APPENDIX H

GROUNDWATER MONITORING DOCUMENTS AND DUSRs

FPM

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSCC

Project No.:

Location: 90-30 metro

Well No.: A-06 Well Diameter: 2 in

Date: 7/23/07 Start Time:

Weather: Rain 73°F Finish Time:

Sampled By: JB/BC

Depth to Bottom of Well: 58.50 Feet.

Depth to Water: 42.37 Feet.

Height of Water Column: 16.13 Feet.

Water Volume in Casing: 2.58 Gallons.

Water Volume to be Purged: 7.74 Gallons.

Water Volume Actually Purged: 8 Gallons.

Purge Method: Submersible Pump

Physical Appearance/Comments:

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	2	7.20	737	62.5	120
	4	7.39	840	61.0	37.3
	8	7.43	837	60.7	120

Samplingand Analytical Methods: Disposable bailer cocs

Laboratory Name and Location: STL CT.

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSCo

Project No.: _____

Location: 90-30 Metropolitan

Well No.: A-7 Well Diameter: 2"

Date: 7/23/07 Start Time: _____

Weather: Rain Finish Time: _____

Sampled By: BC/JB

Depth to Bottom of Well: 115 Feet.

Depth to Water: 43.67 Feet.

Height of Water Column: 71.33 Feet.

Water Volume in Casing: 11.41 Gallons.

Water Volume to be Purged: 34.23 Gallons.

Water Volume Actually Purged: 35 Gallons.

Purge Method: _____

Physical Appearance/Comments: Rotten egg odor

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	11	7.51	709	61.2	85
	22	7.12	704	61.0	41
	34	7.08	703	60.8	12

Samplingand Analytical Methods: Bailer VOC's

Laboratory Name and Location: STL CT.

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: PPSCo

Project No.: _____

Location: 9030 Metropolitan Ave

Well No.: A-8 Well Diameter: 2"

Date: 7/23/07 Start Time: _____

Weather: Rain 70° Finish Time: _____

Sampled By: BC/JB

Depth to Bottom of Well: 82.30 Feet.

Depth to Water: 44.10 Feet.

Height of Water Column: 38.20 Feet.

Water Volume in Casing: 6.11 Gallons.

Water Volume to be Purged: 18.33 Gallons.

Water Volume Actually Purged: 19. Gallons.

Purge Method: Submersible Pump

Physical Appearance/Comments: _____

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	6	7.61	1030	62.0	25
	12	7.45	1033	61.6	20
	18	7.42	1031	61.2	10

Samplingand Analytical Methods: Bailer / VOCs

Laboratory Name and Location: STL CT.

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSC

Project No.: _____

Location: Metro

Well No.: A-5 Well Diameter: 2"

Date: 7/23/07 Start Time: _____

Weather: Rain 70° Finish Time: _____

Sampled By: BC/JB

Depth to Bottom of Well: 57.60 Feet.

Depth to Water: 42.32 Feet.

Height of Water Column: 15.28 Feet.

Water Volume in Casing: 2.45 Gallons.

Water Volume to be Purged: 7.33 Gallons.

Water Volume Actually Purged: 8 Gallons.

Purge Method: _____

Physical Appearance/Comments: _____

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	3	7.12	762	62.5	210
	5	7.42	771	61.7	200
	8	7.39	760	61.6	46

Samplingand Analytical Methods: _____

Laboratory Name and Location: _____

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPS W

Project No.: _____

Location: 90-30 Metropolitan

Well No.: A-02 Well Diameter: 2 In

Date: 7/23/07 Start Time: _____

Weather: Rain 73°F Finish Time: _____

Sampled By: JB/BC

Depth to Bottom of Well: 58.75 Feet.

Depth to Water: 42.2 Feet.

Height of Water Column: 16.55 Feet.

Water Volume in Casing: 2.64 Gallons.

Water Volume to be Purged: 7.94 Gallons.

Water Volume Actually Purged: 5 Gallons.

Purge Method: Submersible pump

Physical Appearance/Comments: _____

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	3	7.23	690	61.9	551
	5	7.31	700	60.8	200
	8	7.45	715	60.5	117

Samplingand Disposal Bailer / VOCs Analytical Methods:

Laboratory Name and Location: STL - CT

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSCA

Project No.: _____

Location: 90-30 MoRo

Well No.: A-3 Well Diameter: 2"

Date: 7/23/07 Start Time: _____

Weather: Rain 70° Finish Time: _____

Sampled By: BC/JB

Depth to Bottom of Well: 57.40 Feet.

Depth to Water: 42.44 Feet.

Height of Water Column: 15.04 Feet.

Water Volume in Casing: 2.41 Gallons.

Water Volume to be Purged: 7.21 Gallons.

Water Volume Actually Purged: 8 Gallons.

Purge Method: ED Sub pump

Physical Appearance/Comments: _____

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	3	7.24	595	60.1	63
	6	7.26	597	59.8	41
	8	7.26	594	59.6	12

Samplingand Analytical Methods: Bailer / VOC

Laboratory Name and Location: STL CT

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSCC

Project No.:

Location: 9030 Mapo

Well No.: A-11 Well Diameter: 2"

Date: 7/23/07 Start Time:

Weather: Rain 70° Finish Time:

Sampled By: BC / JH

Depth to Bottom of Well: 110 Feet.

Depth to Water: 37.4 Feet.

Height of Water Column: 72.6 Feet.

Water Volume in Casing: 11.62 Gallons.

Water Volume to be Purged: 34.85 Gallons.

Water Volume Actually Purged: 36 Gallons.

Purge Method: Sub. pump

Physical Appearance/Comments:

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	12	7.38	85	67.4	76
	24	7.43	87	67.0	69
	36	7.40	87	67.1	91

Samplingand Analytical Methods: Baker / WDC's

Laboratory Name and Location: STL CT

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSA

Project No.: _____

Location: 90 30 Marks.

Well No.: MW-12 Well Diameter: 2"

Date: 7/23/03 Start Time: _____

Weather: rain 70° Finish Time: _____

Sampled By: BC/JB

Depth to Bottom of Well: 28.17 Feet.

Depth to Water: 41.03 Feet.

Height of Water Column: 32.14 Feet.

Water Volume in Casing: 5.34 Gallons.

Water Volume to be Purged: 17.8 Gallons.

Water Volume Actually Purged: 18 Gallons.

Purge Method: Sub. Pump

Physical Appearance/Comments: _____

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	6	7.45	1190	63.0	80
	12	7.41	1149	62.1	43
	18	7.40	1142	62.2	17

Samplingand Baler / VOC Analytical Methods:

Laboratory Name and Location: STL CT

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSCC

Project No.: _____

Location: 90 30 marcojohns

Well No.: mw-7 Well Diameter: 2"

Date: 7/23/07 Start Time: _____

Weather: rain 70° Finish Time: _____

Sampled By: BC/JB

Depth to Bottom of Well: 50.66 Feet.

Depth to Water: 41.03 Feet.

Height of Water Column: 9.63 Feet.

Water Volume in Casing: 1.54 Gallons.

Water Volume to be Purged: 4.62 Gallons.

Water Volume Actually Purged: 5 Gallons.

Purge Method: _____

Physical Appearance/Comments: _____

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	1	7.23	905	63.0	75
	3	7.21	900	61.2	41
	5	7.22	901	61.1	31

Samplingand Bailer VOCs Analytical Methods:

Laboratory Name and Location: STL CT

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPW

Project No.:

Location: 90-30 Metro

Well No.: ME-15 Well Diameter: 4"

Date: 7/23/07 Start Time:

Weather: Rain 70° Finish Time:

Sampled By: BC/JB

Depth to Bottom of Well: 54.86 Feet.

Depth to Water: 40.30 Feet.

Height of Water Column: 14.56 Feet.

Water Volume in Casing: 9.46 Gallons.

Water Volume to be Purged: 28.4 Gallons.

Water Volume Actually Purged: 30 Gallons.

Purge Method: Sub pump

Physical Appearance/Comments:

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	10	7.20	971	63.4	114
	20	7.13	966	62.3	40
	30	7.15	964	61.9	12

Samplingand Baiker/VOCs Analytical Methods:

Laboratory Name and Location: STL CT

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSCW

Project No.: _____

Location: 9030 Mapeo

Well No.: Well Well Diameter: 4"

Date: 7/23/07 Start Time: _____

Weather: Rain 70° Finish Time: _____

Sampled By: EC/JB

Depth to Bottom of Well: 54.68 Feet.

Depth to Water: 40.32 Feet.

Height of Water Column: 14.36 Feet.

Water Volume in Casing: 9.47 Gallons.

Water Volume to be Purged: 28.35 Gallons.

Water Volume Actually Purged: 30 Gallons.

Purge Method: Sub. Pump

Physical Appearance/Comments: _____

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	10	7.47	511	66.1	71
	20	7.39	705	63.5	48
	30	7.35	706	62.4	31

Samplingand Baker/USGS Analytical Methods:

Laboratory Name and Location: STL ET

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSW

Project No.:

Location: METROPOLITAN

Well No.: MW-110 Well Diameter: 2"

Date: 7/26/07 Start Time:

Weather: SUNNY HUMID Finish Time:

Sampled By: BC/AM

Depth to Bottom of Well: 110 Feet.

Depth to Water: 41.83 Feet.

Height of Water Column: 68.17 Feet.

Water Volume in Casing: 11 Gallons.

Water Volume to be Purged: 33 Gallons.

Water Volume Actually Purged: 35 Gallons.

Purge Method: SUBMERSIBLE

Physical Appearance/Comments:

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	12	7.56	978	60.9	
	24	7.67	982	60.3	
	35	7.47	991	59.9	

Samplingand Analytical Methods: VOC's

Laboratory Name and Location: STL Et.

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSW

Project No.:

Location: METROPOLITAN

Well No.: MW-109 Well Diameter: 2"

Date: 7/26/07 Start Time:

Weather: SUNNY HUMID Finish Time:

Sampled By: BC/AM

Depth to Bottom of Well: 85 Feet.

Depth to Water: 41.9 Feet.

Height of Water Column: 43.1 Feet.

Water Volume in Casing: 7 Gallons.

Water Volume to be Purged: 21 Gallons.

Water Volume Actually Purged: 21 Gallons.

Purge Method: SUBMERSIBLE

Physical Appearance/Comments:

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	7	7.14	1014	64.1	65
	14	7.37	1028	61.6	40
	21	7.22	1034	67.8	21

Samplingand Analytical Methods:

Laboratory Name and Location: STL CT

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSW

Project No.:

Location: METROPOLITAN

Well No.: A-18 Well Diameter: 2"

Date: 7/26/07 Start Time:

Weather: SUNNY HUMID Finish Time:

Sampled By: BC/AM

Depth to Bottom of Well: 50.50 Feet.

Depth to Water: 39.58 Feet.

Height of Water Column: 10.92 Feet.

Water Volume in Casing: 2 Gallons.

Water Volume to be Purged: 6 Gallons.

Water Volume Actually Purged: 6 Gallons.

Purge Method: SUBMERSIBLE

Physical Appearance/Comments:

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	2	7.66	332	66.5	340
	4	7.15	410	64.0	400
	6	7.41	421	62.8	48.

Samplingand VOC's Analytical Methods:

Laboratory Name and Location: STL Lt.

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSCE

Project No.: _____

Location: Metropolitan

Well No.: A-14 Well Diameter: 2"

Date: 7/26/07 Start Time: _____

Weather: Sunny Humid Finish Time: _____

Sampled By: BC/AM

Depth to Bottom of Well: 53.20 Feet.

Depth to Water: 40.41 Feet.

Height of Water Column: 12.79 Feet.

Water Volume in Casing: 2.04 Gallons.

Water Volume to be Purged: 6.12 Gallons.

Water Volume Actually Purged: 7 Gallons.

Purge Method: Sub. pump

Physical Appearance/Comments: _____

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	2	7.09	1116	66.5	400
	4	7.13	1149	65.1	150
	6	7.16	1171	65.0	90

Samplingand Analytical Methods: VOC's

Laboratory Name and Location: STL CT.

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSW

Project No.:

Location: METROPOLITAN

Well No.: A-15 Well Diameter: 2"

Date: 7/26/07 Start Time:

Weather: Sunny Humid Finish Time:

Sampled By: BC/AM

Depth to Bottom of Well: 52.60 Feet.

Depth to Water: 40.23 Feet.

Height of Water Column: 12.37 Feet.

Water Volume in Casing: 2 Gallons.

Water Volume to be Purged: 6 Gallons.

Water Volume Actually Purged: 6 Gallons.

Purge Method: Submersible Pump

Physical Appearance/Comments:

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	2	7.22	1048	65.7	320
	4	7.45	1164	65.5	170
	6	7.53	1151	68.2	70

Samplingand Analytical Methods:

Laboratory Name and Location:

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSW

Project No.:

Location: Metropolitan Ave

Well No.: A - 9 Well Diameter: 6"

Date: 7/26/07 Start Time:

Weather: Sunny Humid Finish Time:

Sampled By: BC/AM

Depth to Bottom of Well: 58.75 Feet.

Depth to Water: 42.00 Feet.

Height of Water Column: 16.75 Feet.

Water Volume in Casing: 2.68 Gallons.

Water Volume to be Purged: 8 Gallons.

Water Volume Actually Purged: 8 Gallons.

Purge Method: Sub. pump

Physical Appearance/Comments: Duplicate A-90

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	2.5	7.36	553	65.0	500
	5	7.85	574	62.7	321
	7.5	7.67	572	63.1	48

Samplingand VOC's Analytical Methods:

Laboratory Name and Location: STL

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSW

Project No.:

Location: Metropolitan Ave

Well No.: MW-113 Well Diameter: 2"

Date: 7/26/07 Start Time:

Weather: Sunny Humid Finish Time:

Sampled By: BC/AM

Depth to Bottom of Well: 131 Feet.

Depth to Water: 42.13 Feet.

Height of Water Column: 88.87 Feet.

Water Volume in Casing: 14.21 Gallons.

Water Volume to be Purged: 42.65 Gallons.

Water Volume Actually Purged: 43 Gallons.

Purge Method:

Physical Appearance/Comments:

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	14	7.36	937	64.8	130
	28	7.69	991	63.0	90
	42	7.59	999	62.5	40

Samplingand Analytical Methods:

Laboratory Name and Location:

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSC

Project No.:

Location: Metropolitan

Well No.: MW-112 Well Diameter: 2"

Date: 7/26/07 Start Time:

Weather: Sunny Hemp Finish Time:

Sampled By: BC/ATM

Depth to Bottom of Well: 129.15 Feet.

Depth to Water: 44.52 Feet.

Height of Water Column: 84.63 Feet.

Water Volume in Casing: 13.5 Gallons.

Water Volume to be Purged: 40.62 Gallons.

Water Volume Actually Purged: 41 Gallons.

Purge Method: Sub. pump

Physical Appearance/Comments:

MS - MSD

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	12	7.51	857	63.3	80
	24	7.62	889	62.9	30
	40	7.77	900	62.7	10

Samplingand VOC's Analytical Methods:

Laboratory Name and Location: STL CT.

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSW

Project No.:

Location: METROPOLITAN AVE

Well No.: ME - 8 Well Diameter: 4"

Date: 7/26/07 Start Time:

Weather: SUNNY HUMID Finish Time:

Sampled By: BC/AM

Depth to Bottom of Well: 55 Feet.

Depth to Water: 40.27 Feet.

Height of Water Column: 14.73 Feet.

Water Volume in Casing: 10 Gallons. .65

Water Volume to be Purged: 30 Gallons.

Water Volume Actually Purged: 30 Gallons.

Purge Method: SUBMERSIBLE PUMP

Physical Appearance/Comments:

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	10	7.28	627	64.2	146
	20	7.58	644	63.7	35
	30	7.52	654	63.5	5

Samplingand Analytical Methods: VOC's

Laboratory Name and Location: STL CT

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DSCC

Project No.:

Location: Methadone Island

Well No.: A-17 Well Diameter: 2"

Date: 7/24/07 Start Time:

Weather: Sunny Humid Finish Time:

Sampled By:

Depth to Bottom of Well: 110.00 Feet.

Depth to Water: 41.28 Feet.

Height of Water Column: 68.72 Feet.

Water Volume in Casing: 11. Gallons.

Water Volume to be Purged: 33 Gallons.

Water Volume Actually Purged: 35 Gallons.

Purge Method: Schmersible

Physical Appearance/Comments:

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	12	7.41	756	62.4	144
	24	7.39	917	59.8	70
	35	7.65	951	60.4	11

Samplingand VOC's Analytical Methods:

Laboratory Name and Location: STL CT

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSW

Project No.:

Location: METROPOLITAN

Well No.: A-16 Well Diameter: 2"

Date: 7/26/07 Start Time:

Weather: SUNNY HUMID Finish Time:

Sampled By:

Depth to Bottom of Well: 57.80 Feet.

Depth to Water: 39.62 Feet.

Height of Water Column: 12.18 Feet.

Water Volume in Casing: 2 Gallons.

Water Volume to be Purged: 6 Gallons.

Water Volume Actually Purged: 6 Gallons.

Purge Method: Submersible

Physical Appearance/Comments:

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	2	8.08	537	65.4	110
	4	7.77	476	64.2	73
	6	7.73	452	63.9	41

Samplingand VOC's Analytical Methods:

Laboratory Name and Location: STL CT

FPM group
Engineering and Environmental Science

WELL SAMPLING DATA FORM

Client: DPSCU

Project No.:

Location: Metropolitan

Well No.: ME-13 Well Diameter: 4"

Date: 7/26/07 Start Time:

Weather: Sunny/ Humid Finish Time:

Sampled By: BC/AM

Depth to Bottom of Well: 55.15 Feet.

Depth to Water: 40.38 Feet.

Height of Water Column: 14.77 Feet.

Water Volume in Casing: 9.6 Gallons.

Water Volume to be Purged: 28.8 Gallons.

Water Volume Actually Purged: 30 Gallons.

Purge Method: Sub. Pump

Physical Appearance/Comments:

FIELD MEASUREMENTS

Time	Volume (gal)	pH	Conductivity (uS)	Temperature (°F)	Turbidity (FTU)
	10	7.85	587	64.9	100
	20	7.76	606	65.8	40
	30	7.64	604	64.3	5

Samplingand Analytical Methods: VOCs

Laboratory Name and Location: STL CT

TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

ANALYTICAL REPORT

Job Number: 220-2277-1

SDG Number: 220-2277

Job Description: DPSW Metropolitan Avenue

For:
FPM Group Limited
909 Marconi Avenue
Ronkonkoma, NY 11779

Attention: Mr. Ben Cancemi

Mary Widomski

Designee for
Johanna Dubauskas
Project Manager I
johanna.dubauskas@testamericainc.com
08/09/2007

STL Connecticut Certifications and Approvals: CTDOH PH-047, MADEP CT023, RIDOH A43, NYDOH 10602, NY NELAP 10602, NHDES 2528, NJDEP CT410, ME DOH CT023, UT DOH 2032614458

TestAmerica Laboratories, Inc.
TestAmerica Connecticut 128 Long Hill Cross Road, Shelton, CT 06484
Tel (203) 929-8140 Fax (203) 929-8142 www.testamericainc.com

ACCREDITED IN ACCORDANCE WITH nelac 08/09/2007

Case Narrative For Job: 220-J2277-1

Client: FPM Group Limited
Date: August 9, 2007

I certify that this data package is in compliance with the terms and conditions of this contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or his designee, as verified by the following signature.

Peter Frick
Peter Frick
Laboratory Director

August 9, 2007
Date

Comments
No additional comments.

Receipt
All samples were received in good condition within temperature requirements.

GC/MS VOA
No analytical or quality issues were noted.

