr)	20.0	22.1	110.45
\$ 7 4-Bromofluorobenzene (Surr)	20.0	18.2	91.25

TestAmerica Canton

Data File: \\chromna\Canton\ChromData\A3UX11\20190314-85154.b\UXJ8807.D

Injection Date: 14-Mar-2019 20:28:30

Instrument ID: A3UX11

Lims ID: 240-109102-A-6

Lab Sample ID: 240-109102-6

Client ID: TRIP BLANK

Operator ID: 43582

ALS Bottle#: 29 Worklist Smp#: 57

Purge Vol: 5.000 mL

Dil. Factor: 1.0000

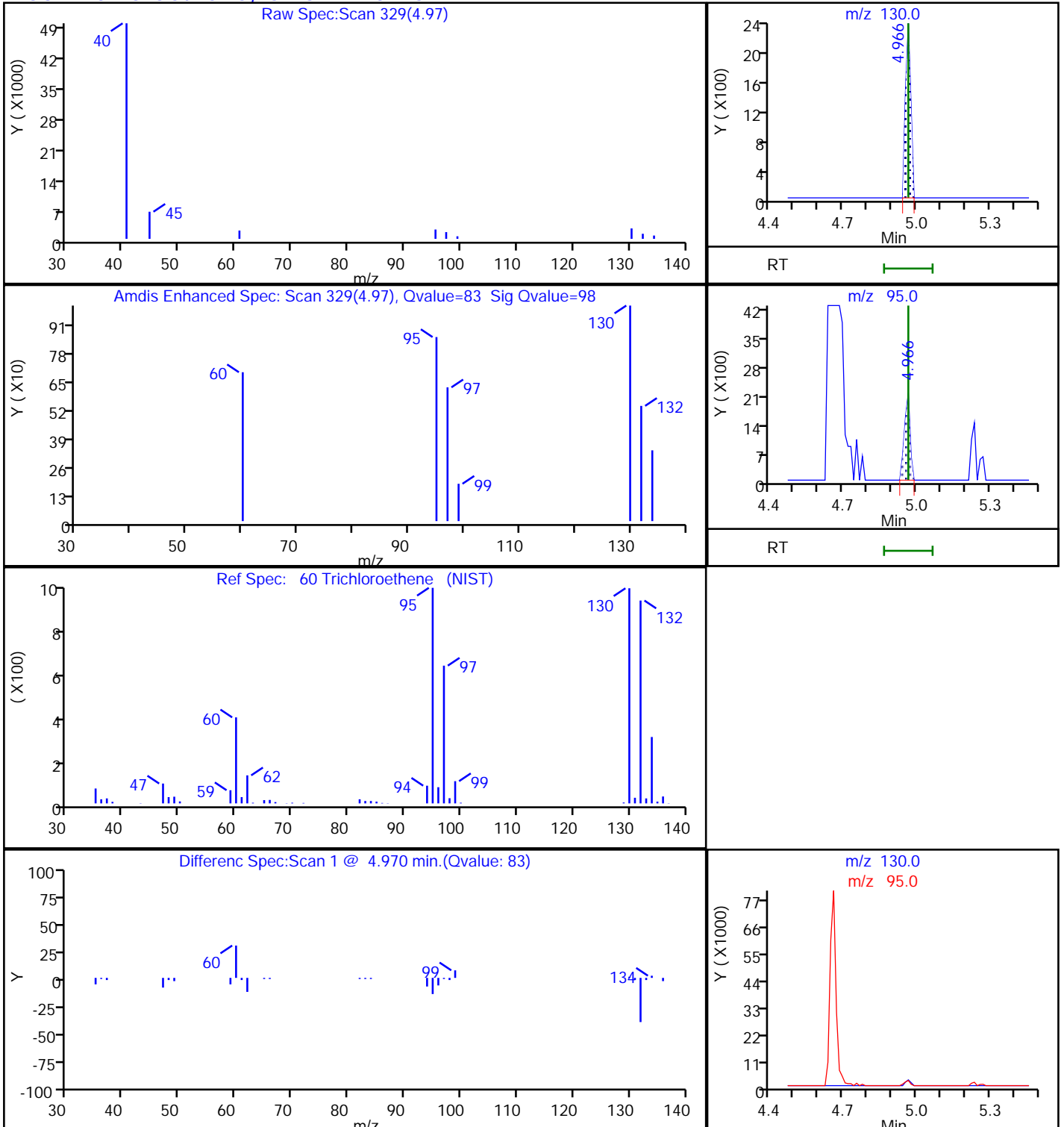
Method: 8260_11

Limit Group: MSV 8260B ICAL

Column: DB-624 (0.18 mm)