FORMULAS FOR NYSDEC SAMPLE CALCULATIONS

Volatiles

(Ax)(IS)(DF)
(AIS)(RRF)(V)(% solids) = C

(AX)(IS)(VT)(1000)(DF)
(AIS)(RRF)(VA)(V)(% solids) = C (for medium level soils)

SemiVolatiles

(AX)(IS)(VE)(DF)(GPC factor is 2 if needed)
(AIS)(RRF)(volume injected)(V)(% solids) = C

Pesticides

(AX)(VE)(DF)
(RRF)(V)(% solids)(volume injected) = C

PCBs for compound/retention time

(AX)(VE)(DF)
(RRF of compound at the stated retention time)(V)(% solids)(volume injected) = C

DRO/CTETPH

(AX)(VE)(DF)
(RRF)(V)(% solids)(volume injected) = C

AX = area of the target Ion
AIS = Area of Internal standard
C = concentration as ug/L or ug/Kg
DF = dilution
IS = Internal standard concentration (ng)
RRF = average RF (from initial cal except CLP methods from continuing cal)
V = sample volume for liquids in mls or sample weight for solids in grams
VA = volume of aliquot for medium level soils
VE = volume of concentrated extract
VT = volume of methanol for volatile medium level soils

Sample Summary Package

METHOD SUMMARY

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Description	Lab Location	Method	Preparation Method
Matrix: Water			
Volatile Organic Compounds by GC/MS	STL CT	SW846 8260B	
Purge-and-Trap	STL CT		SW846 5030B

LAB REFERENCES:

STL CT = TestAmerica Connecticut

METHOD REFERENCES:

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

METHOD / ANALYST SUMMARY

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Method	Analyst	Analyst ID
SW846 8260B	Humbert, Dave	DH

SAMPLE SUMMARY

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Lab Sample ID	Client Sample ID	Client Matrix	Date/Time Sampled	Date/Time Received
220-2277-1	A-2	Water	07/23/2007 0900	07/28/2007 1000
220-2277-2	A-6	Water	07/23/2007 0930	07/28/2007 1000
220-2277-3	A-7	Water	07/23/2007 1000	07/28/2007 1000
220-2277-4	A-8	Water	07/23/2007 1030	07/28/2007 1000
220-2277-5	A-3	Water	07/23/2007 1100	07/28/2007 1000
220-2277-6	A-11	Water	07/23/2007 1130	07/28/2007 1000
220-2277-7	A-12	Water	07/23/2007 1230	07/28/2007 1000
220-2277-8	ME-7	Water	07/23/2007 1300	07/28/2007 1000
220-2277-9	ME-11	Water	07/23/2007 1330	07/28/2007 1000
220-2277-10FB	FB0723	Water	07/23/2007 1400	07/28/2007 1000
220-2277-11TB	TRIP BLANK	Water	07/23/2007 0000	07/28/2007 1000
220-2277-12	A-16	Water	07/26/2007 0800	07/28/2007 1000
220-2277-13	A-17	Water	07/26/2007 0830	07/28/2007 1000
220-2277-14	MW-109	Water	07/26/2007 0900	07/28/2007 1000
220-2277-15	MW-110	Water	07/26/2007 0930	07/28/2007 1000
220-2277-16	A-18	Water	07/26/2007 1000	07/28/2007 1000
220-2277-17	A-14	Water	07/26/2007 1030	07/28/2007 1000
220-2277-18	A-15	Water	07/26/2007 1100	07/28/2007 1000
220-2277-19	A-9	Water	07/26/2007 1130	07/28/2007 1000
220-2277-20	A-10	Water	07/26/2007 1200	07/28/2007 1000
220-2277-21	MW-113	Water	07/26/2007 1300	07/28/2007 1000
220-2277-22	MW-112	Water	07/26/2007 1330	07/28/2007 1000
220-2277-23	ME-8	Water	07/26/2007 1400	07/28/2007 1000
220-2277-24	ME-13	Water	07/26/2007 1430	07/28/2007 1000
220-2277-25FB	FB0726	Water	07/26/2007 0000	07/28/2007 1000
220-2277-26	A-9D	Water	07/26/2007 1145	07/28/2007 1000
220-2277-27	A-5	Water	07/23/2007 1045	07/28/2007 1000
220-2277-28	ME-15	Water	07/23/2007 1200	07/28/2007 1000

SAMPLE RESULTS

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: A-2

Lab Sample ID: 220-2277-1
Client Matrix: Water

Date Sampled: 07/23/2007 0900
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/01/2007 0348
Date Prepared: 08/01/2007 0348

Analysis Batch: 220-8307

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9328.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	32		0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	74	53 - 125
4-Bromofluorobenzene	118	73 - 127
Dibromofluoromethane	75	54 - 137
Toluene-d8 (Surr)	82	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: A-6

Lab Sample ID: 220-2277-2
Client Matrix: Water

Date Sampled: 07/23/2007 0930
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 2.0
Date Analyzed: 08/02/2007 1648
Date Prepared: 08/02/2007 1648

Analysis Batch: 220-8356

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9386.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	20	U	2.8	20
Benzene	10	U	0.80	10
Bromodichloromethane	10	U	0.80	10
Bromoform	10	U	1.6	10
Bromomethane	10	U	2.4	10
2-Butanone (MEK)	20	U	2.4	20
Carbon disulfide	10	U	1.8	10
Carbon tetrachloride	10	U	2.0	10
Chlorobenzene	10	U	0.80	10
Chloroethane	10	U	1.6	10
Chloroform	10	U	1.4	10
Chloromethane	10	U	1.0	10
Dibromochloromethane	10	U	1.0	10
1,1-Dichloroethane	10	U	1.2	10
1,2-Dichloroethane	10	U	1.2	10
1,1-Dichloroethene	10	U	1.4	10
1,2-Dichloropropane	10	U	1.8	10
cis-1,3-Dichloropropene	10	U	1.0	10
trans-1,3-Dichloropropene	10	U	1.6	10
Ethylbenzene	10	U	2.0	10
2-Hexanone	20	U	1.6	20
Methylene Chloride	10	U	0.80	10
4-Methyl-2-pentanone (MIBK)	20	U	1.4	20
Styrene	10	U	1.0	10
1,1,2,2-Tetrachloroethane	10	U	0.80	10
Tetrachloroethene	290		1.0	10
Toluene	10	U	0.60	10
1,1,1-Trichloroethane	10	U	0.80	10
1,1,2-Trichloroethane	10	U	1.2	10
Trichloroethene	10	U	1.4	10
Vinyl chloride	10	U	1.6	10
Xylenes, Total	10	U	2.0	10
cis-1,2-Dichloroethene	10	U	1.2	10
trans-1,2-Dichloroethene	10	U	1.0	10

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	76	53 - 125
4-Bromofluorobenzene	120	73 - 127
Dibromofluoromethane	78	54 - 137
Toluene-d8 (Surr)	83	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: A-7

Lab Sample ID: 220-2277-3
Client Matrix: Water

Date Sampled: 07/23/2007 1000
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/01/2007 0438
Date Prepared: 08/01/2007 0438

Analysis Batch: 220-8307

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9330.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	3.8	J	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	26		1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	41		0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	74	53 - 125
4-Bromofluorobenzene	112	73 - 127
Dibromofluoromethane	75	54 - 137
Toluene-d8 (Surr)	83	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: A-8

Lab Sample ID: 220-2277-4
Client Matrix: Water

Date Sampled: 07/23/2007 1030
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 50
Date Analyzed: 08/02/2007 1508
Date Prepared: 08/02/2007 1508

Analysis Batch: 220-8356

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9382.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	97	J	70	500
Benzene	250	U	20	250
Bromodichloromethane	250	U	20	250
Bromoform	250	U	40	250
Bromomethane	250	U	60	250
2-Butanone (MEK)	500	U	60	500
Carbon disulfide	250	U	45	250
Carbon tetrachloride	250	U	50	250
Chlorobenzene	250	U	20	250
Chloroethane	250	U	40	250
Chloroform	250	U	35	250
Chloromethane	250	U	25	250
Dibromochloromethane	250	U	25	250
1,1-Dichloroethane	250	U	30	250
1,2-Dichloroethane	250	U	30	250
1,1-Dichloroethene	250	U	35	250
1,2-Dichloropropane	250	U	45	250
cis-1,3-Dichloropropene	250	U	25	250
trans-1,3-Dichloropropene	250	U	40	250
Ethylbenzene	250	U	50	250
2-Hexanone	500	U	40	500
Methylene Chloride	250	U	20	250
4-Methyl-2-pentanone (MIBK)	500	U	35	500
Styrene	250	U	25	250
1,1,2,2-Tetrachloroethane	250	U	20	250
Tetrachloroethene	4400		25	250
Toluene	250	U	15	250
1,1,1-Trichloroethane	250	U	20	250
1,1,2-Trichloroethane	250	U	30	250
Trichloroethene	250	U	35	250
Vinyl chloride	250	U	40	250
Xylenes, Total	250	U	50	250
cis-1,2-Dichloroethene	250	U	30	250
trans-1,2-Dichloroethene	250	U	25	250

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	74	53 - 125
4-Bromofluorobenzene	117	73 - 127
Dibromofluoromethane	75	54 - 137
Toluene-d8 (Surr)	84	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: A-3

Lab Sample ID: 220-2277-5
Client Matrix: Water

Date Sampled: 07/23/2007 1100
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/02/2007 1232
Date Prepared: 08/02/2007 1232

Analysis Batch: 220-8356

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9376.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	1.4	J	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	14		0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	71	53 - 125
4-Bromofluorobenzene	112	73 - 127
Dibromofluoromethane	72	54 - 137
Toluene-d8 (Surr)	78	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: A-11

Lab Sample ID: 220-2277-6
Client Matrix: Water

Date Sampled: 07/23/2007 1130
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/02/2007 1257
Date Prepared: 08/02/2007 1257

Analysis Batch: 220-8356

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9377.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	8.5	J	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	0.78	J	0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	73	53 - 125
4-Bromofluorobenzene	116	73 - 127
Dibromofluoromethane	76	54 - 137
Toluene-d8 (Surr)	81	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: A-12

Lab Sample ID: 220-2277-7
Client Matrix: Water

Date Sampled: 07/23/2007 1230
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 20
Date Analyzed: 08/02/2007 1533
Date Prepared: 08/02/2007 1533

Analysis Batch: 220-8356

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9383.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	200	U	28	200
Benzene	100	U	8.0	100
Bromodichloromethane	100	U	8.0	100
Bromoform	100	U	16	100
Bromomethane	100	U	24	100
2-Butanone (MEK)	200	U	24	200
Carbon disulfide	100	U	18	100
Carbon tetrachloride	100	U	20	100
Chlorobenzene	100	U	8.0	100
Chloroethane	100	U	16	100
Chloroform	100	U	14	100
Chloromethane	100	U	10	100
Dibromochloromethane	100	U	10	100
1,1-Dichloroethane	100	U	12	100
1,2-Dichloroethane	100	U	12	100
1,1-Dichloroethene	100	U	14	100
1,2-Dichloropropane	100	U	18	100
cis-1,3-Dichloropropene	100	U	10	100
trans-1,3-Dichloropropene	100	U	16	100
Ethylbenzene	100	U	20	100
2-Hexanone	200	U	16	200
Methylene Chloride	100	U	8.0	100
4-Methyl-2-pentanone (MIBK)	200	U	14	200
Styrene	100	U	10	100
1,1,2,2-Tetrachloroethane	100	U	8.0	100
Tetrachloroethene	1800		10	100
Toluene	100	U	6.0	100
1,1,1-Trichloroethane	100	U	8.0	100
1,1,2-Trichloroethane	100	U	12	100
Trichloroethene	100	U	14	100
Vinyl chloride	100	U	16	100
Xylenes, Total	100	U	20	100
cis-1,2-Dichloroethene	100	U	12	100
trans-1,2-Dichloroethene	100	U	10	100

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	75	53 - 125
4-Bromofluorobenzene	112	73 - 127
Dibromofluoromethane	72	54 - 137
Toluene-d8 (Surr)	81	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: ME-7

Lab Sample ID: 220-2277-8
Client Matrix: Water

Date Sampled: 07/23/2007 1300
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/02/2007 1322
Date Prepared: 08/02/2007 1322

Analysis Batch: 220-8356

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9378.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	0.88	J	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	41		0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	77	53 - 125
4-Bromofluorobenzene	116	73 - 127
Dibromofluoromethane	75	54 - 137
Toluene-d8 (Surr)	82	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: ME-11

Lab Sample ID: 220-2277-9
Client Matrix: Water

Date Sampled: 07/23/2007 1330
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 2.0
Date Analyzed: 08/02/2007 1713
Date Prepared: 08/02/2007 1713

Analysis Batch: 220-8356

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9387.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	20	U	2.8	20
Benzene	10	U	0.80	10
Bromodichloromethane	10	U	0.80	10
Bromoform	10	U	1.6	10
Bromomethane	10	U	2.4	10
2-Butanone (MEK)	20	U	2.4	20
Carbon disulfide	10	U	1.8	10
Carbon tetrachloride	10	U	2.0	10
Chlorobenzene	10	U	0.80	10
Chloroethane	10	U	1.6	10
Chloroform	10	U	1.4	10
Chloromethane	10	U	1.0	10
Dibromochloromethane	10	U	1.0	10
1,1-Dichloroethane	10	U	1.2	10
1,2-Dichloroethane	10	U	1.2	10
1,1-Dichloroethene	10	U	1.4	10
1,2-Dichloropropane	10	U	1.8	10
cis-1,3-Dichloropropene	10	U	1.0	10
trans-1,3-Dichloropropene	10	U	1.6	10
Ethylbenzene	10	U	2.0	10
2-Hexanone	20	U	1.6	20
Methylene Chloride	10	U	0.80	10
4-Methyl-2-pentanone (MIBK)	20	U	1.4	20
Styrene	10	U	1.0	10
1,1,2,2-Tetrachloroethane	10	U	0.80	10
Tetrachloroethene	290		1.0	10
Toluene	10	U	0.60	10
1,1,1-Trichloroethane	10	U	0.80	10
1,1,2-Trichloroethane	10	U	1.2	10
Trichloroethene	10	U	1.4	10
Vinyl chloride	10	U	1.6	10
Xylenes, Total	10	U	2.0	10
cis-1,2-Dichloroethene	10	U	1.2	10
trans-1,2-Dichloroethene	10	U	1.0	10

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	73	53 - 125
4-Bromofluorobenzene	111	73 - 127
Dibromofluoromethane	75	54 - 137
Toluene-d8 (Surr)	80	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: FB0723

Lab Sample ID: 220-2277-10FB
Client Matrix: Water

Date Sampled: 07/23/2007 1400
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/01/2007 0259
Date Prepared: 08/01/2007 0259

Analysis Batch: 220-8307

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9326.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	0.54	J	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	5.0	U	0.50	5.0
Toluene	0.40	J	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	71	53 - 125
4-Bromofluorobenzene	110	73 - 127
Dibromofluoromethane	75	54 - 137
Toluene-d8 (Surr)	80	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: TRIP BLANK

Lab Sample ID: 220-2277-11TB
Client Matrix: Water

Date Sampled: 07/23/2007 0000
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/01/2007 0323
Date Prepared: 08/01/2007 0323

Analysis Batch: 220-8307

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9327.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	3.1	J	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	3.6	J	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	2.8	J	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	5.0	U	0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	77	53 - 125
4-Bromofluorobenzene	121	73 - 127
Dibromofluoromethane	74	54 - 137
Toluene-d8 (Surr)	86	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: A-16

Lab Sample ID: 220-2277-12
Client Matrix: Water

Date Sampled: 07/26/2007 0800
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/02/2007 1347
Date Prepared: 08/02/2007 1347

Analysis Batch: 220-8356

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9379.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	5.0	U	0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	76	53 - 125
4-Bromofluorobenzene	115	73 - 127
Dibromofluoromethane	78	54 - 137
Toluene-d8 (Surr)	84	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: A-17

Lab Sample ID: 220-2277-13
Client Matrix: Water

Date Sampled: 07/26/2007 0830
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/01/2007 1445
Date Prepared: 08/01/2007 1445

Analysis Batch: 220-8321

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9351.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	0.79	J	0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	72	53 - 125
4-Bromofluorobenzene	107	73 - 127
Dibromofluoromethane	74	54 - 137
Toluene-d8 (Surr)	83	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: **MW-109**

Lab Sample ID: 220-2277-14
Client Matrix: Water

Date Sampled: 07/26/2007 0900
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/01/2007 1510
Date Prepared: 08/01/2007 1510

Analysis Batch: 220-8321

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9352.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	2.6	J	0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	71	53 - 125
4-Bromofluorobenzene	114	73 - 127
Dibromofluoromethane	72	54 - 137
Toluene-d8 (Surr)	81	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: **MW-110**

Lab Sample ID: 220-2277-15
Client Matrix: Water

Date Sampled: 07/26/2007 0930
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/01/2007 1535
Date Prepared: 08/01/2007 1535

Analysis Batch: 220-8321

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9353.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	5.0	U	0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	75	53 - 125
4-Bromofluorobenzene	112	73 - 127
Dibromofluoromethane	78	54 - 137
Toluene-d8 (Surr)	85	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: **A-18**

Lab Sample ID: 220-2277-16
Client Matrix: Water

Date Sampled: 07/26/2007 1000
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/01/2007 1600
Date Prepared: 08/01/2007 1600

Analysis Batch: 220-8321

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9354.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	5.0	U	0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	71	53 - 125
4-Bromofluorobenzene	107	73 - 127
Dibromofluoromethane	73	54 - 137
Toluene-d8 (Surr)	78	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: A-14

Lab Sample ID: 220-2277-17
Client Matrix: Water

Date Sampled: 07/26/2007 1030
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 4.0
Date Analyzed: 08/02/2007 1623
Date Prepared: 08/02/2007 1623

Analysis Batch: 220-8356

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9385.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	40	U	5.6	40
Benzene	20	U	1.6	20
Bromodichloromethane	20	U	1.6	20
Bromoform	20	U	3.2	20
Bromomethane	20	U	4.8	20
2-Butanone (MEK)	40	U	4.8	40
Carbon disulfide	20	U	3.6	20
Carbon tetrachloride	20	U	4.0	20
Chlorobenzene	20	U	1.6	20
Chloroethane	20	U	3.2	20
Chloroform	20	U	2.8	20
Chloromethane	20	U	2.0	20
Dibromochloromethane	20	U	2.0	20
1,1-Dichloroethane	20	U	2.4	20
1,2-Dichloroethane	20	U	2.4	20
1,1-Dichloroethene	20	U	2.8	20
1,2-Dichloropropane	20	U	3.6	20
cis-1,3-Dichloropropene	20	U	2.0	20
trans-1,3-Dichloropropene	20	U	3.2	20
Ethylbenzene	20	U	4.0	20
2-Hexanone	40	U	3.2	40
Methylene Chloride	20	U	1.6	20
4-Methyl-2-pentanone (MIBK)	40	U	2.8	40
Styrene	20	U	2.0	20
1,1,2,2-Tetrachloroethane	20	U	1.6	20
Tetrachloroethene	470		2.0	20
Toluene	20	U	1.2	20
1,1,1-Trichloroethane	20	U	1.6	20
1,1,2-Trichloroethane	20	U	2.4	20
Trichloroethene	20	U	2.8	20
Vinyl chloride	20	U	3.2	20
Xylenes, Total	20	U	4.0	20
cis-1,2-Dichloroethene	20	U	2.4	20
trans-1,2-Dichloroethene	20	U	2.0	20

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	76	53 - 125
4-Bromofluorobenzene	115	73 - 127
Dibromofluoromethane	76	54 - 137
Toluene-d8 (Surr)	83	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: **A-15**

Lab Sample ID: 220-2277-18
Client Matrix: Water

Date Sampled: 07/26/2007 1100
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/02/2007 1412
Date Prepared: 08/02/2007 1412

Analysis Batch: 220-8356

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9380.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	10		0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	75	53 - 125
4-Bromofluorobenzene	113	73 - 127
Dibromofluoromethane	75	54 - 137
Toluene-d8 (Surr)	80	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: A-9

Lab Sample ID: 220-2277-19
Client Matrix: Water

Date Sampled: 07/26/2007 1130
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/01/2007 1715
Date Prepared: 08/01/2007 1715

Analysis Batch: 220-8321

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9357.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	23		0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	72	53 - 125
4-Bromofluorobenzene	110	73 - 127
Dibromofluoromethane	73	54 - 137
Toluene-d8 (Surr)	80	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: A-10

Lab Sample ID: 220-2277-20
Client Matrix: Water

Date Sampled: 07/26/2007 1200
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/01/2007 1740
Date Prepared: 08/01/2007 1740

Analysis Batch: 220-8321

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9358.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	10		0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	73	53 - 125
4-Bromofluorobenzene	109	73 - 127
Dibromofluoromethane	74	54 - 137
Toluene-d8 (Surr)	80	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: **MW-113**

Lab Sample ID: 220-2277-21
Client Matrix: Water

Date Sampled: 07/26/2007 1300
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/01/2007 1804
Date Prepared: 08/01/2007 1804

Analysis Batch: 220-8321

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9359.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	3.4	J B	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	8.7		0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	73	53 - 125
4-Bromofluorobenzene	110	73 - 127
Dibromofluoromethane	75	54 - 137
Toluene-d8 (Surr)	81	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: MW-112

Lab Sample ID: 220-2277-22
Client Matrix: Water

Date Sampled: 07/26/2007 1330
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/02/2007 1436
Date Prepared: 08/02/2007 1436

Analysis Batch: 220-8356

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9381.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	1.3	J	0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	75	53 - 125
4-Bromofluorobenzene	115	73 - 127
Dibromofluoromethane	77	54 - 137
Toluene-d8 (Surr)	84	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: ME-8

Lab Sample ID: 220-2277-23
Client Matrix: Water

Date Sampled: 07/26/2007 1400
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 2.0
Date Analyzed: 08/02/2007 1737
Date Prepared: 08/02/2007 1737

Analysis Batch: 220-8356

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9388.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	20	U	2.8	20
Benzene	10	U	0.80	10
Bromodichloromethane	10	U	0.80	10
Bromoform	10	U	1.6	10
Bromomethane	10	U	2.4	10
2-Butanone (MEK)	20	U	2.4	20
Carbon disulfide	10	U	1.8	10
Carbon tetrachloride	10	U	2.0	10
Chlorobenzene	10	U	0.80	10
Chloroethane	10	U	1.6	10
Chloroform	10	U	1.4	10
Chloromethane	10	U	1.0	10
Dibromochloromethane	10	U	1.0	10
1,1-Dichloroethane	10	U	1.2	10
1,2-Dichloroethane	10	U	1.2	10
1,1-Dichloroethene	10	U	1.4	10
1,2-Dichloropropane	10	U	1.8	10
cis-1,3-Dichloropropene	10	U	1.0	10
trans-1,3-Dichloropropene	10	U	1.6	10
Ethylbenzene	10	U	2.0	10
2-Hexanone	20	U	1.6	20
Methylene Chloride	10	U	0.80	10
4-Methyl-2-pentanone (MIBK)	20	U	1.4	20
Styrene	10	U	1.0	10
1,1,2,2-Tetrachloroethane	10	U	0.80	10
Tetrachloroethene	250		1.0	10
Toluene	10	U	0.60	10
1,1,1-Trichloroethane	10	U	0.80	10
1,1,2-Trichloroethane	10	U	1.2	10
Trichloroethene	10	U	1.4	10
Vinyl chloride	10	U	1.6	10
Xylenes, Total	10	U	2.0	10
cis-1,2-Dichloroethene	10	U	1.2	10
trans-1,2-Dichloroethene	10	U	1.0	10

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	72	53 - 125
4-Bromofluorobenzene	108	73 - 127
Dibromofluoromethane	74	54 - 137
Toluene-d8 (Surr)	80	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: ME-13

Lab Sample ID: 220-2277-24
Client Matrix: Water

Date Sampled: 07/26/2007 1430
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/01/2007 1919
Date Prepared: 08/01/2007 1919

Analysis Batch: 220-8321

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9362.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	54		0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	80	53 - 125
4-Bromofluorobenzene	120	73 - 127
Dibromofluoromethane	80	54 - 137
Toluene-d8 (Surr)	85	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: FB0726

Lab Sample ID: 220-2277-25FB
Client Matrix: Water

Date Sampled: 07/26/2007 0000
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/01/2007 1944
Date Prepared: 08/01/2007 1944

Analysis Batch: 220-8321

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9363.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	3.0	J B	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	0.76	J B	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	5.0	U	0.50	5.0
Toluene	0.37	J	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	75	53 - 125
4-Bromofluorobenzene	113	73 - 127
Dibromofluoromethane	74	54 - 137
Toluene-d8 (Surr)	84	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: **A-9D**

Lab Sample ID: 220-2277-26
Client Matrix: Water

Date Sampled: 07/26/2007 1145
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/01/2007 2008
Date Prepared: 08/01/2007 2008

Analysis Batch: 220-8321

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9364.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	22		0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	72	53 - 125
4-Bromofluorobenzene	119	73 - 127
Dibromofluoromethane	77	54 - 137
Toluene-d8 (Surr)	83	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: **A-5**

Lab Sample ID: 220-2277-27
Client Matrix: Water

Date Sampled: 07/23/2007 1045
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 08/01/2007 2033
Date Prepared: 08/01/2007 2033

Analysis Batch: 220-8321

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9365.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	39		0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	71	53 - 125
4-Bromofluorobenzene	107	73 - 127
Dibromofluoromethane	75	54 - 137
Toluene-d8 (Surr)	80	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Client Sample ID: **ME-15**

Lab Sample ID: 220-2277-28
Client Matrix: Water

Date Sampled: 07/23/2007 1200
Date Received: 07/28/2007 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 5.0
Date Analyzed: 08/02/2007 1558
Date Prepared: 08/02/2007 1558

Analysis Batch: 220-8356

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9384.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Acetone	50	U	7.0	50
Benzene	25	U	2.0	25
Bromodichloromethane	25	U	2.0	25
Bromoform	25	U	4.0	25
Bromomethane	25	U	6.0	25
2-Butanone (MEK)	50	U	6.0	50
Carbon disulfide	25	U	4.5	25
Carbon tetrachloride	25	U	5.0	25
Chlorobenzene	25	U	2.0	25
Chloroethane	25	U	4.0	25
Chloroform	25	U	3.5	25
Chloromethane	25	U	2.5	25
Dibromochloromethane	25	U	2.5	25
1,1-Dichloroethane	25	U	3.0	25
1,2-Dichloroethane	25	U	3.0	25
1,1-Dichloroethene	25	U	3.5	25
1,2-Dichloropropane	25	U	4.5	25
cis-1,3-Dichloropropene	25	U	2.5	25
trans-1,3-Dichloropropene	25	U	4.0	25
Ethylbenzene	25	U	5.0	25
2-Hexanone	50	U	4.0	50
Methylene Chloride	25	U	2.0	25
4-Methyl-2-pentanone (MIBK)	50	U	3.5	50
Styrene	25	U	2.5	25
1,1,2,2-Tetrachloroethane	25	U	2.0	25
Tetrachloroethene	580		2.5	25
Toluene	25	U	1.5	25
1,1,1-Trichloroethane	25	U	2.0	25
1,1,2-Trichloroethane	25	U	3.0	25
Trichloroethene	25	U	3.5	25
Vinyl chloride	25	U	4.0	25
Xylenes, Total	25	U	5.0	25
cis-1,2-Dichloroethene	25	U	3.0	25
trans-1,2-Dichloroethene	25	U	2.5	25

Surrogate	%Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	75	53 - 125
4-Bromofluorobenzene	119	73 - 127
Dibromofluoromethane	77	54 - 137
Toluene-d8 (Surr)	85	63 - 121

DATA REPORTING QUALIFIERS

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Lab Section	Qualifier	Description
GC/MS VOA		
	U	Analyzed for but not detected.
	J	Indicates an estimated value.
	B	The analyte was found in an associated blank, as well as in the sample.

QUALITY CONTROL RESULTS

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

QC Association Summary

Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
GC/MS VOA					
Analysis Batch:220-8307					
LCS 220-8307/2	Lab Control Spike	T	Water	8260B	
MB 220-8307/3	Method Blank	T	Water	8260B	
220-2277-1	A-2	T	Water	8260B	
220-2277-3	A-7	T	Water	8260B	
220-2277-10FB	FB0723	T	Water	8260B	
220-2277-11TB	TRIP BLANK	T	Water	8260B	
Analysis Batch:220-8321					
LCS 220-8321/2	Lab Control Spike	T	Water	8260B	
MB 220-8321/3	Method Blank	T	Water	8260B	
220-2277-13	A-17	T	Water	8260B	
220-2277-14	MW-109	T	Water	8260B	
220-2277-15	MW-110	T	Water	8260B	
220-2277-16	A-18	T	Water	8260B	
220-2277-19	A-9	T	Water	8260B	
220-2277-20	A-10	T	Water	8260B	
220-2277-21	MW-113	T	Water	8260B	
220-2277-24	ME-13	T	Water	8260B	
220-2277-25FB	FB0726	T	Water	8260B	
220-2277-26	A-9D	T	Water	8260B	
220-2277-27	A-5	T	Water	8260B	
Analysis Batch:220-8356					
LCS 220-8356/2	Lab Control Spike	T	Water	8260B	
MSB 220-8356/24	Matrix Spike Blank	T	Water	8260B	
MB 220-8356/4	Method Blank	T	Water	8260B	
220-2277-2	A-6	T	Water	8260B	
220-2277-4	A-8	T	Water	8260B	
220-2277-5	A-3	T	Water	8260B	
220-2277-6	A-11	T	Water	8260B	
220-2277-7	A-12	T	Water	8260B	
220-2277-8	ME-7	T	Water	8260B	
220-2277-9	ME-11	T	Water	8260B	
220-2277-12	A-16	T	Water	8260B	
220-2277-17	A-14	T	Water	8260B	
220-2277-18	A-15	T	Water	8260B	
220-2277-22	MW-112	T	Water	8260B	
220-2277-22MS	Matrix Spike	T	Water	8260B	
220-2277-22MSD	Matrix Spike Duplicate	T	Water	8260B	
220-2277-23	ME-8	T	Water	8260B	
220-2277-28	ME-15	T	Water	8260B	

Report Basis
T = Total

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Surrogate Recovery Report

8260B Volatile Organic Compounds by GC/MS

Client Matrix: Water

Lab Sample ID	Client Sample ID	(BFB) (%Rec)	(DCE) (%Rec)	(DFM) (%Rec)	(TOL) (%Rec)
220-2277-1	A-2	118	74	75	82
220-2277-2	A-6	120	76	78	83
220-2277-3	A-7	112	74	75	83
220-2277-4	A-8	117	74	75	84
220-2277-5	A-3	112	71	72	78
220-2277-6	A-11	116	73	76	81
220-2277-7	A-12	112	75	72	81
220-2277-8	ME-7	116	77	75	82
220-2277-9	ME-11	111	73	75	80
220-2277-10	FB0723	110	71	75	80
220-2277-11	TRIP BLANK	121	77	74	86
220-2277-12	A-16	115	76	78	84
220-2277-13	A-17	107	72	74	83
220-2277-14	MW-109	114	71	72	81
220-2277-15	MW-110	112	75	78	85
220-2277-16	A-18	107	71	73	78
220-2277-17	A-14	115	76	76	83
220-2277-18	A-15	113	75	75	80

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Surrogate Recovery Report

8260B Volatile Organic Compounds by GC/MS

Client Matrix: Water

		(BFB) (%Rec)	(DCE) (%Rec)	(DFM) (%Rec)	(TOL) (%Rec)
220-2277-19	A-9	110	72	73	80
220-2277-20	A-10	109	73	74	80
220-2277-21	MW-113	110	73	75	81
220-2277-22	MW-112	115	75	77	84
220-2277-23	ME-8	108	72	74	80
220-2277-24	ME-13	120	80	80	85
220-2277-25	FB0726	113	75	74	84
220-2277-26	A-9D	119	72	77	83
220-2277-27	A-5	107	71	75	80
220-2277-28	ME-15	119	75	77	85
220-2277-22 MS	MW-112	94	73	76	79
220-2277-22 MSD	MW-112	94	73	75	79
LCS 220-8307/2		101	74	76	86
LCS 220-8321/2		94	73	73	80
LCS 220-8356/2		97	74	76	80
MB 220-8307/3		108	76	77	83
MB 220-8321/3		113	74	76	82
MB 220-8356/4		109	74	73	80

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Surrogate Recovery Report

8260B Volatile Organic Compounds by GC/MS

Client Matrix: Water

	(BFB) (%Rec)	(DCE) (%Rec)	(DFM) (%Rec)	(TOL) (%Rec)
MSB 220-8356/24	95	75	77	81

Surrogate		Acceptance Limits
(BFB)	4-Bromofluorobenzene	73 - 127
(DCE)	1,2-Dichloroethane-d4 (Surr)	53 - 125
(DFM)	Dibromofluoromethane	54 - 137
(TOL)	Toluene-d8 (Surr)	63 - 121

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Method Blank - Batch: 220-8307

Method: 8260B
Preparation: 5030B

Lab Sample ID: MB 220-8307/3
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 07/31/2007 2136
Date Prepared: 07/31/2007 2136

Analysis Batch: 220-8307
Prep Batch: N/A
Units: ug/L

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9313.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result	Qual	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	5.0	U	0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	% Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	76	53 - 125
4-Bromofluorobenzene	108	73 - 127
Dibromofluoromethane	77	54 - 137
Toluene-d8 (Surr)	83	63 - 121

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

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Lab Control Spike - Batch: 220-8307

Method: 8260B
Preparation: 5030B

Lab Sample ID: LCS 220-8307/2
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 07/31/2007 2046
Date Prepared: 07/31/2007 2046

Analysis Batch: 220-8307
Prep Batch: N/A
Units: ug/L

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9311.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Spike Amount	Result	% Rec.	Limit	Qual
Acetone	20.0	32.5	163	18 - 263	
Benzene	20.0	20.0	100	68 - 126	
Bromodichloromethane	20.0	19.6	98	67 - 118	
Bromoform	20.0	17.8	89	63 - 115	
Bromomethane	20.0	22.3	112	27 - 171	
2-Butanone (MEK)	20.0	24.9	124	30 - 222	
Carbon disulfide	20.0	17.8	89	44 - 142	
Carbon tetrachloride	20.0	21.7	108	56 - 131	
Chlorobenzene	20.0	20.9	105	71 - 114	
Chloroethane	20.0	27.3	137	53 - 167	
Chloroform	20.0	21.1	106	70 - 124	
Chloromethane	20.0	22.5	112	43 - 134	
Dibromochloromethane	20.0	18.4	92	65 - 114	
1,1-Dichloroethane	20.0	20.7	103	67 - 121	
1,2-Dichloroethane	20.0	20.7	103	68 - 124	
1,1-Dichloroethene	20.0	21.7	109	57 - 137	
1,2-Dichloropropane	20.0	20.9	104	69 - 122	
cis-1,3-Dichloropropene	20.0	19.1	95	60 - 122	
trans-1,3-Dichloropropene	20.0	19.2	96	55 - 126	
Ethylbenzene	20.0	21.0	105	71 - 115	
2-Hexanone	20.0	22.2	111	54 - 179	
Methylene Chloride	20.0	19.8	99	61 - 129	
4-Methyl-2-pentanone (MIBK)	20.0	19.7	99	61 - 140	
Styrene	20.0	18.6	93	69 - 112	
1,1,2,2-Tetrachloroethane	20.0	20.1	101	66 - 129	
Tetrachloroethene	20.0	20.4	102	62 - 118	
Toluene	20.0	20.7	103	70 - 116	
1,1,1-Trichloroethane	20.0	20.1	100	60 - 128	
1,1,2-Trichloroethane	20.0	20.0	100	70 - 119	
Trichloroethene	20.0	21.0	105	58 - 125	
Vinyl chloride	20.0	22.3	111	51 - 139	
Xylenes, Total	60.0	61.9	103	66 - 118	
cis-1,2-Dichloroethene	20.0	20.9	105	65 - 120	
trans-1,2-Dichloroethene	20.0	20.3	101	57 - 129	

Surrogate	% Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	74	53 - 125
4-Bromofluorobenzene	101	73 - 127
Dibromofluoromethane	76	54 - 137
Toluene-d8 (Surr)	86	63 - 121

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

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Method Blank - Batch: 220-8321

Method: 8260B
Preparation: 5030B

Lab Sample ID: MB 220-8321/3
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/01/2007 1141
Date Prepared: 08/01/2007 1141

Analysis Batch: 220-8321
Prep Batch: N/A
Units: ug/L

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9344.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result	Qual	MDL	RL
Acetone	3.1	J	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	0.56	J	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	5.0	U	0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	% Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	74	53 - 125
4-Bromofluorobenzene	113	73 - 127
Dibromofluoromethane	76	54 - 137
Toluene-d8 (Surr)	82	63 - 121

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

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Lab Control Spike - Batch: 220-8321

Method: 8260B
Preparation: 5030B

Lab Sample ID: LCS 220-8321/2
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/01/2007 1034
Date Prepared: 08/01/2007 1034

Analysis Batch: 220-8321
Prep Batch: N/A
Units: ug/L

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9342.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Spike Amount	Result	% Rec.	Limit	Qual
Acetone	20.0	33.9	170	18 - 263	B
Benzene	20.0	20.3	101	68 - 126	
Bromodichloromethane	20.0	19.3	96	67 - 118	
Bromoform	20.0	18.1	91	63 - 115	
Bromomethane	20.0	29.0	145	27 - 171	
2-Butanone (MEK)	20.0	29.7	148	30 - 222	
Carbon disulfide	20.0	18.6	93	44 - 142	
Carbon tetrachloride	20.0	21.5	107	56 - 131	
Chlorobenzene	20.0	20.2	101	71 - 114	
Chloroethane	20.0	28.7	143	53 - 167	
Chloroform	20.0	20.3	102	70 - 124	
Chloromethane	20.0	24.7	124	43 - 134	
Dibromochloromethane	20.0	18.9	94	65 - 114	
1,1-Dichloroethane	20.0	20.7	104	67 - 121	
1,2-Dichloroethane	20.0	21.1	105	68 - 124	
1,1-Dichloroethene	20.0	22.9	115	57 - 137	
1,2-Dichloropropane	20.0	21.6	108	69 - 122	
cis-1,3-Dichloropropene	20.0	19.5	98	60 - 122	
trans-1,3-Dichloropropene	20.0	19.1	96	55 - 126	
Ethylbenzene	20.0	20.0	100	71 - 115	
2-Hexanone	20.0	25.0	125	54 - 179	
Methylene Chloride	20.0	19.3	96	61 - 129	B
4-Methyl-2-pentanone (MIBK)	20.0	21.8	109	61 - 140	
Styrene	20.0	18.7	94	69 - 112	
1,1,2,2-Tetrachloroethane	20.0	20.8	104	66 - 129	
Tetrachloroethene	20.0	21.0	105	62 - 118	
Toluene	20.0	20.0	100	70 - 116	
1,1,1-Trichloroethane	20.0	20.7	103	60 - 128	
1,1,2-Trichloroethane	20.0	20.9	104	70 - 119	
Trichloroethene	20.0	20.7	104	58 - 125	
Vinyl chloride	20.0	24.5	123	51 - 139	
Xylenes, Total	60.0	60.6	101	66 - 118	
cis-1,2-Dichloroethene	20.0	20.6	103	65 - 120	
trans-1,2-Dichloroethene	20.0	20.8	104	57 - 129	

Surrogate	% Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	73	53 - 125
4-Bromofluorobenzene	94	73 - 127
Dibromofluoromethane	73	54 - 137
Toluene-d8 (Surr)	80	63 - 121

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

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Method Blank - Batch: 220-8356

Method: 8260B
Preparation: 5030B

Lab Sample ID: MB 220-8356/4
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/02/2007 1130
Date Prepared: 08/02/2007 1130