Detector: MS SCAN

60 Trichloroethene, CAS: 79-01-6



Surrogate Summary

Client: ARCADIS U.S., Inc.
Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

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

Matrix: Water

Prep Type: Total/NA

Lab Sample ID	Client Sample ID	Percent Surrogate Recovery (Acceptance Limits)			
		DCA (70-121)	BFB (59-120)	TOL (70-123)	DBFM (75-128)
240-108941-H-1 MS	Matrix Spike	103	112	124 X	93
240-108941-K-1 MSD	Matrix Spike Duplicate	102	102	118	92
240-108954-D-4 MS	Matrix Spike	84	85	87	94
240-108954-D-4 MSD	Matrix Spike Duplicate	85	84	86	93
240-109102-1	MW-72_030519	89	68	79	91
240-109102-2	MW-72S_030519	113	100	114	98
240-109102-3	MW-83_030519	109	94	113	98
240-109102-4	MW-83S_030519	105	88	104	95
240-109102-5	MW-82SR_030519	110	90	109	96
240-109102-6	TRIP BLANK	113	91	110	97
LCS 240-371553/4	Lab Control Sample	101	109	121	89
LCS 240-372032/4	Lab Control Sample	90	90	92	97
MB 240-371553/6	Method Blank	106	94	112	97
MB 240-372032/6	Method Blank	89	69	80	92

Surrogate Legend

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

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

Matrix: Water

Prep Type: Total/NA

Lab Sample ID	Client Sample ID	Percent Surrogate Recovery (Acceptance Limits)
		DCA (63-125)
240-109102-1	MW-72_030519	98
240-109102-2	MW-72S_030519	99
240-109102-3	MW-83_030519	101
240-109102-4	MW-83S_030519	98
240-109102-5	MW-82SR_030519	101
240-109195-F-7 MS	Matrix Spike	104
240-109195-F-7 MSD	Matrix Spike Duplicate	102
LCS 240-371373/4	Lab Control Sample	96
MB 240-371373/5	Method Blank	102

Surrogate Legend

DCA = 1,2-Dichloroethane-d4 (Surr)

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

Matrix: Water

Prep Type: Total/NA

Lab Sample ID	Client Sample ID	Percent Surrogate Recovery (Acceptance Limits)
		DCA (10-150)
MRL 240-371373/6	Lab Control Sample	99

Surrogate Legend

DCA = 1,2-Dichloroethane-d4 (Surr)

TestAmerica Canton

QC Sample Results

Client: ARCADIS U.S., Inc.
 Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

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

Lab Sample ID: MB 240-371553/6

Matrix: Water

Analysis Batch: 371553

Client Sample ID: Method Blank

Prep Type: Total/NA

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/14/19 11:35	1
cis-1,2-Dichloroethene	0.237	J	1.0	0.16	ug/L			03/14/19 11:35	1
Tetrachloroethene	1.0	U	1.0	0.15	ug/L			03/14/19 11:35	1
trans-1,2-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/14/19 11:35	1
Trichloroethene	0.319	J	1.0	0.10	ug/L			03/14/19 11:35	1
Vinyl chloride	1.0	U	1.0	0.20	ug/L			03/14/19 11:35	1

Surrogate	MB %Recovery	MB Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	106		70 - 121		03/14/19 11:35	1
4-Bromofluorobenzene (Surr)	94		59 - 120		03/14/19 11:35	1
Toluene-d8 (Surr)	112		70 - 123		03/14/19 11:35	1
Dibromofluoromethane (Surr)	97		75 - 128		03/14/19 11:35	1

Lab Sample ID: LCS 240-371553/4

Matrix: Water

Analysis Batch: 371553

Client Sample ID: Lab Control Sample

Prep Type: Total/NA

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec. Limits
1,1-Dichloroethene	10.0	11.4		ug/L		114	65 - 139
cis-1,2-Dichloroethene	10.0	10.6		ug/L		106	76 - 128
Tetrachloroethene	10.0	8.77		ug/L		88	74 - 130
trans-1,2-Dichloroethene	10.0	10.1		ug/L		101	78 - 133
Trichloroethene	10.0	8.77		ug/L		88	76 - 125
Vinyl chloride	10.0	16.3	*	ug/L		163	58 - 143

Surrogate	LCS %Recovery	LCS Qualifier	Limits
1,2-Dichloroethane-d4 (Surr)	101		70 - 121
4-Bromofluorobenzene (Surr)	109		59 - 120
Toluene-d8 (Surr)	121		70 - 123
Dibromofluoromethane (Surr)	89		75 - 128

Lab Sample ID: MRL 240-371553/5

Matrix: Water

Analysis Batch: 371553

Client Sample ID: Lab Control Sample

Prep Type: Total/NA

Analyte	Spike Added	MRL Result	MRL Qualifier	Unit	D	%Rec	%Rec. Limits
Vinyl chloride	0.00100	0.00154	^	ng/uL		154	10 - 150

Lab Sample ID: 240-108941-H-1 MS

Matrix: Water

Analysis Batch: 371553

Client Sample ID: Matrix Spike

Prep Type: Total/NA

Analyte	Sample Result	Sample Qualifier	Spike Added	MS Result	MS Qualifier	Unit	D	%Rec	%Rec. Limits
1,1-Dichloroethene	1.0	U	10.0	10.1		ug/L		101	53 - 140
cis-1,2-Dichloroethene	0.30	J B	10.0	10.6		ug/L		103	64 - 130
Tetrachloroethene	1.0	U	10.0	7.44		ug/L		74	51 - 136
trans-1,2-Dichloroethene	1.0	U	10.0	10.1		ug/L		101	68 - 133
Trichloroethene	0.19	J B	10.0	8.00		ug/L		78	55 - 131

TestAmerica Canton

QC Sample Results

Client: ARCADIS U.S., Inc.
Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

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

Lab Sample ID: 240-108941-H-1 MS

Matrix: Water

Analysis Batch: 371553

Client Sample ID: Matrix Spike

Prep Type: Total/NA

Analyte	Sample Result	Sample Qualifier	Spike Added	MS Result	MS Qualifier	Unit	D	%Rec	%Rec. Limits
Vinyl chloride	0.71	J*	10.0	14.2		ug/L		135	43 - 154
Surrogate	MS %Recovery	MS Qualifier	Limits						
1,2-Dichloroethane-d4 (Surr)	103		70 - 121						
4-Bromofluorobenzene (Surr)	112		59 - 120						
Toluene-d8 (Surr)	124	X	70 - 123						
Dibromofluoromethane (Surr)	93		75 - 128						

Lab Sample ID: 240-108941-K-1 MSD

Matrix: Water

Analysis Batch: 371553

Client Sample ID: Matrix Spike Duplicate

Prep Type: Total/NA

Analyte	Sample Result	Sample Qualifier	Spike Added	MSD Result	MSD Qualifier	Unit	D	%Rec	%Rec. Limits	RPD	RPD Limit
1,1-Dichloroethene	1.0	U	10.0	11.5		ug/L		115	53 - 140	13	35
cis-1,2-Dichloroethene	0.30	J B	10.0	10.3		ug/L		100	64 - 130	2	21
Tetrachloroethene	1.0	U	10.0	8.04		ug/L		80	51 - 136	8	23
trans-1,2-Dichloroethene	1.0	U	10.0	10.6		ug/L		106	68 - 133	5	24
Trichloroethene	0.19	J B	10.0	8.17		ug/L		80	55 - 131	2	23
Vinyl chloride	0.71	J*	10.0	15.3		ug/L		146	43 - 154	8	29
Surrogate	MSD %Recovery	MSD Qualifier	Limits								
1,2-Dichloroethane-d4 (Surr)	102		70 - 121								
4-Bromofluorobenzene (Surr)	102		59 - 120								
Toluene-d8 (Surr)	118		70 - 123								
Dibromofluoromethane (Surr)	92		75 - 128								