Analysis Batch: 220-8356
Prep Batch: N/A
Units: ug/L

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9375.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result	Qual	MDL	RL
Acetone	10	U	1.4	10
Benzene	5.0	U	0.40	5.0
Bromodichloromethane	5.0	U	0.40	5.0
Bromoform	5.0	U	0.80	5.0
Bromomethane	5.0	U	1.2	5.0
2-Butanone (MEK)	10	U	1.2	10
Carbon disulfide	5.0	U	0.90	5.0
Carbon tetrachloride	5.0	U	1.0	5.0
Chlorobenzene	5.0	U	0.40	5.0
Chloroethane	5.0	U	0.80	5.0
Chloroform	5.0	U	0.70	5.0
Chloromethane	5.0	U	0.50	5.0
Dibromochloromethane	5.0	U	0.50	5.0
1,1-Dichloroethane	5.0	U	0.60	5.0
1,2-Dichloroethane	5.0	U	0.60	5.0
1,1-Dichloroethene	5.0	U	0.70	5.0
1,2-Dichloropropane	5.0	U	0.90	5.0
cis-1,3-Dichloropropene	5.0	U	0.50	5.0
trans-1,3-Dichloropropene	5.0	U	0.80	5.0
Ethylbenzene	5.0	U	1.0	5.0
2-Hexanone	10	U	0.80	10
Methylene Chloride	5.0	U	0.40	5.0
4-Methyl-2-pentanone (MIBK)	10	U	0.70	10
Styrene	5.0	U	0.50	5.0
1,1,2,2-Tetrachloroethane	5.0	U	0.40	5.0
Tetrachloroethene	5.0	U	0.50	5.0
Toluene	5.0	U	0.30	5.0
1,1,1-Trichloroethane	5.0	U	0.40	5.0
1,1,2-Trichloroethane	5.0	U	0.60	5.0
Trichloroethene	5.0	U	0.70	5.0
Vinyl chloride	5.0	U	0.80	5.0
Xylenes, Total	5.0	U	1.0	5.0
cis-1,2-Dichloroethene	5.0	U	0.60	5.0
trans-1,2-Dichloroethene	5.0	U	0.50	5.0

Surrogate	% Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	74	53 - 125
4-Bromofluorobenzene	109	73 - 127
Dibromofluoromethane	73	54 - 137
Toluene-d8 (Surr)	80	63 - 121

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

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Lab Control Spike - Batch: 220-8356

Method: 8260B
Preparation: 5030B

Lab Sample ID: LCS 220-8356/2
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/02/2007 0959
Date Prepared: 08/02/2007 0959

Analysis Batch: 220-8356
Prep Batch: N/A
Units: ug/L

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9372.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Spike Amount	Result	% Rec.	Limit	Qual
Acetone	20.0	33.8	169	18 - 263	
Benzene	20.0	20.2	101	68 - 126	
Bromodichloromethane	20.0	19.1	95	67 - 118	
Bromoform	20.0	17.6	88	63 - 115	
Bromomethane	20.0	24.2	121	27 - 171	
2-Butanone (MEK)	20.0	27.9	139	30 - 222	
Carbon disulfide	20.0	19.0	95	44 - 142	
Carbon tetrachloride	20.0	21.6	108	56 - 131	
Chlorobenzene	20.0	19.7	99	71 - 114	
Chloroethane	20.0	26.9	134	53 - 167	
Chloroform	20.0	20.1	101	70 - 124	
Chloromethane	20.0	24.3	121	43 - 134	
Dibromochloromethane	20.0	18.6	93	65 - 114	
1,1-Dichloroethane	20.0	20.8	104	67 - 121	
1,2-Dichloroethane	20.0	20.9	105	68 - 124	
1,1-Dichloroethene	20.0	22.7	114	57 - 137	
1,2-Dichloropropane	20.0	20.6	103	69 - 122	
cis-1,3-Dichloropropene	20.0	19.3	97	60 - 122	
trans-1,3-Dichloropropene	20.0	19.8	99	55 - 126	
Ethylbenzene	20.0	20.4	102	71 - 115	
2-Hexanone	20.0	24.0	120	54 - 179	
Methylene Chloride	20.0	20.2	101	61 - 129	
4-Methyl-2-pentanone (MIBK)	20.0	20.3	101	61 - 140	
Styrene	20.0	18.2	91	69 - 112	
1,1,2,2-Tetrachloroethane	20.0	21.1	106	66 - 129	
Tetrachloroethene	20.0	20.2	101	62 - 118	
Toluene	20.0	19.4	97	70 - 116	
1,1,1-Trichloroethane	20.0	20.5	102	60 - 128	
1,1,2-Trichloroethane	20.0	20.1	101	70 - 119	
Trichloroethene	20.0	20.0	100	58 - 125	
Vinyl chloride	20.0	23.6	118	51 - 139	
Xylenes, Total	60.0	60.7	101	66 - 118	
cis-1,2-Dichloroethene	20.0	19.9	99	65 - 120	
trans-1,2-Dichloroethene	20.0	20.2	101	57 - 129	

Surrogate	% Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	74	53 - 125
4-Bromofluorobenzene	97	73 - 127
Dibromofluoromethane	76	54 - 137
Toluene-d8 (Surr)	80	63 - 121

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

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Matrix Spike Blank - Batch: 220-8356

Method: 8260B
Preparation: 5030B

Lab Sample ID: MSB 220-8356/24
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/02/2007 2032
Date Prepared: 08/02/2007 2032

Analysis Batch: 220-8356
Prep Batch: N/A
Units: ug/L

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9395.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Spike Amount	Result	% Rec.	Limit	Qual
Acetone	50.0	48.9	98	18 - 263	
Benzene	50.0	53.2	106	68 - 126	
Bromodichloromethane	50.0	52.0	104	67 - 118	
Bromoform	50.0	47.0	94	63 - 115	
Bromomethane	50.0	53.7	107	27 - 171	
2-Butanone (MEK)	50.0	54.0	108	30 - 222	
Carbon disulfide	50.0	50.3	101	44 - 142	
Carbon tetrachloride	50.0	57.7	115	56 - 131	
Chlorobenzene	50.0	53.5	107	71 - 114	
Chloroethane	50.0	64.7	129	53 - 167	
Chloroform	50.0	54.8	110	70 - 124	
Chloromethane	50.0	50.7	101	43 - 134	
Dibromochloromethane	50.0	48.1	96	65 - 114	
1,1-Dichloroethane	50.0	55.0	110	67 - 121	
1,2-Dichloroethane	50.0	54.2	108	68 - 124	
1,1-Dichloroethene	50.0	56.8	114	57 - 137	
1,2-Dichloropropane	50.0	55.5	111	69 - 122	
cis-1,3-Dichloropropene	50.0	49.5	99	60 - 122	
trans-1,3-Dichloropropene	50.0	49.7	99	55 - 126	
Ethylbenzene	50.0	52.8	106	71 - 115	
2-Hexanone	50.0	55.3	111	54 - 179	
Methylene Chloride	50.0	53.9	108	61 - 129	
4-Methyl-2-pentanone (MIBK)	50.0	56.3	113	61 - 140	
Styrene	50.0	54.2	108	69 - 112	
1,1,2,2-Tetrachloroethane	50.0	52.2	104	66 - 129	
Tetrachloroethene	50.0	54.6	109	62 - 118	
Toluene	50.0	52.8	106	70 - 116	
1,1,1-Trichloroethane	50.0	53.9	108	60 - 128	
1,1,2-Trichloroethane	50.0	55.0	110	70 - 119	
Trichloroethene	50.0	53.4	107	58 - 125	
Vinyl chloride	50.0	53.7	107	51 - 139	
Xylenes, Total	150	162	108	66 - 118	
cis-1,2-Dichloroethene	50.0	55.1	110	65 - 120	
trans-1,2-Dichloroethene	50.0	56.0	112	57 - 129	

Surrogate	% Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	75	53 - 125
4-Bromofluorobenzene	95	73 - 127
Dibromofluoromethane	77	54 - 137
Toluene-d8 (Surr)	81	63 - 121

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

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Matrix Spike/
Matrix Spike Duplicate Recovery Report - Batch: 220-8356

Method: 8260B
Preparation: 5030B

MS Lab Sample ID: 220-2277-22
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/02/2007 1942
Date Prepared: 08/02/2007 1942

Analysis Batch: 220-8356
Prep Batch: N/A

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9393.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

MSD Lab Sample ID: 220-2277-22
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/02/2007 2007
Date Prepared: 08/02/2007 2007

Analysis Batch: 220-8356
Prep Batch: N/A

Instrument ID: HP 5890/5971 GC/MS
Lab File ID: L9394.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	% Rec. MS	% Rec. MSD	Limit	RPD	RPD Limit	MS Qual	MSD Qual
Acetone	80	89	18 - 263	11	20		
Benzene	99	102	68 - 126	3	20		
Bromodichloromethane	96	99	67 - 118	3	20		
Bromoform	90	95	63 - 115	6	20		
Bromomethane	83	94	27 - 171	12	20		
2-Butanone (MEK)	101	99	30 - 222	2	20		
Carbon disulfide	90	94	44 - 142	4	20		
Carbon tetrachloride	92	96	56 - 131	4	20		
Chlorobenzene	99	103	71 - 114	4	20		
Chloroethane	122	121	53 - 167	1	20		
Chloroform	98	101	70 - 124	3	20		
Chloromethane	92	96	43 - 134	4	20		
Dibromochloromethane	89	94	65 - 114	5	20		
1,1-Dichloroethane	104	103	67 - 121	1	20		
1,2-Dichloroethane	99	104	68 - 124	5	20		
1,1-Dichloroethene	104	105	57 - 137	1	20		
1,2-Dichloropropane	101	103	69 - 122	2	20		
cis-1,3-Dichloropropene	94	96	60 - 122	3	20		
trans-1,3-Dichloropropene	94	96	55 - 126	3	20		
Ethylbenzene	99	105	71 - 115	6	20		
2-Hexanone	98	103	54 - 179	6	20		
Methylene Chloride	99	101	61 - 129	3	20		
4-Methyl-2-pentanone (MIBK)	101	109	61 - 140	7	20		
Styrene	97	103	69 - 112	7	20		
1,1,2,2-Tetrachloroethane	98	106	66 - 129	7	20		
Tetrachloroethene	97	110	62 - 118	12	20		
Toluene	99	103	70 - 116	3	20		
1,1,1-Trichloroethane	100	103	60 - 128	2	20		
1,1,2-Trichloroethane	104	104	70 - 119	0	20		
Trichloroethene	98	101	58 - 125	3	20		

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

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Matrix Spike/
Matrix Spike Duplicate Recovery Report - Batch: 220-8356

Method: 8260B
Preparation: 5030B

MS Lab Sample ID:	220-2277-22	Analysis Batch: 220-8356	Instrument ID: HP 5890/5971 GC/MS
Client Matrix:	Water	Prep Batch: N/A	Lab File ID: L9393.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	08/02/2007 1942		Final Weight/Volume: 5 mL
Date Prepared:	08/02/2007 1942		

MSD Lab Sample ID:	220-2277-22	Analysis Batch: 220-8356	Instrument ID: HP 5890/5971 GC/MS
Client Matrix:	Water	Prep Batch: N/A	Lab File ID: L9394.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	08/02/2007 2007		Final Weight/Volume: 5 mL
Date Prepared:	08/02/2007 2007		

	% Rec.						
Analyte	MS	MSD	Limit	RPD	RPD Limit	MS Qual	MSD Qual
Vinyl chloride	100	100	51 - 139	0	20		
Xylenes, Total	99	104	66 - 118	5	20		
cis-1,2-Dichloroethene	100	105	65 - 120	5	20		
trans-1,2-Dichloroethene	99	102	57 - 129	3	20		

Surrogate	MS % Rec	MSD % Rec	Acceptance Limits
1,2-Dichloroethane-d4 (Surr)	73	73	53 - 125
4-Bromofluorobenzene	94	94	73 - 127
Dibromofluoromethane	76	75	54 - 137
Toluene-d8 (Surr)	79	79	63 - 121

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

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Matrix Spike/
Matrix Spike Duplicate Data Report - Batch: 220-8356

Method: 8260B
Preparation: 5030B

MS Lab Sample ID: 220-2277-22
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/02/2007 1942
Date Prepared: 08/02/2007 1942

Units: ug/L

MSD Lab Sample ID: 220-2277-22
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/02/2007 2007
Date Prepared: 08/02/2007 2007

Analyte	Sample Result/Qual		MS Spike Amount	MSD Spike Amount	MS Result/Qual	MSD Result/Qual
Acetone	10	U	50.0	50.0	39.8	44.4
Benzene	5.0	U	50.0	50.0	49.4	50.8
Bromodichloromethane	5.0	U	50.0	50.0	47.8	49.3
Bromoform	5.0	U	50.0	50.0	44.8	47.5
Bromomethane	5.0	U	50.0	50.0	41.7	47.2
2-Butanone (MEK)	10	U	50.0	50.0	50.3	49.4
Carbon disulfide	5.0	U	50.0	50.0	45.2	47.2
Carbon tetrachloride	5.0	U	50.0	50.0	46.1	47.8
Chlorobenzene	5.0	U	50.0	50.0	49.3	51.3
Chloroethane	5.0	U	50.0	50.0	60.9	60.3
Chloroform	5.0	U	50.0	50.0	49.1	50.7
Chloromethane	5.0	U	50.0	50.0	46.2	48.2
Dibromochloromethane	5.0	U	50.0	50.0	44.7	47.2
1,1-Dichloroethane	5.0	U	50.0	50.0	52.0	51.6
1,2-Dichloroethane	5.0	U	50.0	50.0	49.3	52.1
1,1-Dichloroethene	5.0	U	50.0	50.0	52.2	52.6
1,2-Dichloropropane	5.0	U	50.0	50.0	50.6	51.6
cis-1,3-Dichloropropene	5.0	U	50.0	50.0	46.9	48.2
trans-1,3-Dichloropropene	5.0	U	50.0	50.0	46.9	48.1
Ethylbenzene	5.0	U	50.0	50.0	49.4	52.4
2-Hexanone	10	U	50.0	50.0	48.9	51.7
Methylene Chloride	5.0	U	50.0	50.0	49.3	50.6
4-Methyl-2-pentanone (MIBK)	10	U	50.0	50.0	50.5	54.4
Styrene	5.0	U	50.0	50.0	48.3	51.7
1,1,2,2-Tetrachloroethane	5.0	U	50.0	50.0	49.2	52.8
Tetrachloroethene	1.3	J	50.0	50.0	49.9	56.1
Toluene	5.0	U	50.0	50.0	49.7	51.3
1,1,1-Trichloroethane	5.0	U	50.0	50.0	50.2	51.4
1,1,2-Trichloroethane	5.0	U	50.0	50.0	51.8	52.1
Trichloroethene	5.0	U	50.0	50.0	49.0	50.4
Vinyl chloride	5.0	U	50.0	50.0	49.8	49.9
Xylenes, Total	5.0	U	150	150	149	156
cis-1,2-Dichloroethene	5.0	U	50.0	50.0	49.8	52.3
trans-1,2-Dichloroethene	5.0	U	50.0	50.0	49.7	51.2

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

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Laboratory Chronicle

Client Samples:

Lab ID: 220-2277-1 **Client ID:** A-2

Sample Date/Time: 07/23/2007 0900 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-B-1	1	220-8307		08/01/2007 0348	1.0	STL CT	DH
P-5030B		1			08/01/2007 0348	1.0	STL CT	DH

Lab ID: 220-2277-2 **Client ID:** A-6

Sample Date/Time: 07/23/2007 0930 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-A-2	1	220-8356		08/02/2007 1648	2.0	STL CT	DH
P-5030B		1			08/02/2007 1648	2.0	STL CT	DH

Lab ID: 220-2277-3 **Client ID:** A-7

Sample Date/Time: 07/23/2007 1000 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-A-3	1	220-8307		08/01/2007 0438	1.0	STL CT	DH
P-5030B		1			08/01/2007 0438	1.0	STL CT	DH

Lab ID: 220-2277-4 **Client ID:** A-8

Sample Date/Time: 07/23/2007 1030 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-B-4	1	220-8356		08/02/2007 1508	50	STL CT	DH
P-5030B		1			08/02/2007 1508	50	STL CT	DH

Lab ID: 220-2277-5 **Client ID:** A-3

Sample Date/Time: 07/23/2007 1100 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-A-5	1	220-8356		08/02/2007 1232	1.0	STL CT	DH
P-5030B		1			08/02/2007 1232	1.0	STL CT	DH

A = Analytical Method P = Prep Method

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Laboratory Chronicle

Client Samples:

Lab ID: 220-2277-6 **Client ID: A-11**

Sample Date/Time: 07/23/2007 1130 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-B-6	1	220-8356		08/02/2007 1257	1.0	STL CT	DH
P-5030B		1			08/02/2007 1257	1.0	STL CT	DH

Lab ID: 220-2277-7 **Client ID: A-12**

Sample Date/Time: 07/23/2007 1230 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-B-7	1	220-8356		08/02/2007 1533	20	STL CT	DH
P-5030B		1			08/02/2007 1533	20	STL CT	DH

Lab ID: 220-2277-8 **Client ID: ME-7**

Sample Date/Time: 07/23/2007 1300 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-B-8	1	220-8356		08/02/2007 1322	1.0	STL CT	DH
P-5030B		1			08/02/2007 1322	1.0	STL CT	DH

Lab ID: 220-2277-9 **Client ID: ME-11**

Sample Date/Time: 07/23/2007 1330 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-A-9	1	220-8356		08/02/2007 1713	2.0	STL CT	DH
P-5030B		1			08/02/2007 1713	2.0	STL CT	DH

Lab ID: 220-2277-10 **Client ID: FB0723**

Sample Date/Time: 07/23/2007 1400 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-B-10	1	220-8307		08/01/2007 0259	1.0	STL CT	DH
P-5030B		1			08/01/2007 0259	1.0	STL CT	DH

A = Analytical Method P = Prep Method

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Laboratory Chronicle

Client Samples:

Lab ID: 220-2277-11 **Client ID: TRIP BLANK**

Sample Date/Time: 07/23/2007 0000 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-B-11	1	220-8307		08/01/2007 0323	1.0	STL CT	DH
P-5030B		1			08/01/2007 0323	1.0	STL CT	DH

Lab ID: 220-2277-12 **Client ID: A-16**

Sample Date/Time: 07/26/2007 0800 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-A-12	1	220-8356		08/02/2007 1347	1.0	STL CT	DH
P-5030B		1			08/02/2007 1347	1.0	STL CT	DH

Lab ID: 220-2277-13 **Client ID: A-17**

Sample Date/Time: 07/26/2007 0830 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-A-13	1	220-8321		08/01/2007 1445	1.0	STL CT	DH
P-5030B		1			08/01/2007 1445	1.0	STL CT	DH

Lab ID: 220-2277-14 **Client ID: MW-109**

Sample Date/Time: 07/26/2007 0900 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-A-14	1	220-8321		08/01/2007 1510	1.0	STL CT	DH
P-5030B		1			08/01/2007 1510	1.0	STL CT	DH

Lab ID: 220-2277-15 **Client ID: MW-110**

Sample Date/Time: 07/26/2007 0930 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-B-15	1	220-8321		08/01/2007 1535	1.0	STL CT	DH
P-5030B		1			08/01/2007 1535	1.0	STL CT	DH

A = Analytical Method P = Prep Method

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Laboratory Chronicle

Client Samples:

Lab ID: 220-2277-16 **Client ID:** A-18

Sample Date/Time: 07/26/2007 1000 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-A-16	1	220-8321		08/01/2007 1600	1.0	STL CT	DH
P-5030B		1			08/01/2007 1600	1.0	STL CT	DH

Lab ID: 220-2277-17 **Client ID:** A-14

Sample Date/Time: 07/26/2007 1030 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-B-17	1	220-8356		08/02/2007 1623	4.0	STL CT	DH
P-5030B		1			08/02/2007 1623	4.0	STL CT	DH

Lab ID: 220-2277-18 **Client ID:** A-15

Sample Date/Time: 07/26/2007 1100 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-B-18	1	220-8356		08/02/2007 1412	1.0	STL CT	DH
P-5030B		1			08/02/2007 1412	1.0	STL CT	DH

Lab ID: 220-2277-19 **Client ID:** A-9

Sample Date/Time: 07/26/2007 1130 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-A-19	1	220-8321		08/01/2007 1715	1.0	STL CT	DH
P-5030B		1			08/01/2007 1715	1.0	STL CT	DH

Lab ID: 220-2277-20 **Client ID:** A-10

Sample Date/Time: 07/26/2007 1200 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-A-20	1	220-8321		08/01/2007 1740	1.0	STL CT	DH
P-5030B		1			08/01/2007 1740	1.0	STL CT	DH

A = Analytical Method P = Prep Method

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Laboratory Chronicle

Client Samples:

Lab ID: 220-2277-21 **Client ID: MW-113**

Sample Date/Time: 07/26/2007 1300 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-A-21	1	220-8321		08/01/2007 1804	1.0	STL CT	DH
P-5030B		1			08/01/2007 1804	1.0	STL CT	DH

Lab ID: 220-2277-22 **Client ID: MW-112**

Sample Date/Time: 07/26/2007 1330 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-B-22	1	220-8356		08/02/2007 1436	1.0	STL CT	DH
P-5030B		1			08/02/2007 1436	1.0	STL CT	DH

Lab ID: 220-2277-22MS **Client ID: MW-112**

Sample Date/Time: 07/26/2007 1330 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-C-22 MS	1	220-8356		08/02/2007 1942	1.0	STL CT	DH
P-5030B		1			08/02/2007 1942	1.0	STL CT	DH

Lab ID: 220-2277-22MSD **Client ID: MW-112**

Sample Date/Time: 07/26/2007 1330 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-C-22 MSD	1	220-8356		08/02/2007 2007	1.0	STL CT	DH
P-5030B		1			08/02/2007 2007	1.0	STL CT	DH

Lab ID: 220-2277-23 **Client ID: ME-8**

Sample Date/Time: 07/26/2007 1400 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-B-23	1	220-8356		08/02/2007 1737	2.0	STL CT	DH
P-5030B		1			08/02/2007 1737	2.0	STL CT	DH

A = Analytical Method P = Prep Method

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Laboratory Chronicle

Client Samples:

Lab ID: 220-2277-24 **Client ID: ME-13**

Sample Date/Time: 07/26/2007 1430 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-A-24	1	220-8321		08/01/2007 1919	1.0	STL CT	DH
P-5030B		1			08/01/2007 1919	1.0	STL CT	DH

Lab ID: 220-2277-25 **Client ID: FB0726**

Sample Date/Time: 07/26/2007 0000 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-A-25	1	220-8321		08/01/2007 1944	1.0	STL CT	DH
P-5030B		1			08/01/2007 1944	1.0	STL CT	DH

Lab ID: 220-2277-26 **Client ID: A-9D**

Sample Date/Time: 07/26/2007 1145 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-A-26	1	220-8321		08/01/2007 2008	1.0	STL CT	DH
P-5030B		1			08/01/2007 2008	1.0	STL CT	DH

Lab ID: 220-2277-27 **Client ID: A-5**

Sample Date/Time: 07/23/2007 1045 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-A-27	1	220-8321		08/01/2007 2033	1.0	STL CT	DH
P-5030B		1			08/01/2007 2033	1.0	STL CT	DH

Lab ID: 220-2277-28 **Client ID: ME-15**

Sample Date/Time: 07/23/2007 1200 Received Date/Time: 07/28/2007 1000

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B	220-2277-B-28	1	220-8356		08/02/2007 1558	5.0	STL CT	DH
P-5030B		1			08/02/2007 1558	5.0	STL CT	DH

A = Analytical Method P = Prep Method

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Laboratory Chronicle

Client Samples:

Lab ID: MB **Client ID:** MB

Sample Date/Time: NA Received Date/Time: NA

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B		1	220-8307		07/31/2007 2136	1.0	STL CT	DH
P-5030B		1			07/31/2007 2136	1.0	STL CT	DH
A-8260B		1	220-8321		08/01/2007 1141	1.0	STL CT	DH
P-5030B		1			08/01/2007 1141	1.0	STL CT	DH
A-8260B		1	220-8356		08/02/2007 1130	1.0	STL CT	DH
P-5030B		1			08/02/2007 1130	1.0	STL CT	DH

Lab ID: LCS **Client ID:** LCS

Sample Date/Time: NA Received Date/Time: NA

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B		1	220-8307		07/31/2007 2046	1.0	STL CT	DH
P-5030B		1			07/31/2007 2046	1.0	STL CT	DH
A-8260B		1	220-8321		08/01/2007 1034	1.0	STL CT	DH
P-5030B		1			08/01/2007 1034	1.0	STL CT	DH
A-8260B		1	220-8356		08/02/2007 0959	1.0	STL CT	DH
P-5030B		1			08/02/2007 0959	1.0	STL CT	DH

Lab ID: MSB **Client ID:** MSB

Sample Date/Time: NA Received Date/Time: NA

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
A-8260B		1	220-8356		08/02/2007 2032	1.0	STL CT	DH
P-5030B		1			08/02/2007 2032	1.0	STL CT	DH

A = Analytical Method P = Prep Method

MISCELLANEOUS DOCUMENTS

Chain of Custody Record

-2277

SEVERN TRENT **STL**
Severn Trent Laboratories, Inc.

STL-4124 (0901)

Client: FPM	Project Manager: Bob Cancemi	Date: 7/26/07	Chain of Custody Number: 184462
Address: 1000 Koukoma 909 Marconi Ave	Telephone Number (Area Code)/Fax Number: 631-737-6200 x2410	Lab Number:	Page 2 of 3
City: Ronkonkoma	State: NY	Zip Code: 11779	Site Contact: Bob C	Lab Contact: Johanna D.
Project Name and Location (State): Metropolitan Ave / NY	Carrier/Waybill Number:		
Contract/Purchase Order/Quote No.			

Sample I.D. No. and Description (Containers for each sample may be combined on one line)	Date	Time	Matrix: Air	Aqueous	Sed.	Soil	Containers & Preservatives: Unpres.	H2SO4	HNO3	HCl	NaOH	ZnAc/NaOH	Analysis (Attach list if more space is needed): CAS VI	Special Instructions/ Conditions of Receipt
11 TRIP BLANK	7/23/07			X						2				
12 A-16	7/25/07	800												
13 A-17		830												
14 MW-109		900												
15 MW-110		930												
16 A-18		1000												
17 A-14		1035												
18 A-15		1100												
19 A-9		1130												
20 A-10		1200												
21 MW-113		1300												
22 MW-112		1330								6				MS/MSD

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Possible Hazard Identification
☒ Non-Hazard ☐ Flammable ☐ Skin Irritant ☐ Poison B ☐ Unknown

Sample Disposal
☐ Return To Client ☒ Disposal By Lab ☐ Archive For _____ Months (A fee may be assessed if samples are retained longer than 1 month)

Turn Around Time Required
☐ 24 Hours ☐ 48 Hours ☐ 7 Days ☐ 14 Days ☐ 21 Days ☐ Other STD

QC Requirements (Specify): NYSDEC ASP CAT B

	Date	Time		Date	Time
1. Relinquished By:	7/27/07	1700	1. Received By:	7/28/07	10:00
2. Relinquished By:			2. Received By:		
3. Relinquished By:			3. Received By:		

Comments: Will call / Hold Samples 0.8°C PASSED RAD SCREEN

DISTRIBUTION: WHITE - Returned to Client with Report; CANARY - Stays with the Sample; PINK - Field Copy

08/09/2007

Chain of Custody Record

STL-4124 (0605)

STL Connecticut
128 Long Hill Cross Road
Shelton, CT 06484
Tel: 203-929-8140
Fax: 203-929-8142

-2277

SEVERN TRENT STL
Severn Trent Laboratories, Inc.

158

Client: FPM	Project Manager: Bob Cancemi	Date: 7/23/07	Chain of Custody Number: 10411
Address: 909 Marconi Ave.	Telephone Number (Area Code)/Fax Number: 631-737-2410/200	Lab Number:	Page 1 of 3
City: Ronkonkoma	State: NY	Zip Code: 11779	Site Contact: Bob C
Project Name and Location (State): Metropolitan Ave (NY)	Rad. Screen:	Lab Contact: Johanna D	Analysis (Attach list if more space is needed): 8260
Contract/Purchase Order/Project No.			

Sample I.D. and Location/Description (Containers for each sample may be combined on one line)	Date	Time	Lab ID	Matrix: Aqueous	Solid	Other	Containers & Preservatives: Unpres.	H2SO4	HNO3	HCl	NaOH	ZnAc/NaOH	Comments
1 A-2	7/23/07	900		X						2			
2 A6		930											
3 A-7		1000											
4 A-8		1030											
5 A-3		1100											
A-11		1130											
A-12		1230											
Me 7		1300											
Me 11		1330											
FB0723	↓	1400								↓			

Possible Hazard Identification
☒ Non-Hazard ☐ Flammable ☐ Skin Irritant ☐ Poison B ☐ Unknown

Sample Disposal ☒ Disposal By Lab ☐ Return To Client ☐ Archive For _____ Months (A fee may be assessed if samples are retained longer than 1 month)

Turn Around Time Required (business days)
☐ 24 Hours ☐ 48 Hours ☐ 5 Days ☐ 10 Days ☐ 15 Days ☐ Other STD

QC Requirements (Specify): NYSDEC ASP CAT B.

1. Relinquished By: Ron Hu	Date: 7/27/07	Time: 1700	1. Received By: Tracy R	Date: 7/28/07	Time: 10:00
2. Relinquished By:	Date:	Time:	2. Received By:	Date:	Time:
3. Received By:	Date:	Time:	Cooler Temps: 0.8°C		

Comments: Hold Samples

DISTRIBUTION: WHITE – Stays with the Samples; CANARY – Returned to Client with Report; PINK – Field Copy

"PASSED RAD SCREEN"

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08/09/2007

Chain of Custody Record

-2277

SEVERN TRENT **STL**

Severn Trent Laboratories, Inc.

STL-4124 (0901)

Client: FPM	Project Manager: Ben Carcemi	Date: 7/26/07	Chain of Custody Number: 343924
Address: 909 Marconi Ave	Telephone Number (Area Code)/Fax Number: 631-737-6200/2410	Lab Number:	Page 3 of 3
City: Ronkonkoma / State: NY / Zip Code: 11779	Site Contact:	Lab Contact:	
Project Name and Location (State): Metropolitan Ave (NY)	Carrier/Waybill Number:		
Contract/Purchase Order/Quote No.			

Sample I.D. No. and Description (Containers for each sample may be combined on one line)	Date	Time	Matrix: Air	Aqueous	Sed.	Soil	Containers & Preservatives: Unpres.	H2SO4	HNO3	HCl	NaOH	ZnAc/ NaOH	Analysis (Attach list if more space is needed): VOA's	Special Instructions/ Conditions of Receipt
23 ME-8	7/25/07	1400		X						2			X BC	
24 ME-13		1430												
25 FB0726														
26 A-9D	7/25/07	1145												
27 A-5	7/23/07													
28 Me-5	7/23/07													

Added by Sample Control 7/3

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Possible Hazard Identification: [X] Non-Hazard [ ] Flammable [ ] Skin Irritant [ ] Poison B [ ] Unknown

Sample Disposal: [ ] Return To Client [X] Disposal By Lab [ ] Archive For _____ Months (A fee may be assessed if samples are retained longer than 1 month)

Turn Around Time Required: [ ] 24 Hours [ ] 48 Hours [ ] 7 Days [ ] 14 Days [ ] 21 Days [ ] Other STD

QC Requirements (Specify): NYSDEC ASP CAT B

	Date	Time		Date	Time
1. Relinquished By: [illegible]	7/27/07	1200	1. Received By: [signature]	7/28/07	10:00
2. Relinquished By			2. Received By		
3. Relinquished By			3. Received By		

Comments: +10 Samples

0.8°C

"PASSED RAD SCREEN"

DISTRIBUTION: WHITE - Returned to Client with Report; CANARY - Stays with the Sample; PINK - Field Copy

08/09/2007

STL - Connecticut
Internal Chain-of-Custody

220- 2277
FPM metropolitan aveNY

Trip Blank: # 11

QC: ✓ Air: ✓

FB: #10,25

Soil: ✓ Water: 1-28

Date Received: 7/28/07

Sample #s: 1-28

Locations: REF 35L

Laboratory Sample #	Relinquished by	Accepted by	Date	Time	Reason	Relinquished by	Accepted by	Date	Time
1-12	UB	D.C.	7/31	1800	VOA				
13-28	UB	D.C.	8/1/07	13:50	VOA	Used			
2,4,5,6-9,	12,17-18,22-23,28	D.C.	8/2/07	12:00	VOA	Used			

STL Form# SMF00507.CT

LOGIN SAMPLE RECEIPT CHECK LIST

Client: FPM Group Limited

Job Number: 220-2277-1
Sdg Number: 220-2277

Login Number: 2277

Question	T/F/NA	Comment
Radioactivity either was not measured or, if measured, is at or below background	True	
The cooler's custody seal, if present, is intact.	True	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
There are no discrepancies between the sample IDs on the containers and the COC.	True	
Samples are received within Holding Time.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
VOA sample vials do not have headspace or bubble is <6mm (1/4") in diameter.	True	
If necessary, staff have been informed of any short hold time or quick TAT needs	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	

VOLATILE DATA

FORM II
GC/MS VOA SURROGATE RECOVERY

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Matrix: Water Level: Low

GC Column (1): RTX-VMS ID: 0.18(mm)

Lab Sample ID	Client Sample ID	12DCE #	BFB #	DBFM #	TOL #
220-2277-1	A-2	74	118	75	82
220-2277-2	A-6	76	120	78	83
220-2277-3	A-7	74	112	75	83
220-2277-4	A-8	74	117	75	84
220-2277-5	A-3	71	112	72	78
220-2277-6	A-11	73	116	76	81
220-2277-7	A-12	75	112	72	81
220-2277-8	ME-7	77	116	75	82
220-2277-9	ME-11	73	111	75	80
220-2277-10	FB0723	71	110	75	80
220-2277-11	TRIP BLANK	77	121	74	86
220-2277-12	A-16	76	115	78	84
220-2277-13	A-17	72	107	74	83
220-2277-14	MW-109	71	114	72	81
220-2277-15	MW-110	75	112	78	85
220-2277-16	A-18	71	107	73	78
220-2277-17	A-14	76	115	76	83
220-2277-18	A-15	75	113	75	80
220-2277-19	A-9	72	110	73	80
220-2277-20	A-10	73	109	74	80
220-2277-21	MW-113	73	110	75	81
220-2277-22	MW-112	75	115	77	84
220-2277-23	ME-8	72	108	74	80
220-2277-24	ME-13	80	120	80	85
220-2277-25	FB0726	75	113	74	84
220-2277-26	A-9D	72	119	77	83
220-2277-27	A-5	71	107	75	80
220-2277-28	ME-15	75	119	77	85
MB 220-8307/3		76	108	77	83
MB 220-8321/3		74	113	76	82
MB 220-8356/4		74	109	73	80
LCS 220-8307/2		74	101	76	86
LCS 220-8321/2		73	94	73	80
LCS 220-8356/2		74	97	76	80
MSB 220-8356/24		75	95	77	81
220-2277-22 MS	MW-112	73	94	76	79
220-2277-22 MSD	MW-112	73	94	75	79

Column to be used to flag recovery values

FORM II 8260B

FORM II
GC/MS VOA SURROGATE RECOVERY

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Matrix: Water Level: Low

GC Column (1): RTX-VMS ID: 0.18(mm)

Lab Sample ID	Client Sample ID	12DCE #	BFB #	DBFM #	TOL #

	QC LIMITS
12DCE = 1,2-Dichloroethane-d4 (Surr)	53 = 125
BFB = 4-Bromofluorobenzene	73 = 127
DBFM = Dibromofluoromethane	54 = 137
TOL = Toluene-d8 (Surr)	63 = 121

Column to be used to flag recovery values

FORM II 8260B

LAB CONTROL SPIKE
VOLATILE ORGANIC COMPOUNDS BY GC/MS

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Matrix: Water Level: Low Lab File ID: L9311.D

Lab ID: LCS 220-8307/2 Client ID:

COMPOUND	SPIKE ADDED (ug/L)	LCS CONCENTRATION (ug/L)	LCS % REC	QC LIMITS REC	#
Acetone	20.0	32.5	163	18-263	
Benzene	20.0	20.0	100	68-126	
Bromodichloromethane	20.0	19.6	98	67-118	
Bromoform	20.0	17.8	89	63-115	
Bromomethane	20.0	22.3	112	27-171	
2-Butanone (MEK)	20.0	24.9	124	30-222	
Carbon disulfide	20.0	17.8	89	44-142	
Carbon tetrachloride	20.0	21.7	108	56-131	
Chlorobenzene	20.0	20.9	105	71-114	
Chloroethane	20.0	27.3	137	53-167	
Chloroform	20.0	21.1	106	70-124	
Chloromethane	20.0	22.5	112	43-134	
Dibromochloromethane	20.0	18.4	92	65-114	
1,1-Dichloroethane	20.0	20.7	103	67-121	
1,2-Dichloroethane	20.0	20.7	103	68-124	
1,1-Dichloroethene	20.0	21.7	109	57-137	
1,2-Dichloropropane	20.0	20.9	104	69-122	
cis-1,3-Dichloropropene	20.0	19.1	95	60-122	
trans-1,3-Dichloropropene	20.0	19.2	96	55-126	
Ethylbenzene	20.0	21.0	105	71-115	
2-Hexanone	20.0	22.2	111	54-179	
Methylene Chloride	20.0	19.8	99	61-129	
4-Methyl-2-pentanone (MIBK)	20.0	19.7	99	61-140	
Styrene	20.0	18.6	93	69-112	
1,1,2,2-Tetrachloroethane	20.0	20.1	101	66-129	
Tetrachloroethene	20.0	20.4	102	62-118	
Toluene	20.0	20.7	103	70-116	
1,1,1-Trichloroethane	20.0	20.1	100	60-128	
1,1,2-Trichloroethane	20.0	20.0	100	70-119	
Trichloroethene	20.0	21.0	105	58-125	
Vinyl chloride	20.0	22.3	111	51-139	
Xylenes, Total	60.0	61.9	103	66-118	
cis-1,2-Dichloroethene	20.0	20.9	105	65-120	
trans-1,2-Dichloroethene	20.0	20.3	101	57-129	

Calculations are performed before rounding

Column to be used to flag recovery and RPD values

LAB CONTROL SPIKE
VOLATILE ORGANIC COMPOUNDS BY GC/MS

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Matrix: Water Level: Low Lab File ID: L9342.D

Lab ID: LCS 220-8321/2 Client ID:

COMPOUND	SPIKE ADDED (ug/L)	LCS CONCENTRATION (ug/L)	LCS % REC	QC LIMITS REC	#
Acetone	20.0	33.9	170	18-263	
Benzene	20.0	20.3	101	68-126	
Bromodichloromethane	20.0	19.3	96	67-118	
Bromoform	20.0	18.1	91	63-115	
Bromomethane	20.0	29.0	145	27-171	
2-Butanone (MEK)	20.0	29.7	148	30-222	
Carbon disulfide	20.0	18.6	93	44-142	
Carbon tetrachloride	20.0	21.5	107	56-131	
Chlorobenzene	20.0	20.2	101	71-114	
Chloroethane	20.0	28.7	143	53-167	
Chloroform	20.0	20.3	102	70-124	
Chloromethane	20.0	24.7	124	43-134	
Dibromochloromethane	20.0	18.9	94	65-114	
1,1-Dichloroethane	20.0	20.7	104	67-121	
1,2-Dichloroethane	20.0	21.1	105	68-124	
1,1-Dichloroethene	20.0	22.9	115	57-137	
1,2-Dichloropropane	20.0	21.6	108	69-122	
cis-1,3-Dichloropropene	20.0	19.5	98	60-122	
trans-1,3-Dichloropropene	20.0	19.1	96	55-126	
Ethylbenzene	20.0	20.0	100	71-115	
2-Hexanone	20.0	25.0	125	54-179	
Methylene Chloride	20.0	19.3	96	61-129	
4-Methyl-2-pentanone (MIBK)	20.0	21.8	109	61-140	
Styrene	20.0	18.7	94	69-112	
1,1,2,2-Tetrachloroethane	20.0	20.8	104	66-129	
Tetrachloroethene	20.0	21.0	105	62-118	
Toluene	20.0	20.0	100	70-116	
1,1,1-Trichloroethane	20.0	20.7	103	60-128	
1,1,2-Trichloroethane	20.0	20.9	104	70-119	
Trichloroethene	20.0	20.7	104	58-125	
Vinyl chloride	20.0	24.5	123	51-139	
Xylenes, Total	60.0	60.6	101	66-118	
cis-1,2-Dichloroethene	20.0	20.6	103	65-120	
trans-1,2-Dichloroethene	20.0	20.8	104	57-129	