Lab Sample ID: MB 240-372032/6

Matrix: Water

Analysis Batch: 372032

Client Sample ID: Method Blank

Prep Type: Total/NA

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/18/19 13:10	1
cis-1,2-Dichloroethene	1.0	U	1.0	0.16	ug/L			03/18/19 13:10	1
Tetrachloroethene	1.0	U	1.0	0.15	ug/L			03/18/19 13:10	1
trans-1,2-Dichloroethene	1.0	U	1.0	0.19	ug/L			03/18/19 13:10	1
Trichloroethene	1.0	U	1.0	0.10	ug/L			03/18/19 13:10	1
Vinyl chloride	1.0	U	1.0	0.20	ug/L			03/18/19 13:10	1
Surrogate	MB %Recovery	MB Qualifier	Limits				Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	89		70 - 121					03/18/19 13:10	1
4-Bromofluorobenzene (Surr)	69		59 - 120					03/18/19 13:10	1
Toluene-d8 (Surr)	80		70 - 123					03/18/19 13:10	1
Dibromofluoromethane (Surr)	92		75 - 128					03/18/19 13:10	1

TestAmerica Canton

QC Sample Results

Client: ARCADIS U.S., Inc.
Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

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

Lab Sample ID: LCS 240-372032/4

Matrix: Water

Analysis Batch: 372032

Client Sample ID: Lab Control Sample

Prep Type: Total/NA

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec. Limits
1,1-Dichloroethene	10.0	9.75		ug/L		97	65 - 139
cis-1,2-Dichloroethene	10.0	11.4		ug/L		114	76 - 128
Tetrachloroethene	10.0	11.6		ug/L		116	74 - 130
trans-1,2-Dichloroethene	10.0	11.9		ug/L		119	78 - 133
Trichloroethene	10.0	10.6		ug/L		106	76 - 125
Vinyl chloride	10.0	9.89		ug/L		99	58 - 143

Surrogate	LCS %Recovery	LCS Qualifier	Limits
1,2-Dichloroethane-d4 (Surr)	90		70 - 121
4-Bromofluorobenzene (Surr)	90		59 - 120
Toluene-d8 (Surr)	92		70 - 123
Dibromofluoromethane (Surr)	97		75 - 128

Lab Sample ID: MRL 240-372032/5

Matrix: Water

Analysis Batch: 372032

Client Sample ID: Lab Control Sample

Prep Type: Total/NA

Analyte	Spike Added	MRL Result	MRL Qualifier	Unit	D	%Rec	%Rec. Limits
Vinyl chloride	0.00100	0.000830	J	ng/uL		83	10 - 150

Lab Sample ID: 240-108954-D-4 MS

Matrix: Water

Analysis Batch: 372032

Client Sample ID: Matrix Spike

Prep Type: Total/NA

Analyte	Sample Result	Sample Qualifier	Spike Added	MS Result	MS Qualifier	Unit	D	%Rec	%Rec. Limits
1,1-Dichloroethene	6.8	F1	10.0	9.42	F1	ug/L		26	53 - 140
cis-1,2-Dichloroethene	2.0	J	10.0	10.3		ug/L		83	64 - 130
Tetrachloroethene	5.0	U	10.0	10.5		ug/L		105	51 - 136
trans-1,2-Dichloroethene	5.0	U	10.0	9.80		ug/L		98	68 - 133
Trichloroethene	0.87	J	10.0	8.74		ug/L		79	55 - 131
Vinyl chloride	19	F1	10.0	11.0	F1	ug/L		-85	43 - 154

Surrogate	MS %Recovery	MS Qualifier	Limits
1,2-Dichloroethane-d4 (Surr)	84		70 - 121
4-Bromofluorobenzene (Surr)	85		59 - 120
Toluene-d8 (Surr)	87		70 - 123
Dibromofluoromethane (Surr)	94		75 - 128

Lab Sample ID: 240-108954-D-4 MSD

Matrix: Water

Analysis Batch: 372032

Client Sample ID: Matrix Spike Duplicate

Prep Type: Total/NA

Analyte	Sample Result	Sample Qualifier	Spike Added	MSD Result	MSD Qualifier	Unit	D	%Rec	%Rec. Limits	RPD	RPD Limit
1,1-Dichloroethene	6.8	F1	10.0	10.0	F1	ug/L		32	53 - 140	6	35
cis-1,2-Dichloroethene	2.0	J	10.0	11.1		ug/L		91	64 - 130	7	21
Tetrachloroethene	5.0	U	10.0	10.8		ug/L		108	51 - 136	3	23
trans-1,2-Dichloroethene	5.0	U	10.0	10.7		ug/L		107	68 - 133	9	24
Trichloroethene	0.87	J	10.0	9.67		ug/L		88	55 - 131	10	23

TestAmerica Canton

QC Sample Results

Client: ARCADIS U.S., Inc.
 Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

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

Lab Sample ID: 240-108954-D-4 MSD
Matrix: Water
Analysis Batch: 372032

Client Sample ID: Matrix Spike Duplicate
Prep Type: Total/NA

Analyte	Sample Result	Sample Qualifier	Spike Added	MSD Result	MSD Qualifier	Unit	D	%Rec	%Rec. Limits	RPD	RPD Limit
Vinyl chloride	19	F1	10.0	12.0	F1	ug/L		-75	43 - 154	9	29
Surrogate	%Recovery	MSD Qualifier	MSD Limits								
1,2-Dichloroethane-d4 (Surr)	85		70 - 121								
4-Bromofluorobenzene (Surr)	84		59 - 120								
Toluene-d8 (Surr)	86		70 - 123								
Dibromofluoromethane (Surr)	93		75 - 128								

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

Lab Sample ID: MB 240-371373/5
Matrix: Water
Analysis Batch: 371373

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

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,4-Dioxane	2.0	U	2.0	0.86	ug/L			03/13/19 11:11	1
Surrogate	%Recovery	MB Qualifier	MB Limits						
1,2-Dichloroethane-d4 (Surr)	102		63 - 125						

Lab Sample ID: LCS 240-371373/4
Matrix: Water
Analysis Batch: 371373

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

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec. Limits
1,4-Dioxane	10.0	10.8	J	ug/L		108	59 - 131
Surrogate	%Recovery	LCS Qualifier	LCS Limits				
1,2-Dichloroethane-d4 (Surr)	96		63 - 125				

Lab Sample ID: MRL 240-371373/6
Matrix: Water
Analysis Batch: 371373

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

Analyte	Spike Added	MRL Result	MRL Qualifier	Unit	D	%Rec	%Rec. Limits
1,4-Dioxane	0.00100	0.00132	J	ng/uL		132	10 - 150
Surrogate	%Recovery	MRL Qualifier	MRL Limits				
1,2-Dichloroethane-d4 (Surr)	99		10 - 150				

Lab Sample ID: 240-109195-F-7 MS
Matrix: Water
Analysis Batch: 371373

Client Sample ID: Matrix Spike
Prep Type: Total/NA

Analyte	Sample Result	Sample Qualifier	Spike Added	MS Result	MS Qualifier	Unit	D	%Rec	%Rec. Limits
1,4-Dioxane	2.0	U F2	10.0	10.0		ug/L		100	52 - 129

TestAmerica Canton

QC Sample Results

Client: ARCADIS U.S., Inc.
 Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

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

Lab Sample ID: 240-109195-F-7 MS
 Matrix: Water
 Analysis Batch: 371373

Client Sample ID: Matrix Spike
 Prep Type: Total/NA

Surrogate	MS %Recovery	MS Qualifier	Limits
1,2-Dichloroethane-d4 (Surr)	104		63 - 125

Lab Sample ID: 240-109195-F-7 MSD
 Matrix: Water
 Analysis Batch: 371373

Client Sample ID: Matrix Spike Duplicate
 Prep Type: Total/NA

Analyte	Sample Result	Sample Qualifier	Spike Added	MSD Result	MSD Qualifier	Unit	D	%Rec	%Rec. Limits	RPD	RPD Limit
1,4-Dioxane	2.0	U F2	10.0	11.8	F2	ug/L		118	52 - 129	16	13

Surrogate	MSD %Recovery	MSD Qualifier	Limits
1,2-Dichloroethane-d4 (Surr)	102		63 - 125

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QC Association Summary

Client: ARCADIS U.S., Inc.
 Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