Calculations are performed before rounding

Column to be used to flag recovery and RPD values

LAB CONTROL SPIKE
VOLATILE ORGANIC COMPOUNDS BY GC/MS

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Matrix: Water Level: Low Lab File ID: L9372.D

Lab ID: LCS 220-8356/2 Client ID:

COMPOUND	SPIKE ADDED (ug/L)	LCS CONCENTRATION (ug/L)	LCS % REC	QC LIMITS REC	#
Acetone	20.0	33.8	169	18-263	
Benzene	20.0	20.2	101	68-126	
Bromodichloromethane	20.0	19.1	95	67-118	
Bromoform	20.0	17.6	88	63-115	
Bromomethane	20.0	24.2	121	27-171	
2-Butanone (MEK)	20.0	27.9	139	30-222	
Carbon disulfide	20.0	19.0	95	44-142	
Carbon tetrachloride	20.0	21.6	108	56-131	
Chlorobenzene	20.0	19.7	99	71-114	
Chloroethane	20.0	26.9	134	53-167	
Chloroform	20.0	20.1	101	70-124	
Chloromethane	20.0	24.3	121	43-134	
Dibromochloromethane	20.0	18.6	93	65-114	
1,1-Dichloroethane	20.0	20.8	104	67-121	
1,2-Dichloroethane	20.0	20.9	105	68-124	
1,1-Dichloroethene	20.0	22.7	114	57-137	
1,2-Dichloropropane	20.0	20.6	103	69-122	
cis-1,3-Dichloropropene	20.0	19.3	97	60-122	
trans-1,3-Dichloropropene	20.0	19.8	99	55-126	
Ethylbenzene	20.0	20.4	102	71-115	
2-Hexanone	20.0	24.0	120	54-179	
Methylene Chloride	20.0	20.2	101	61-129	
4-Methyl-2-pentanone (MIBK)	20.0	20.3	101	61-140	
Styrene	20.0	18.2	91	69-112	
1,1,2,2-Tetrachloroethane	20.0	21.1	106	66-129	
Tetrachloroethene	20.0	20.2	101	62-118	
Toluene	20.0	19.4	97	70-116	
1,1,1-Trichloroethane	20.0	20.5	102	60-128	
1,1,2-Trichloroethane	20.0	20.1	101	70-119	
Trichloroethene	20.0	20.0	100	58-125	
Vinyl chloride	20.0	23.6	118	51-139	
Xylenes, Total	60.0	60.7	101	66-118	
cis-1,2-Dichloroethene	20.0	19.9	99	65-120	
trans-1,2-Dichloroethene	20.0	20.2	101	57-129	

Calculations are performed before rounding

Column to be used to flag recovery and RPD values

MATRIX SPIKE
VOLATILE ORGANIC COMPOUNDS BY GC/MS

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Matrix: Water Level: Low Lab File ID: L9393.D

Lab ID: 220-2277-22 MS Client ID: MW-112

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE CONCENTRATION (ug/L)		MS CONCENTRATION (ug/L)	MS % REC	QC LIMITS REC	#
Acetone	50.0	10	U	39.8	80	18-263	
Benzene	50.0	5.0	U	49.4	99	68-126	
Bromodichloromethane	50.0	5.0	U	47.8	96	67-118	
Bromoform	50.0	5.0	U	44.8	90	63-115	
Bromomethane	50.0	5.0	U	41.7	83	27-171	
2-Butanone (MEK)	50.0	10	U	50.3	101	30-222	
Carbon disulfide	50.0	5.0	U	45.2	90	44-142	
Carbon tetrachloride	50.0	5.0	U	46.1	92	56-131	
Chlorobenzene	50.0	5.0	U	49.3	99	71-114	
Chloroethane	50.0	5.0	U	60.9	122	53-167	
Chloroform	50.0	5.0	U	49.1	98	70-124	
Chloromethane	50.0	5.0	U	46.2	92	43-134	
Dibromochloromethane	50.0	5.0	U	44.7	89	65-114	
1,1-Dichloroethane	50.0	5.0	U	52.0	104	67-121	
1,2-Dichloroethane	50.0	5.0	U	49.3	99	68-124	
1,1-Dichloroethene	50.0	5.0	U	52.2	104	57-137	
1,2-Dichloropropane	50.0	5.0	U	50.6	101	69-122	
cis-1,3-Dichloropropene	50.0	5.0	U	46.9	94	60-122	
trans-1,3-Dichloropropene	50.0	5.0	U	46.9	94	55-126	
Ethylbenzene	50.0	5.0	U	49.4	99	71-115	
2-Hexanone	50.0	10	U	48.9	98	54-179	
Methylene Chloride	50.0	5.0	U	49.3	99	61-129	
4-Methyl-2-pentanone (MIBK)	50.0	10	U	50.5	101	61-140	
Styrene	50.0	5.0	U	48.3	97	69-112	
1,1,2,2-Tetrachloroethane	50.0	5.0	U	49.2	98	66-129	
Tetrachloroethene	50.0	1.3	J	49.9	97	62-118	
Toluene	50.0	5.0	U	49.7	99	70-116	
1,1,1-Trichloroethane	50.0	5.0	U	50.2	100	60-128	
1,1,2-Trichloroethane	50.0	5.0	U	51.8	104	70-119	
Trichloroethene	50.0	5.0	U	49.0	98	58-125	
Vinyl chloride	50.0	5.0	U	49.8	100	51-139	
Xylenes, Total	150	5.0	U	149	99	66-118	
cis-1,2-Dichloroethene	50.0	5.0	U	49.8	100	65-120	
trans-1,2-Dichloroethene	50.0	5.0	U	49.7	99	57-129	

Calculations are performed before rounding

Column to be used to flag recovery and RPD values

MATRIX SPIKE BLANK
VOLATILE ORGANIC COMPOUNDS BY GC/MS

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Matrix: Water Level: Low Lab File ID: L9395.D

Lab ID: MSB 220-8356/24 Client ID:

COMPOUND	SPIKE ADDED (ug/L)	MSB CONCENTRATION (ug/L)	MSB % REC	QC LIMITS REC	#
Acetone	50.0	48.9	98	18-263	
Benzene	50.0	53.2	106	68-126	
Bromodichloromethane	50.0	52.0	104	67-118	
Bromoform	50.0	47.0	94	63-115	
Bromomethane	50.0	53.7	107	27-171	
2-Butanone (MEK)	50.0	54.0	108	30-222	
Carbon disulfide	50.0	50.3	101	44-142	
Carbon tetrachloride	50.0	57.7	115	56-131	
Chlorobenzene	50.0	53.5	107	71-114	
Chloroethane	50.0	64.7	129	53-167	
Chloroform	50.0	54.8	110	70-124	
Chloromethane	50.0	50.7	101	43-134	
Dibromochloromethane	50.0	48.1	96	65-114	
1,1-Dichloroethane	50.0	55.0	110	67-121	
1,2-Dichloroethane	50.0	54.2	108	68-124	
1,1-Dichloroethene	50.0	56.8	114	57-137	
1,2-Dichloropropane	50.0	55.5	111	69-122	
cis-1,3-Dichloropropene	50.0	49.5	99	60-122	
trans-1,3-Dichloropropene	50.0	49.7	99	55-126	
Ethylbenzene	50.0	52.8	106	71-115	
2-Hexanone	50.0	55.3	111	54-179	
Methylene Chloride	50.0	53.9	108	61-129	
4-Methyl-2-pentanone (MIBK)	50.0	56.3	113	61-140	
Styrene	50.0	54.2	108	69-112	
1,1,2,2-Tetrachloroethane	50.0	52.2	104	66-129	
Tetrachloroethene	50.0	54.6	109	62-118	
Toluene	50.0	52.8	106	70-116	
1,1,1-Trichloroethane	50.0	53.9	108	60-128	
1,1,2-Trichloroethane	50.0	55.0	110	70-119	
Trichloroethene	50.0	53.4	107	58-125	
Vinyl chloride	50.0	53.7	107	51-139	
Xylenes, Total	150	162	108	66-118	
cis-1,2-Dichloroethene	50.0	55.1	110	65-120	
trans-1,2-Dichloroethene	50.0	56.0	112	57-129	

Calculations are performed before rounding

Column to be used to flag recovery and RPD values

MATRIX SPIKE DUPLICATE
VOLATILE ORGANIC COMPOUNDS BY GC/MS

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Matrix: Water Level: Low Lab File ID: L9394.D

Lab ID: 220-2277-22 MSD Client ID: MW-112

COMPOUND	SPIKE ADDED (ug/L)	MSD CONCENTRATION (ug/L)	MSD % REC	% RPD	QC LIMITS RPD	QC LIMITS REC	#
Acetone	50.0	44.4	89	11	20	18-263	
Benzene	50.0	50.8	102	3	20	68-126	
Bromodichloromethane	50.0	49.3	99	3	20	67-118	
Bromoform	50.0	47.5	95	6	20	63-115	
Bromomethane	50.0	47.2	94	12	20	27-171	
2-Butanone (MEK)	50.0	49.4	99	2	20	30-222	
Carbon disulfide	50.0	47.2	94	4	20	44-142	
Carbon tetrachloride	50.0	47.8	96	4	20	56-131	
Chlorobenzene	50.0	51.3	103	4	20	71-114	
Chloroethane	50.0	60.3	121	1	20	53-167	
Chloroform	50.0	50.7	101	3	20	70-124	
Chloromethane	50.0	48.2	96	4	20	43-134	
Dibromochloromethane	50.0	47.2	94	5	20	65-114	
1,1-Dichloroethane	50.0	51.6	103	1	20	67-121	
1,2-Dichloroethane	50.0	52.1	104	5	20	68-124	
1,1-Dichloroethene	50.0	52.6	105	1	20	57-137	
1,2-Dichloropropane	50.0	51.6	103	2	20	69-122	
cis-1,3-Dichloropropene	50.0	48.2	96	3	20	60-122	
trans-1,3-Dichloropropene	50.0	48.1	96	3	20	55-126	
Ethylbenzene	50.0	52.4	105	6	20	71-115	
2-Hexanone	50.0	51.7	103	6	20	54-179	
Methylene Chloride	50.0	50.6	101	3	20	61-129	
4-Methyl-2-pentanone (MIBK)	50.0	54.4	109	7	20	61-140	
Styrene	50.0	51.7	103	7	20	69-112	
1,1,2,2-Tetrachloroethane	50.0	52.8	106	7	20	66-129	
Tetrachloroethene	50.0	56.1	110	12	20	62-118	
Toluene	50.0	51.3	103	3	20	70-116	
1,1,1-Trichloroethane	50.0	51.4	103	2	20	60-128	
1,1,2-Trichloroethane	50.0	52.1	104	0	20	70-119	
Trichloroethene	50.0	50.4	101	3	20	58-125	
Vinyl chloride	50.0	49.9	100	0	20	51-139	
Xylenes, Total	150	156	104	5	20	66-118	
cis-1,2-Dichloroethene	50.0	52.3	105	5	20	65-120	
trans-1,2-Dichloroethene	50.0	51.2	102	3	20	57-129	

Calculations are performed before rounding

Column to be used to flag recovery and RPD values

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Lab File ID: L9313.D Lab Sample ID: MB 220-8307/3

Instrument ID: MSL Date Analyzed: 07/31/2007 21:36

Matrix: Water Heated Purge:(Y/N) N

GC Column: RTX-VMS ID: 0.18 (mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
	LCS 220-8307/2	L9311.D	07/31/2007 20:46
FB0723	220-2277-10	L9326.D	08/01/2007 02:59
TRIP BLANK	220-2277-11	L9327.D	08/01/2007 03:23
A-2	220-2277-1	L9328.D	08/01/2007 03:48
A-7	220-2277-3	L9330.D	08/01/2007 04:38

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Lab File ID: L9344.D Lab Sample ID: MB 220-8321/3

Instrument ID: MSL Date Analyzed: 08/01/2007 11:41

Matrix: Water Heated Purge:(Y/N) N

GC Column: RTX-VMS ID: 0.18 (mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
	LCS 220-8321/2	L9342.D	08/01/2007 10:34
A-17	220-2277-13	L9351.D	08/01/2007 14:45
MW-109	220-2277-14	L9352.D	08/01/2007 15:10
MW-110	220-2277-15	L9353.D	08/01/2007 15:35
A-18	220-2277-16	L9354.D	08/01/2007 16:00
A-9	220-2277-19	L9357.D	08/01/2007 17:15
A-10	220-2277-20	L9358.D	08/01/2007 17:40
MW-113	220-2277-21	L9359.D	08/01/2007 18:04
ME-13	220-2277-24	L9362.D	08/01/2007 19:19
FB0726	220-2277-25	L9363.D	08/01/2007 19:44
A-9D	220-2277-26	L9364.D	08/01/2007 20:08
A-5	220-2277-27	L9365.D	08/01/2007 20:33

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Lab File ID: L9375.D Lab Sample ID: MB 220-8356/4

Instrument ID: MSL Date Analyzed: 08/02/2007 11:30

Matrix: Water Heated Purge: (Y/N) N

GC Column: RTX-VMS ID: 0.18 (mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
	LCS 220-8356/2	L9372.D	08/02/2007 09:59
A-3	220-2277-5	L9376.D	08/02/2007 12:32
A-11	220-2277-6	L9377.D	08/02/2007 12:57
ME-7	220-2277-8	L9378.D	08/02/2007 13:22
A-16	220-2277-12	L9379.D	08/02/2007 13:47
A-15	220-2277-18	L9380.D	08/02/2007 14:12
MW-112	220-2277-22	L9381.D	08/02/2007 14:36
A-8	220-2277-4	L9382.D	08/02/2007 15:08
A-12	220-2277-7	L9383.D	08/02/2007 15:33
ME-15	220-2277-28	L9384.D	08/02/2007 15:58
A-14	220-2277-17	L9385.D	08/02/2007 16:23
A-6	220-2277-2	L9386.D	08/02/2007 16:48
ME-11	220-2277-9	L9387.D	08/02/2007 17:13
ME-8	220-2277-23	L9388.D	08/02/2007 17:37
MW-112 MS	220-2277-22 MS	L9393.D	08/02/2007 19:42
MW-112 MSD	220-2277-22 MSD	L9394.D	08/02/2007 20:07
	MSB 220-8356/24	L9395.D	08/02/2007 20:32

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VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Lab File ID: LB417.D BFB Injection Date: 07/26/2007

Instrument ID: MSL BFB Injection Time: 13:09

Analy. Batch No.: 8212

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE	
50	15.0 - 40.0 % of mass 95	19.9	
75	30.0 - 60.0 % of m/z 95	54.9	
95	Base Peak, 100% relative abundance	100.0	
96	5.0 - 9.0 % of mass 95	6.2	
173	Less than 2.0 % of mass 174	0.0	(0.0)1
174	Greater than 50.0 % of mass 95	87.4	
175	5.0 - 9.0 % of mass 174	6.2	(7.0)1
176	95.0 - 101.0 % of mass 174	87.9	(100.6)1
177	5.0 - 9.0 % of mass 176	5.5	(6.2)2

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
	IC 220-8212/1	L9152.D	07/26/2007	13:32
	IC 220-8212/2	L9153.D	07/26/2007	13:57
	IC 220-8212/3	L9154.D	07/26/2007	14:22
	IC 220-8212/4	L9155.D	07/26/2007	14:47
	IC 220-8212/5	L9156.D	07/26/2007	15:11

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Lab File ID: LB425.D BFB Injection Date: 07/31/2007

Instrument ID: MSL BFB Injection Time: 20:13

Analy. Batch No.: 8307

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE	
50	15.0 - 40.0 % of mass 95	21.2	
75	30.0 - 60.0 % of m/z 95	55.1	
95	Base Peak, 100% relative abundance	100.0	
96	5.0 - 9.0 % of mass 95	7.3	
173	Less than 2.0 % of mass 174	0.3	(0.3)1
174	Greater than 50.0 % of mass 95	89.3	
175	5.0 - 9.0 % of mass 174	6.8	(7.6)1
176	95.0 - 101.0 % of mass 174	89.3	(99.9)1
177	5.0 - 9.0 % of mass 176	5.7	(6.3)2

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
	CCVIS 220-8307/1	L9310.D	07/31/2007	20:21
	LCS 220-8307/2	L9311.D	07/31/2007	20:46
	MB 220-8307/3	L9313.D	07/31/2007	21:36
FB0723	220-2277-10	L9326.D	08/01/2007	02:59
TRIP BLANK	220-2277-11	L9327.D	08/01/2007	03:23
A-2	220-2277-1	L9328.D	08/01/2007	03:48
A-7	220-2277-3	L9330.D	08/01/2007	04:38

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Lab File ID: LB426.D BFB Injection Date: 08/01/2007

Instrument ID: MSL BFB Injection Time: 09:13

Analy. Batch No.: 8321

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE	
50	15.0 - 40.0 % of mass 95	20.6	
75	30.0 - 60.0 % of m/z 95	53.6	
95	Base Peak, 100% relative abundance	100.0	
96	5.0 - 9.0 % of mass 95	6.5	
173	Less than 2.0 % of mass 174	0.5	(0.6)1
174	Greater than 50.0 % of mass 95	89.8	
175	5.0 - 9.0 % of mass 174	6.7	(7.5)1
176	95.0 - 101.0 % of mass 174	88.8	(98.9)1
177	5.0 - 9.0 % of mass 176	5.8	(6.5)2

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
	CCVIS 220-8321/1	L9341.D	08/01/2007	09:47
	LCS 220-8321/2	L9342.D	08/01/2007	10:34
	MB 220-8321/3	L9344.D	08/01/2007	11:41
A-17	220-2277-13	L9351.D	08/01/2007	14:45
MW-109	220-2277-14	L9352.D	08/01/2007	15:10
MW-110	220-2277-15	L9353.D	08/01/2007	15:35
A-18	220-2277-16	L9354.D	08/01/2007	16:00
A-9	220-2277-19	L9357.D	08/01/2007	17:15
A-10	220-2277-20	L9358.D	08/01/2007	17:40
MW-113	220-2277-21	L9359.D	08/01/2007	18:04
ME-13	220-2277-24	L9362.D	08/01/2007	19:19
FB0726	220-2277-25	L9363.D	08/01/2007	19:44
A-9D	220-2277-26	L9364.D	08/01/2007	20:08
A-5	220-2277-27	L9365.D	08/01/2007	20:33

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VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Lab File ID: LB428.D BFB Injection Date: 08/02/2007

Instrument ID: MSL BFB Injection Time: 08:50

Analy. Batch No.: 8356

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE	
50	15.0 - 40.0 % of mass 95	22.0	
75	30.0 - 60.0 % of m/z 95	56.5	
95	Base Peak, 100% relative abundance	100.0	
96	5.0 - 9.0 % of mass 95	6.7	
173	Less than 2.0 % of mass 174	0.3	(0.3)1
174	Greater than 50.0 % of mass 95	93.4	
175	5.0 - 9.0 % of mass 174	7.1	(7.6)1
176	95.0 - 101.0 % of mass 174	90.5	(96.9)1
177	5.0 - 9.0 % of mass 176	5.7	(6.3)2

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
	CCVIS 220-8356/1	L9371.D	08/02/2007	09:18
	LCS 220-8356/2	L9372.D	08/02/2007	09:59
	MB 220-8356/4	L9375.D	08/02/2007	11:30
A-3	220-2277-5	L9376.D	08/02/2007	12:32
A-11	220-2277-6	L9377.D	08/02/2007	12:57
ME-7	220-2277-8	L9378.D	08/02/2007	13:22
A-16	220-2277-12	L9379.D	08/02/2007	13:47
A-15	220-2277-18	L9380.D	08/02/2007	14:12
MW-112	220-2277-22	L9381.D	08/02/2007	14:36
A-8	220-2277-4	L9382.D	08/02/2007	15:08
A-12	220-2277-7	L9383.D	08/02/2007	15:33
ME-15	220-2277-28	L9384.D	08/02/2007	15:58
A-14	220-2277-17	L9385.D	08/02/2007	16:23
A-6	220-2277-2	L9386.D	08/02/2007	16:48
ME-11	220-2277-9	L9387.D	08/02/2007	17:13
ME-8	220-2277-23	L9388.D	08/02/2007	17:37
MW-112 MS	220-2277-22 MS	L9393.D	08/02/2007	19:42
MW-112 MSD	220-2277-22 MSD	L9394.D	08/02/2007	20:07
	MSB 220-8356/24	L9395.D	08/02/2007	20:32