GC/MS VOA

Analysis Batch: 371373

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
240-109102-1	MW-72_030519	Total/NA	Water	8260B SIM	
240-109102-2	MW-72S_030519	Total/NA	Water	8260B SIM	
240-109102-3	MW-83_030519	Total/NA	Water	8260B SIM	
240-109102-4	MW-83S_030519	Total/NA	Water	8260B SIM	
240-109102-5	MW-82SR_030519	Total/NA	Water	8260B SIM	
MB 240-371373/5	Method Blank	Total/NA	Water	8260B SIM	
LCS 240-371373/4	Lab Control Sample	Total/NA	Water	8260B SIM	
MRL 240-371373/6	Lab Control Sample	Total/NA	Water	8260B SIM	
240-109195-F-7 MS	Matrix Spike	Total/NA	Water	8260B SIM	
240-109195-F-7 MSD	Matrix Spike Duplicate	Total/NA	Water	8260B SIM	

Analysis Batch: 371553

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
240-109102-2	MW-72S_030519	Total/NA	Water	8260B	
240-109102-3	MW-83_030519	Total/NA	Water	8260B	
240-109102-4	MW-83S_030519	Total/NA	Water	8260B	
240-109102-5	MW-82SR_030519	Total/NA	Water	8260B	
240-109102-6	TRIP BLANK	Total/NA	Water	8260B	
MB 240-371553/6	Method Blank	Total/NA	Water	8260B	
LCS 240-371553/4	Lab Control Sample	Total/NA	Water	8260B	
MRL 240-371553/5	Lab Control Sample	Total/NA	Water	8260B	
240-108941-H-1 MS	Matrix Spike	Total/NA	Water	8260B	
240-108941-K-1 MSD	Matrix Spike Duplicate	Total/NA	Water	8260B	

Analysis Batch: 372032

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
240-109102-1	MW-72_030519	Total/NA	Water	8260B	
MB 240-372032/6	Method Blank	Total/NA	Water	8260B	
LCS 240-372032/4	Lab Control Sample	Total/NA	Water	8260B	
MRL 240-372032/5	Lab Control Sample	Total/NA	Water	8260B	
240-108954-D-4 MS	Matrix Spike	Total/NA	Water	8260B	
240-108954-D-4 MSD	Matrix Spike Duplicate	Total/NA	Water	8260B	

Lab Chronicle

Client: ARCADIS U.S., Inc.
Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

Client Sample ID: MW-72_030519

Date Collected: 03/05/19 14:20

Date Received: 03/08/19 08:00

Lab Sample ID: 240-109102-1

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	8260B		1	372032	03/18/19 16:27	LEE	TAL CAN
Total/NA	Analysis	8260B SIM		1	371373	03/13/19 12:28	SAM	TAL CAN

Client Sample ID: MW-72S_030519

Date Collected: 03/05/19 15:50

Date Received: 03/08/19 08:00

Lab Sample ID: 240-109102-2

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	8260B		1	371553	03/14/19 18:59	LEE	TAL CAN
Total/NA	Analysis	8260B SIM		1	371373	03/13/19 12:53	SAM	TAL CAN

Client Sample ID: MW-83_030519

Date Collected: 03/05/19 12:13

Date Received: 03/08/19 08:00

Lab Sample ID: 240-109102-3

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	8260B		1	371553	03/14/19 19:22	LEE	TAL CAN
Total/NA	Analysis	8260B SIM		1	371373	03/13/19 13:19	SAM	TAL CAN

Client Sample ID: MW-83S_030519

Date Collected: 03/05/19 14:07

Date Received: 03/08/19 08:00

Lab Sample ID: 240-109102-4

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	8260B		1	371553	03/14/19 19:44	LEE	TAL CAN
Total/NA	Analysis	8260B SIM		1	371373	03/13/19 13:45	SAM	TAL CAN

Client Sample ID: MW-82SR_030519

Date Collected: 03/05/19 17:16

Date Received: 03/08/19 08:00

Lab Sample ID: 240-109102-5

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	8260B		1	371553	03/14/19 20:06	LEE	TAL CAN
Total/NA	Analysis	8260B SIM		1	371373	03/13/19 14:11	SAM	TAL CAN

Client Sample ID: TRIP BLANK

Date Collected: 03/05/19 00:00

Date Received: 03/08/19 08:00

Lab Sample ID: 240-109102-6

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	8260B		1	371553	03/14/19 20:28	LEE	TAL CAN

TestAmerica Canton

Lab Chronicle

Client: ARCADIS U.S., Inc.
Project/Site: Ford LTP Livonia MI - E203631

TestAmerica Job ID: 240-109102-1

Laboratory References:

TAL CAN = TestAmerica Canton, 4101 Shuffel Street NW, North Canton, OH 44720, TEL (330)497-9396

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Accreditation/Certification Summary

Client: ARCADIS U.S., Inc.

TestAmerica Job ID: 240-109102-1

Project/Site: Ford LTP Livonia MI - E203631

Laboratory: TestAmerica Canton

All accreditations/certifications held by this laboratory are listed. Not all accreditations/certifications are applicable to this report.

Authority	Program	EPA Region	Identification Number	Expiration Date
California	State Program	9	2927	02-23-20
Connecticut	State Program	1	PH-0590	12-31-19
Florida	NELAP	4	E87225	06-30-19
Illinois	NELAP	5	200004	07-31-19
Kansas	NELAP	7	E-10336	04-30-19 *
Kentucky (UST)	State Program	4	58	02-23-20
Kentucky (WW)	State Program	4	98016	12-31-19
Minnesota	NELAP	5	039-999-348	12-31-19 *
Minnesota (Petrofund)	State Program	1	3506	07-31-19
Nevada	State Program	9	OH00048	07-31-19
New Jersey	NELAP	2	OH001	06-30-19
New York	NELAP	2	10975	03-31-19 *
Ohio VAP	State Program	5	CL0024	09-06-19
Oregon	NELAP	10	4062	02-23-20
Pennsylvania	NELAP	3	68-00340	08-31-19 *
Texas	NELAP	6	T104704517-18-10	08-31-19
USDA	Federal		P330-16-00404	12-28-19
Virginia	NELAP	3	460175	09-14-19
Washington	State Program	10	C971	01-12-20 *
West Virginia DEP	State Program	3	210	12-31-19

* Accreditation/Certification renewal pending - accreditation/certification considered valid.

TestAmerica Canton

3.4/C3.2

Chain of Custody Record

TestAmerica Laboratory location: Brighton — 10448 Citation Drive, Suite 200 / Brighton, MI 48116 / 810-229-2763

Client Contact Company Name: Arcadis Address: 28550 Cabot Drive, Suite 500 City/State/Zip: Novi, MI, 48377 Phone: 248-994-2240		Regulatory program: <input type="checkbox"/> DW <input type="checkbox"/> NPDES <input type="checkbox"/> RCRA <input type="checkbox"/> Other	
Client Project Manager: Kris Hinskey Telephone: 248-994-2240 Email: kris@hinskey.com		Lab Contact: Mike DelMonico Telephone: 330-497-9396	
Project Name: Ford LTP Project Number: M1001454.0006.00003 PO # M1001454.0006.00003		TestAmerica Laboratories, Inc. COC No: _____ of _____ COCs For lab use only	
Method of Shipment/Carrier: Shipping/Tracking No:		Analyses Walk-in client Lab sampling Job/SDG No:	
Matrix Air <input type="checkbox"/> Solid <input type="checkbox"/> Other: _____ Aqueous <input type="checkbox"/> Sediment <input type="checkbox"/> Other: _____		Containers & Preservatives TAT if different from below: 10 day <input type="checkbox"/> 3 weeks <input type="checkbox"/> 2 weeks <input type="checkbox"/> 1 week <input type="checkbox"/> 2 days <input type="checkbox"/> 1 day <input type="checkbox"/>	
Sample Identification MW-72-030519 MW-725-030519 MW-83-030519 MW-835-030519 MW-825R-030519 TRIP BLANK	Sample Date 3/5/19 3/5/19 3/5/19 3/5/19 3/5/19	Sample Time 1420 1550 1213 1407 1716	H2SO4 <input type="checkbox"/> HNO3 <input type="checkbox"/> HCl <input type="checkbox"/> NaOH <input type="checkbox"/> ZnAc <input type="checkbox"/> NaOH <input type="checkbox"/> Other: _____ Filtered Sample (Y/N) <input type="checkbox"/> Composite C/Grab <input type="checkbox"/> 1,1-DCE 8260B <input type="checkbox"/> Cis-1,2-DCE 8260B <input type="checkbox"/> Trans-1,2-DCE 8260B <input type="checkbox"/> PCE 8260B <input type="checkbox"/> TCE 8260B <input type="checkbox"/> Vinyl Chloride 8260B <input type="checkbox"/> 1,4-Dioxane 8260B SIM <input type="checkbox"/>
Possible Hazard Identification <input checked="" type="checkbox"/> Non-Hazard <input type="checkbox"/> Flammable <input type="checkbox"/> Irritant <input type="checkbox"/> Poison B <input type="checkbox"/> Unknown		Sample Disposal (A fee may apply) <input type="checkbox"/> Return to Client	
Special Instructions/ QC Requirements & Comments: Submit all results through Cadena at jim.tomalia@cadena.com, Cadena #E203631 Level IV Reporting.			
Relinquished by: RACHEL BIELAK Rachel Biellak		Company: ARCADIS Date/Time: 3/6/19 0945	
Relinquished by: NOVI COLD STORAGE Rodulfo		Company: ARCADIS Date/Time: 03/07/19 12:52	
Relinquished by: [Signature]		Company: TESTAMERICA Date/Time: 3/7/19 14:30	
Received in Laboratory by: [Signature]		Company: TESTAMERICA Date/Time: 3/7/19 1252	
Received by: [Signature]		Company: ARCADIS Date/Time: 3/8/19 800	