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

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Sample No.: CCVIS 220-8307/1 Date Analyzed: 07/31/2007

Lab File ID (Standard): L9310.D Time Analyzed: 20:21

Instrument ID: MSL Heated Purge: (Y/N) N

GC Column: RTX-VMS ID: 0.18

	DCB AREA	#	RT	#	CBZ AREA	#	RT	#	FB AREA	#	RT	#
12 HOUR STD	160938		10.01		442233		7.96		439656		4.90	
UPPER LIMIT	321876		10.51		884466		8.46		879312		5.40	
LOWER LIMIT	80469		9.51		221117		7.46		219828		4.40	
Lab Sample ID												
LCS 220-8307/2	153272		10.02		428646		7.96		440213		4.90	
MB 220-8307/3	130327		10.01		433216		7.96		450138		4.90	
220-2277-10	124592		10.02		437455		7.96		451044		4.89	
220-2277-11	117531		10.01		417702		7.96		446792		4.90	
220-2277-1	114343		10.01		422601		7.96		442574		4.90	
220-2277-3	119044		10.02		425701		7.96		451910		4.90	

QC LIMITS

DCB = 1,4-Dichlorobenzene-d4
CBZ = Chlorobenzene-d5
FB = Fluorobenzene

Area Upper Limit = 200% of Internal Standard Area
Area Lower Limit = 50% of Internal Standard Area
RT Upper Limit = +0.5 minutes of Internal Standard Retention Time
RT Lower Limit = -0.5 minutes of Internal Standard Retention Time

Column used to flag values outside QC limits

FORM VIII 8260B

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

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Sample No.: CCVIS 220-8321/1 Date Analyzed: 08/01/2007

Lab File ID (Standard): L9341.D Time Analyzed: 09:47

Instrument ID: MSL Heated Purge: (Y/N) N

GC Column: RTX-VMS ID: 0.18

	DCB AREA	#	RT	#	CBZ AREA	#	RT	#	FB AREA	#	RT	#
12 HOUR STD	164048		10.01		444659		7.97		451284		4.91	
UPPER LIMIT	328096		10.51		889318		8.47		902568		5.41	
LOWER LIMIT	82024		9.51		222330		7.47		225642		4.41	
Lab Sample ID												
LCS 220-8321/2	156570		10.01		442311		7.97		450523		4.91	
MB 220-8321/3	124435		10.01		431702		7.96		449562		4.90	
220-2277-13	126762		10.01		425019		7.96		448065		4.90	
220-2277-14	121626		10.01		422636		7.96		451047		4.90	
220-2277-15	123577		10.01		411260		7.96		432833		4.90	
220-2277-16	124354		10.01		435439		7.96		443177		4.90	
220-2277-19	117011		10.01		420067		7.96		443072		4.90	
220-2277-20	123039		10.01		428130		7.96		444007		4.90	
220-2277-21	122483		10.01		422393		7.96		435324		4.90	
220-2277-24	114741		10.01		413641		7.96		424369		4.90	
220-2277-25	122252		10.01		421068		7.96		445250		4.90	
220-2277-26	119100		10.02		426544		7.96		445152		4.90	
220-2277-27	122697		10.02		426795		7.96		445548		4.89	

QC LIMITS

DCB = 1,4-Dichlorobenzene-d4
CBZ = Chlorobenzene-d5
FB = Fluorobenzene

Area Upper Limit = 200% of Internal Standard Area
Area Lower Limit = 50% of Internal Standard Area
RT Upper Limit = +0.5 minutes of Internal Standard Retention Time
RT Lower Limit = -0.5 minutes of Internal Standard Retention Time

Column used to flag values outside QC limits

FORM VIII 8260B

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

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Sample No.: CCVIS 220-8356/1 Date Analyzed: 08/02/2007

Lab File ID (Standard): L9371.D Time Analyzed: 09:18

Instrument ID: MSL Heated Purge: (Y/N) N

GC Column: RTX-VMS ID: 0.18

	DCB AREA	#	RT	#	CBZ AREA	#	RT	#	FB AREA	#	RT	#
12 HOUR STD	157752		10.02		448655		7.96		436789		4.90	
UPPER LIMIT	315504		10.52		897310		8.46		873578		5.40	
LOWER LIMIT	78876		9.52		224328		7.46		218395		4.40	
Lab Sample ID												
LCS 220-8356/2	154291		10.01		440917		7.96		441851		4.90	
MB 220-8356/4	128132		10.01		427020		7.96		438738		4.90	
220-2277-5	124058		10.02		454599		7.96		462027		4.90	
220-2277-6	122354		10.01		437111		7.96		447349		4.90	
220-2277-8	119729		10.03		434210		7.97		444846		4.90	
220-2277-12	122772		10.01		434522		7.95		452863		4.89	
220-2277-18	119186		10.01		425077		7.95		441024		4.89	
220-2277-22	121255		10.01		435304		7.96		452327		4.90	
220-2277-4	121030		10.02		428736		7.96		452451		4.90	
220-2277-7	124274		10.01		436646		7.96		452541		4.90	
220-2277-28	118200		10.01		425683		7.95		445725		4.89	
220-2277-17	119467		10.01		427654		7.96		447266		4.90	
220-2277-2	112968		10.01		426839		7.95		436691		4.89	
220-2277-9	123236		10.02		441225		7.96		458179		4.90	
220-2277-23	123622		10.02		436237		7.96		440199		4.89	
220-2277-22 MS	160374		10.01		438582		7.95		440538		4.89	
220-2277-22 MSD	164899		10.01		444075		7.95		455584		4.89	
MSB 220-8356/24	165255		10.01		439881		7.96		444303		4.89	

QC LIMITS

DCB = 1,4-Dichlorobenzene-d4
CBZ = Chlorobenzene-d5
FB = Fluorobenzene

Area Upper Limit = 200% of Internal Standard Area
Area Lower Limit = 50% of Internal Standard Area
RT Upper Limit = +0.5 minutes of Internal Standard Retention Time
RT Lower Limit = -0.5 minutes of Internal Standard Retention Time

Column used to flag values outside QC limits
FORM VIII 8260B

DETECTION LIMIT STUDY

Method.....:8260B
Analyst.....: Patty Mercure
Equipment ID.:,MSL
Analysis Date:01/23/2006(grp 1)

Date..:2006-04-05
Units.:ug/L
Batch.:63780
T-Val.:6.965

COMPOUND/ELEMENT/TEST	Matrix	Det. Lim.	Units	Mean	Std. Dev.	Grp
Dichlorodifluoromethane	Water	3.87	ug/L	5.588490	0.555588	
Raw Data: 6.04253 5.75398 4.96896						
Chloromethane	Water	5.14	ug/L	5.104133	0.738675	
Raw Data: 5.93126 4.87095 4.51019						
Vinyl chloride	Water	2.05	ug/L	5.131690	0.294862	
Raw Data: 5.42400 5.13673 4.83434						
Bromomethane	Water	1.50	ug/L	7.013927	0.215605	
Raw Data: 7.10251 6.76814 7.17113						
Chloroethane	Water	3.00	ug/L	5.286273	0.431266	
Raw Data: 5.67010 5.36913 4.81959						
Trichlorofluoromethane	Water	2.18	ug/L	5.878263	0.312544	
Raw Data: 6.08783 6.02793 5.51903						
Dichlorofluoromethane	Water	3.69	ug/L	7.548820	0.530080	
Raw Data: 8.08982 7.52626 7.03038						
Ethyl ether	Water	0.19	ug/L	5.411697	0.026675	
Raw Data: 5.38095 5.42866 5.42548						
1,1 Dichloro-1-Fluoroethane	Water	3.38	ug/L	6.473440	0.484578	
Raw Data: 6.87768 6.60637 5.93627						
Freon 123	Water	11.62	ug/L	8.301683	1.668383	
Raw Data: 7.57270 10.2105 7.12185						
Trichlorotrifluoroethane	Water	2.35	ug/L	4.829840	0.337914	
Raw Data: 5.18168 4.50783 4.80001						
1,1-Dichloroethene	Water	5.25	ug/L	4.871583	0.753620	
Raw Data: 5.63898 4.84323 4.13254						
Carbon disulfide	Water	2.55	ug/L	4.884507	0.366728	
Raw Data: 5.26135 4.86336 4.52881						
Iodomethane	Water	1.81	ug/L	4.863700	0.259787	
Raw Data: 4.88762 5.11070 4.59278						
3-Chloropropene (Allyl Chloride)	Water	3.04	ug/L	4.943467	0.436435	
Raw Data: 5.41606 4.85872 4.55562						
Methylene chloride	Water	4.25	ug/L	5.470433	0.610813	
Raw Data: 5.59237 6.01108 4.80785						
Acetone	Water	5.53	ug/L	6.607970	0.793837	
Raw Data: 7.37668 6.65604 5.79119						
trans-1,2-Dichloroethene	Water	2.92	ug/L	4.988330	0.419735	
Raw Data: 5.29091 5.16493 4.50915						
Methyl-tert-butyl-ether (MTBE)	Water	2.53	ug/L	5.670983	0.362996	
Raw Data: 6.07797 5.55431 5.38067						
Acrolein	Water	51.41	ug/L	29.166333	7.380836	
Raw Data: 37.5998 26.0146 23.8846						

DETECTION LIMIT STUDY

Method.....:8260B
Analyst.....: Patty Mercure
Equipment ID.:,MSL
Analysis Date:01/23/2006(grp 1)

Date..:2006-04-05
Units.:ug/L
Batch.:63780
T-Val.:6.965

COMPOUND/ELEMENT/TEST	Matrix	Det. Lim.	Units	Mean	Std. Dev.	Grp
tert-Butyl alcohol	Water	22.47	ug/L	29.834333	3.225418	
Raw Data: 33.2374 29.4434 26.8222						
Methyl acetate	Water	2.56	ug/L	5.170977	0.367420	
Raw Data: 5.52415 5.19798 4.79080						
Acetonitrile	Water	53.30	ug/L	50.074000	7.652999	
Raw Data: 58.0100 49.4725 42.7395						
Isopropyl ether	Water	3.27	ug/L	5.300517	0.469032	
Raw Data: 5.80792 5.21082 4.88281						
tert-butyl Ethyl ether	Water	3.32	ug/L	5.225603	0.476166	
Raw Data: 5.55866 5.43794 4.68021						
Acrylonitrile	Water	3.23	ug/L	10.153730	0.463778	
Raw Data: 9.79450 10.6773 9.98939						
2-Chloro-1,3-butadiene (chloroprene)	Water	4.17	ug/L	4.685450	0.599101	
Raw Data: 5.34527 4.53554 4.17554						
1,1-Dichloroethane	Water	1.07	ug/L	5.211267	0.153270	
Raw Data: 5.25566 5.33744 5.04070						
Vinyl acetate	Water	0.65	ug/L	5.421377	0.093255	
Raw Data: 5.50599 5.43675 5.32139						
cis-1,2-Dichloroethene	Water	3.50	ug/L	5.102250	0.503194	
Raw Data: 5.64464 4.65060 5.01151						
2,2-Dichloropropane	Water	2.53	ug/L	6.907253	0.363717	
Raw Data: 7.18886 7.03629 6.49661						
Bromochloromethane	Water	3.90	ug/L	5.057470	0.560150	
Raw Data: 5.29452 5.46012 4.41777						
Chloroform	Water	2.38	ug/L	6.252087	0.341768	
Raw Data: 6.60799 6.22180 5.92647						
Ethyl acetate	Water	11.90	ug/L	11.259840	1.708047	
Raw Data: 10.5796 13.2032 9.99672						
Methyl Acrylate	Water	0.60	ug/L	5.580237	0.086861	
Raw Data: 5.48509 5.65529 5.60033						
Tetrahydrofuran	Water	11.32	ug/L	11.324757	1.625828	
Raw Data: 12.7691 11.6412 9.56397						
1,1,1-Trichloroethane	Water	2.75	ug/L	6.123623	0.394997	
Raw Data: 6.57466 5.95682 5.83939						
Carbon tetrachloride	Water	3.32	ug/L	5.959003	0.477038	
Raw Data: 6.35000 6.09952 5.42749						
2-Butanone (MEK)	Water	5.48	ug/L	5.680403	0.787346	
Raw Data: 6.58340 5.13746 5.32035						
1,1-Dichloropropene	Water	2.29	ug/L	5.318347	0.328462	
Raw Data: 5.68144 5.23172 5.04188						

DETECTION LIMIT STUDY

Method.....:8260B
Analyst.....: Patty Mercure
Equipment ID.:,MSL
Analysis Date:01/23/2006(grp 1)

Date..:2006-04-05
Units.:ug/L
Batch.:63780
T-Val.:6.965

COMPOUND/ELEMENT/TEST	Matrix	Det. Lim.	Units	Mean	Std. Dev.	Grp
Cyclohexane	Water	3.85	ug/L	5.168847	0.552456	
Raw Data: 5.79523 4.96025 4.75106						
tert-Amyl methyl ether	Water	2.53	ug/L	5.108273	0.362767	
Raw Data: 5.45201 5.14373 4.72908						
tert-butyl Formate	Water	4.02	ug/L	2.389730	0.577473	
Raw Data: 2.98316 1.82966 2.35637						
1-Chlorobutane	Water	6.33	ug/L	6.174730	0.908970	
Raw Data: 6.97423 5.18606 6.36390						
Propionitrile	Water	25.21	ug/L	52.138167	3.619673	
Raw Data: 56.0092 51.5677 48.8376						
Isobutyl alcohol	Water	70.51	ug/L	52.616067	10.123939	
Raw Data: 64.0088 49.1887 44.6507						
Benzene	Water	3.22	ug/L	4.983610	0.462666	
Raw Data: 5.40603 5.05565 4.48915						
Methacrylonitrile	Water	4.83	ug/L	5.521243	0.694040	
Raw Data: 6.32023 5.06784 5.17566						
1,2-Dichloroethane	Water	2.74	ug/L	6.114607	0.392781	
Raw Data: 6.39724 6.28048 5.66610						
Methyl cyclohexane	Water	5.42	ug/L	5.772187	0.777669	
Raw Data: 6.08295 6.34642 4.88719						
Trichloroethene	Water	2.63	ug/L	5.040137	0.377248	
Raw Data: 5.47539 4.83773 4.80729						
Dibromomethane	Water	1.40	ug/L	5.341493	0.201323	
Raw Data: 5.37153 5.52611 5.12684						
1,2-Dichloropropane	Water	2.12	ug/L	5.313437	0.303913	
Raw Data: 5.55803 5.40907 4.97321						
Bromodichloromethane	Water	3.04	ug/L	6.037853	0.436468	
Raw Data: 6.35947 6.21309 5.54100						
Methylmethacrylate	Water	23.32	ug/L	11.980643	3.348338	
Raw Data: 10.4132 15.8252 9.70353						
1,4-Dioxane	Water	20.74	ug/L	23.857200	2.977648	
Raw Data: 24.3330 20.6703 26.5683						
2-Chloroethylvinylether	Water	3.43	ug/L	5.165803	0.491914	
Raw Data: 5.60364 5.26026 4.63351						
cis-1,3-Dichloropropene	Water	3.57	ug/L	5.291210	0.512315	
Raw Data: 5.75373 5.37936 4.74054						
2-Nitropropane	Water	11.28	ug/L	14.339267	1.619235	
Raw Data: 15.3907 15.1525 12.4746						
Chloroacetonitrile	Water	41.70	ug/L	51.598967	5.986364	
Raw Data: 58.4248 49.1308 47.2413						

DETECTION LIMIT STUDY

Method.....:8260B
Analyst.....: Patty Mercure
Equipment ID.:,MSL
Analysis Date:01/23/2006(grp 1)

Date..:2006-04-05
Units.:ug/L
Batch.:63780
T-Val.:6.965

COMPOUND/ELEMENT/TEST	Matrix	Det. Lim.	Units	Mean	Std. Dev.	Grp
trans-1,3-Dichloropropene	Water	1.40	ug/L	5.437937	0.200443	
Raw Data: 5.65667 5.26304 5.39410						
1,1,2-Trichloroethane	Water	3.31	ug/L	5.434230	0.475564	
Raw Data: 5.95975 5.03351 5.30943						
Toluene	Water	1.65	ug/L	4.856570	0.237479	
Raw Data: 5.03843 4.94338 4.58790						
1,1-Dichloro-2-propanone	Water	5.76	ug/L	27.944233	0.826705	
Raw Data: 28.5710 28.2544 27.0073						
4-Methyl-2-pentanone (MIBK)	Water	2.74	ug/L	5.568487	0.393826	
Raw Data: 5.40095 6.01838 5.28613						
Tetrachloroethene	Water	1.57	ug/L	4.507767	0.225692	
Raw Data: 4.67014 4.60311 4.25005						
Ethylmethacrylate	Water	2.99	ug/L	5.214813	0.428648	
Raw Data: 5.69086 5.09415 4.85943						
Dibromochloromethane	Water	1.29	ug/L	5.148520	0.184821	
Raw Data: 5.03446 5.36176 5.04934						
1,3-Dichloropropane	Water	2.84	ug/L	5.049937	0.407431	
Raw Data: 5.19039 5.36856 4.59086						
1,2-Dibromoethane (EDB)	Water	1.63	ug/L	4.846843	0.233486	
Raw Data: 4.96600 4.99671 4.57782						
2-Hexanone	Water	2.28	ug/L	5.204017	0.327318	
Raw Data: 5.51068 5.24201 4.85936						
1-Chlorohexane	Water	1.21	ug/L	4.972647	0.173342	
Raw Data: 4.84938 4.89771 5.17085						
Chlorobenzene	Water	1.93	ug/L	4.898137	0.276771	
Raw Data: 5.04538 5.07016 4.57887						
1,1,1,2-Tetrachloroethane	Water	0.73	ug/L	5.237460	0.104161	
Raw Data: 5.35658 5.19230 5.16350						
Ethylbenzene	Water	1.66	ug/L	4.525157	0.238810	
Raw Data: 4.37897 4.80074 4.39576						
m&p-Xylenes	Water	2.73	ug/L	9.793037	0.391277	
Raw Data: 10.0943 9.93400 9.35081						
o-Xylene	Water	2.45	ug/L	4.932860	0.351963	
Raw Data: 5.32412 4.83244 4.64202						
Styrene	Water	1.28	ug/L	4.720847	0.183903	
Raw Data: 4.81010 4.84309 4.50935						
Bromoform	Water	0.20	ug/L	5.389153	0.028664	
Raw Data: 5.35639 5.40147 5.40960						
Isopropylbenzene	Water	3.80	ug/L	4.918950	0.545131	
~~Raw Data: 5.10159 5.34931 4.30595~~						

DETECTION LIMIT STUDY

Method.....:8260B
Analyst.....: Patty Mercure
Equipment ID.:,MSL
Analysis Date:01/23/2006(grp 1)