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TestAmerica Canton Sample Receipt Form/Narrative

Login # : 109102

Canton Facility

Client Acadvis Site Name _____

Cooler unpacked by: [Signature]


Cooler Received on 3/8/14 Opened on 3/8/14

FedEx: 1st Grd Exp UPS FAS Clipper Client Drop Off TestAmerica Courier Other

Receipt After-hours: Drop-off Date/Time

Storage Location

TestAmerica Cooler # 770 Foam Box Client Cooler Box Other _____
Packing material used: Bubble Wrap Foam Plastic Bag None Other _____
COOLANT: Wet Ice Blue Ice Dry Ice Water None

- Cooler temperature upon receipt See Multiple Cooler Form
IR GUN# IR-8 (CF -0.2 °C) Observed Cooler Temp. 3.4 °C Corrected Cooler Temp. 3.2 °C
IR GUN #36 (CF +0.7°C) Observed Cooler Temp. _____ °C Corrected Cooler Temp. _____ °C
- Were tamper/custody seals on the outside of the cooler(s)? If Yes Quantity 1 Yes No
-Were the seals on the outside of the cooler(s) signed & dated? Yes No NA
-Were tamper/custody seals on the bottle(s) or bottle kits (LLHg/MeHg)? Yes No NA
-Were tamper/custody seals intact and uncompromised? Yes No NA
- Shippers' packing slip attached to the cooler(s)? Yes No
- Did custody papers accompany the sample(s)? Yes No
- Were the custody papers relinquished & signed in the appropriate place? Yes No
- Was/were the person(s) who collected the samples clearly identified on the COC? Yes No
- Did all bottles arrive in good condition (Unbroken)? Yes No
- Could all bottle labels be reconciled with the COC? Yes No
- Were correct bottle(s) used for the test(s) indicated? Yes No
- Sufficient quantity received to perform indicated analyses? Yes No
- Are these work share samples? Yes No
If yes, Questions 12-16 have been checked at the originating laboratory.
- Were all preserved sample(s) at the correct pH upon receipt? Yes No NA pH Strip Lot# HC861525
- Were VOAs on the COC? Yes No NA
- Were air bubbles >6 mm in any VOA vials? Yes No NA  ← Larger than this.
- Was a VOA trip blank present in the cooler(s)? Trip Blank Lot # 831701 Yes No
- Was a LL Hg or Me Hg trip blank present? Yes No

Tests that are not checked for pH by Receiving:
VOAs
Oil and Grease
TOC

Contacted PM _____ Date _____ by _____ via Verbal Voice Mail Other

Concerning _____

17. CHAIN OF CUSTODY & SAMPLE DISCREPANCIES

Samples processed by:

AM

18. SAMPLE CONDITION

Sample(s) _____ were received after the recommended holding time had expired.
Sample(s) _____ were received in a broken container.
Sample(s) _____ were received with bubble >6 mm in diameter. (Notify PM)

19. SAMPLE PRESERVATION

Sample(s) _____ were further preserved in the laboratory.

Time preserved: _____ Preservative(s) added/Lot number(s): _____

VOA Sample Preservation - Date/Time VOAs Frozen: _____