Date..:2006-04-05
Units.:ug/L
Batch.:63780
T-Val.:6.965

COMPOUND/ELEMENT/TEST	Matrix	Det. Lim.	Units	Mean	Std. Dev.	Grp
1,1,2,2-Tetrachloroethane	Water	3.29	ug/L	5.287843	0.472162	
Raw Data: 5.82916 5.07348 4.96089						
Bromobenzene	Water	2.09	ug/L	4.984567	0.299740	
Raw Data: 5.23057 5.07241 4.65072						
1,2,3-Trichloropropane	Water	3.84	ug/L	5.035813	0.550872	
Raw Data: 5.65853 4.83683 4.61208						
trans-1,4-Dichloro-2-butene	Water	5.28	ug/L	11.110000	0.758349	
Raw Data: 11.6824 11.3977 10.2499						
n-Propylbenzene	Water	2.37	ug/L	4.845527	0.339678	
Raw Data: 5.17184 4.87084 4.49390						
2-Chlorotoluene	Water	1.27	ug/L	5.046727	0.182293	
Raw Data: 4.88638 5.24500 5.00880						
4-Chlorotoluene	Water	1.98	ug/L	5.103233	0.284529	
Raw Data: 5.22522 5.30643 4.77805						
1,3,5-Trimethylbenzene	Water	2.33	ug/L	4.773503	0.334229	
Raw Data: 4.98611 4.94614 4.38826						
tert-Butylbenzene	Water	1.16	ug/L	4.552073	0.166758	
Raw Data: 4.60582 4.68533 4.36507						
1,2,4-Trimethylbenzene	Water	3.35	ug/L	4.887550	0.480546	
Raw Data: 5.35037 4.92123 4.39105						
sec-Butylbenzene	Water	1.67	ug/L	4.156090	0.239703	
Raw Data: 4.42647 4.07217 3.96963						
p-Isopropyltoluene	Water	2.16	ug/L	4.314183	0.309625	
Raw Data: 4.42742 4.55125 3.96388						
1,3-Dichlorobenzene	Water	2.00	ug/L	4.633493	0.287405	
Raw Data: 4.71807 4.86912 4.31329						
1,4-Dichlorobenzene	Water	1.25	ug/L	4.797680	0.179630	
Raw Data: 4.79112 4.98050 4.62142						
1,2-Dichlorobenzene	Water	1.93	ug/L	4.832130	0.277512	
Raw Data: 4.78106 5.13163 4.58370						
Benzyl chloride	Water	5.06	ug/L	4.706687	0.725906	
Raw Data: 5.15356 5.09739 3.86911						
n-Butylbenzene	Water	2.12	ug/L	4.565733	0.303890	
Raw Data: 4.84762 4.60577 4.24381						
1,2-Dibromo-3-chloropropane	Water	6.32	ug/L	5.417800	0.907570	
Raw Data: 5.70475 6.14721 4.40144						
Nitrobenzene	Water	15.20	ug/L	41.932467	2.183000	
Raw Data: 43.9728 42.1942 39.6304						
1,2,4-Trichlorobenzene	Water	3.01	ug/L	3.141947	0.432088	
~~Raw Data: 2.65321 3.47322 3.29941~~						

DETECTION LIMIT STUDY

Method.....:8260B
Analyst.....: Patty Mercure
Equipment ID.:,MSL
Analysis Date:01/23/2006(grp 1)

Date..:2006-04-05
Units.:ug/L
Batch.:63780
T-Val.:6.965

COMPOUND/ELEMENT/TEST	Matrix	Det. Lim.	Units	Mean	Std. Dev.	Grp
Hexachlorobutadiene	Water	3.11	ug/L	3.466077	0.446742	
Raw Data: 3.97435 3.28824 3.13564						
Naphthalene	Water	2.20	ug/L	3.868153	0.316285	
Raw Data: 4.20508 3.82174 3.57764						
1,2,3-Trichlorobenzene	Water	1.85	ug/L	2.810147	0.264897	
Raw Data: 2.88777 3.02756 2.51511						
1,2-Dichloroethene (total)	Water	5.20	ug/L	10.090563	0.746442	
Raw Data: 10.9355 9.81553 9.52066						
Xylenes (total)	Water	4.97	ug/L	14.725867	0.713664	
Raw Data: 15.4184 14.7664 13.9928						

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-2 | Lab Sample ID: 220-2277-1

Matrix: Water | Lab File ID: L9328.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/01/2007 03:48

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8307 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	32		5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079310.b\L9328.D
Lab Smp Id: 220-2277-B-1 Client Smp ID: A-2
Inj Date : 01-AUG-2007 03:48 MS Autotune Date: 26-APR-2004 14:21
Operator : D. GAYDA Inst ID: msl.i
Smp Info : 220-2277-B-1
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079310.b\L8260BNW.m
Meth Date : 31-Jul-2007 21:07 target Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 19
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

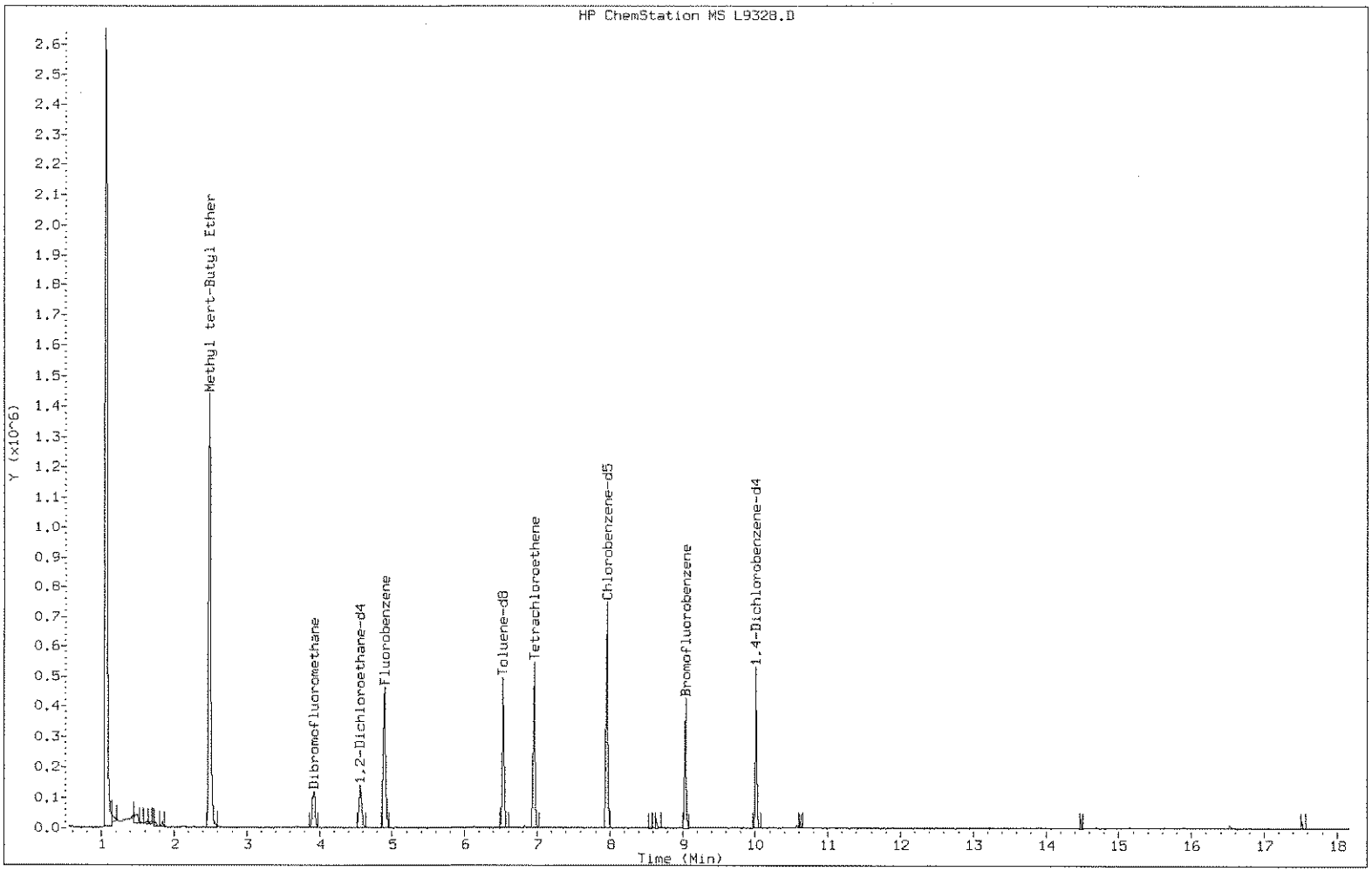
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.898	4.896	(1.000)	442574	25.0000	
	24 Methyl tert-Butyl Ether	73	2.487	2.495	(0.508)	1487367	80.9471	81
$	41 Dibromofluoromethane	111	3.924	3.922	(0.801)	107355	18.8024	19
$	55 1,2-Dichloroethane-d4	65	4.563	4.561	(0.932)	129018	18.5341	18
*	75 Chlorobenzene-d5	117	7.958	7.956	(1.000)	422601	25.0000	
$	77 Toluene-d8	98	6.531	6.529	(0.821)	308126	20.5053	20
	80 Tetrachloroethene	164	6.954	6.952	(0.874)	114486	32.2434	32
*	95 1,4-Dichlorobenzene-d4	152	10.014	10.012	(1.000)	114343	25.0000	
$	125 Bromofluorobenzene	95	9.040	9.038	(0.903)	126324	29.4486	29

Data File: L9328.D

Date: 01-AUG-2007 03:48

Client ID: A-2 Instrument: msl.i

Sample Info: 220-2277-B-1 Operator: D. GAYDA



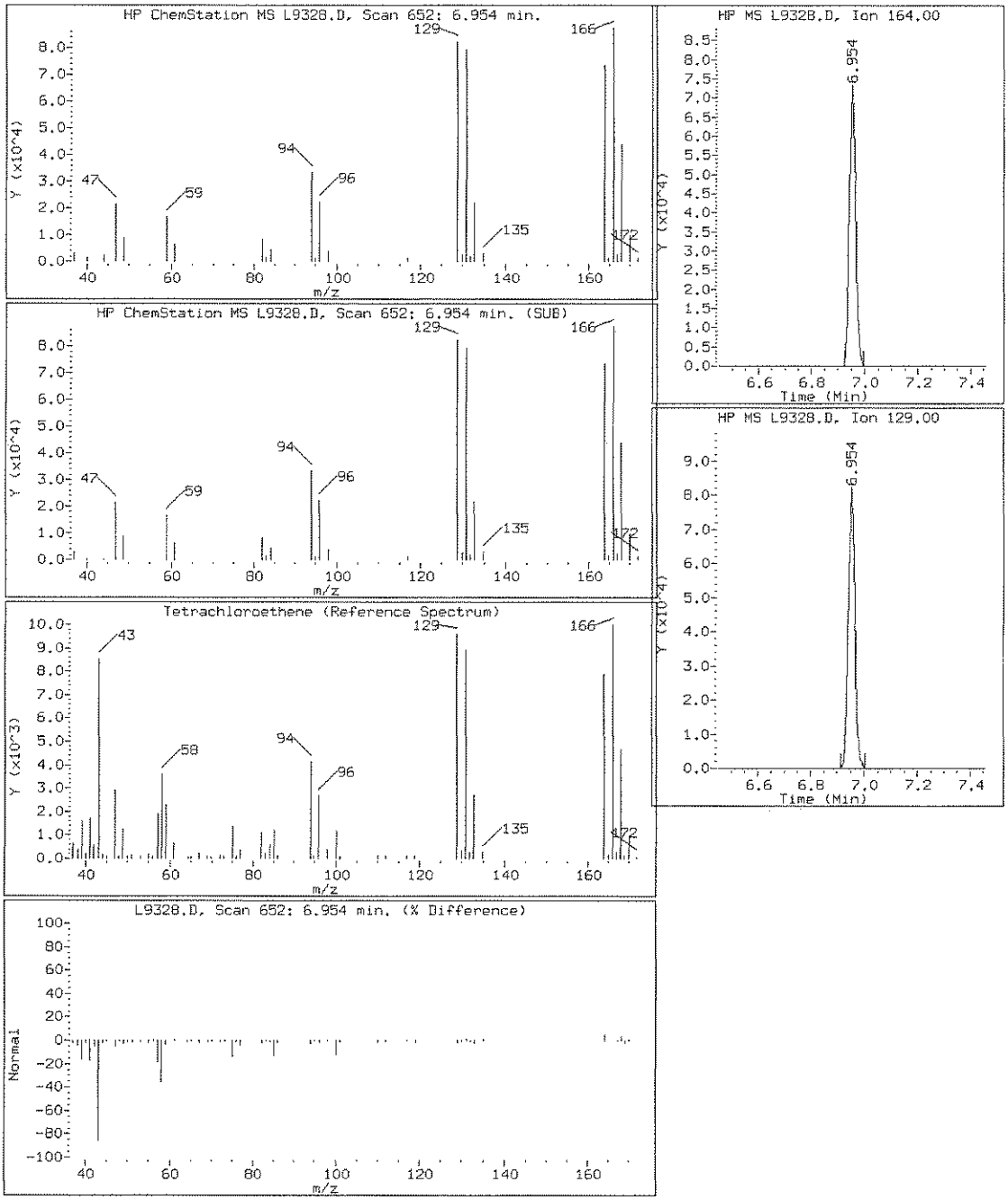
Data File: L9328.D

Date: 01-AUG-2007 03:48

Client ID: A-2 Instrument: msl.i

Sample Info: 220-2277-B-1 Operator: D. GAYDA

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-6 | Lab Sample ID: 220-2277-2

Matrix: Water | Lab File ID: L9386.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/02/2007 16:48

Level: (low/med) Low | Dilution Factor: 2

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8356 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	20	U	20	2.8
71-43-2	Benzene	10	U	10	0.80
75-27-4	Bromodichloromethane	10	U	10	0.80
75-25-2	Bromoform	10	U	10	1.6
74-83-9	Bromomethane	10	U	10	2.4
78-93-3	2-Butanone (MEK)	20	U	20	2.4
75-15-0	Carbon disulfide	10	U	10	1.8
56-23-5	Carbon tetrachloride	10	U	10	2.0
108-90-7	Chlorobenzene	10	U	10	0.80
75-00-3	Chloroethane	10	U	10	1.6
67-66-3	Chloroform	10	U	10	1.4
74-87-3	Chloromethane	10	U	10	1.0
124-48-1	Dibromochloromethane	10	U	10	1.0
75-34-3	1,1-Dichloroethane	10	U	10	1.2
107-06-2	1,2-Dichloroethane	10	U	10	1.2
75-35-4	1,1-Dichloroethene	10	U	10	1.4
78-87-5	1,2-Dichloropropane	10	U	10	1.8
10061-01-5	cis-1,3-Dichloropropene	10	U	10	1.0
10061-02-6	trans-1,3-Dichloropropene	10	U	10	1.6
100-41-4	Ethylbenzene	10	U	10	2.0
591-78-6	2-Hexanone	20	U	20	1.6
75-09-2	Methylene Chloride	10	U	10	0.80
108-10-1	4-Methyl-2-pentanone (MIBK)	20	U	20	1.4
100-42-5	Styrene	10	U	10	1.0
79-34-5	1,1,2,2-Tetrachloroethane	10	U	10	0.80
127-18-4	Tetrachloroethene	290		10	1.0
108-88-3	Toluene	10	U	10	0.60
71-55-6	1,1,1-Trichloroethane	10	U	10	0.80
79-00-5	1,1,2-Trichloroethane	10	U	10	1.2
79-01-6	Trichloroethene	10	U	10	1.4
75-01-4	Vinyl chloride	10	U	10	1.6
1330-20-7	Xylenes, Total	10	U	10	2.0
156-59-2	cis-1,2-Dichloroethene	10	U	10	1.2
156-60-5	trans-1,2-Dichloroethene	10	U	10	1.0

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079370.b\L9386.D
Lab Smp Id: 220-2277-A-2 Client Smp ID: A-6
Inj Date : 02-AUG-2007 16:48 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-A-2
Misc Info : : ;;; ; 8260 ; 2 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 68
Dil Factor: 2.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	2.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

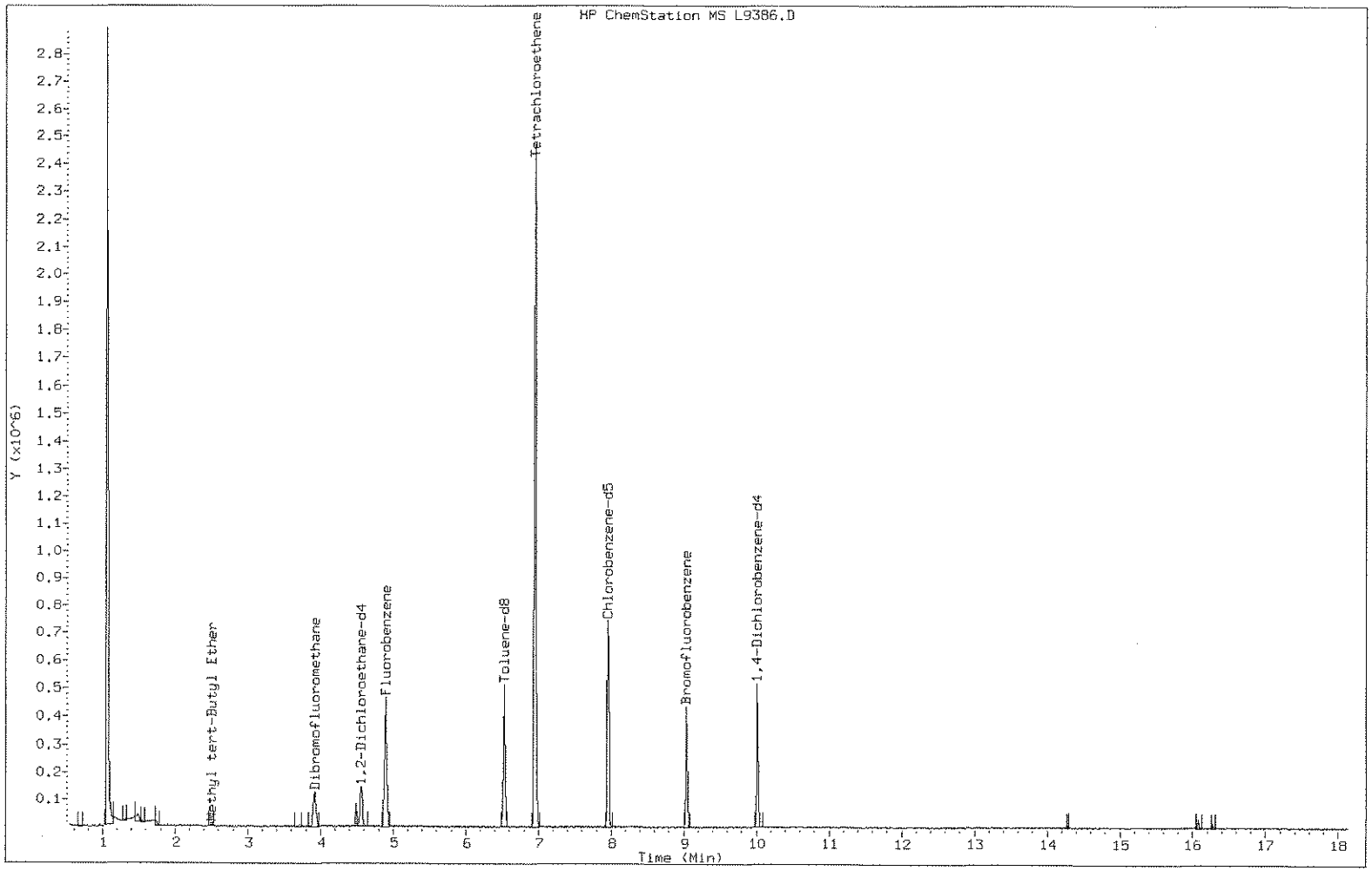
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.892	4.901	(1.000)	436691	25.0000	
	24 Methyl tert-Butyl Ether	73	2.491	2.491	(0.509)	15432	0.85117	2
$	41 Dibromofluoromethane	111	3.918	3.927	(0.801)	109485	19.4338	19
$	55 1,2-Dichloroethane-d4	65	4.558	4.567	(0.932)	130582	19.0115	19
*	75 Chlorobenzene-d5	117	7.952	7.961	(1.000)	426839	25.0000	
$	77 Toluene-d8	98	6.525	6.535	(0.821)	315186	20.7668	21
	80 Tetrachloroethene	164	6.949	6.958	(0.874)	526255	146.741	290
*	95 1,4-Dichlorobenzene-d4	152	10.008	10.018	(1.000)	112968	25.0000	
$	125 Bromofluorobenzene	95	9.034	9.043	(0.903)	127611	30.1107	30

Data File: L9386.D

Date: 02-AUG-2007 16:48

Client ID: A-6 Instrument: msl.i

Sample Info: 220-2277-A-2 Operator: D. HUMBERT



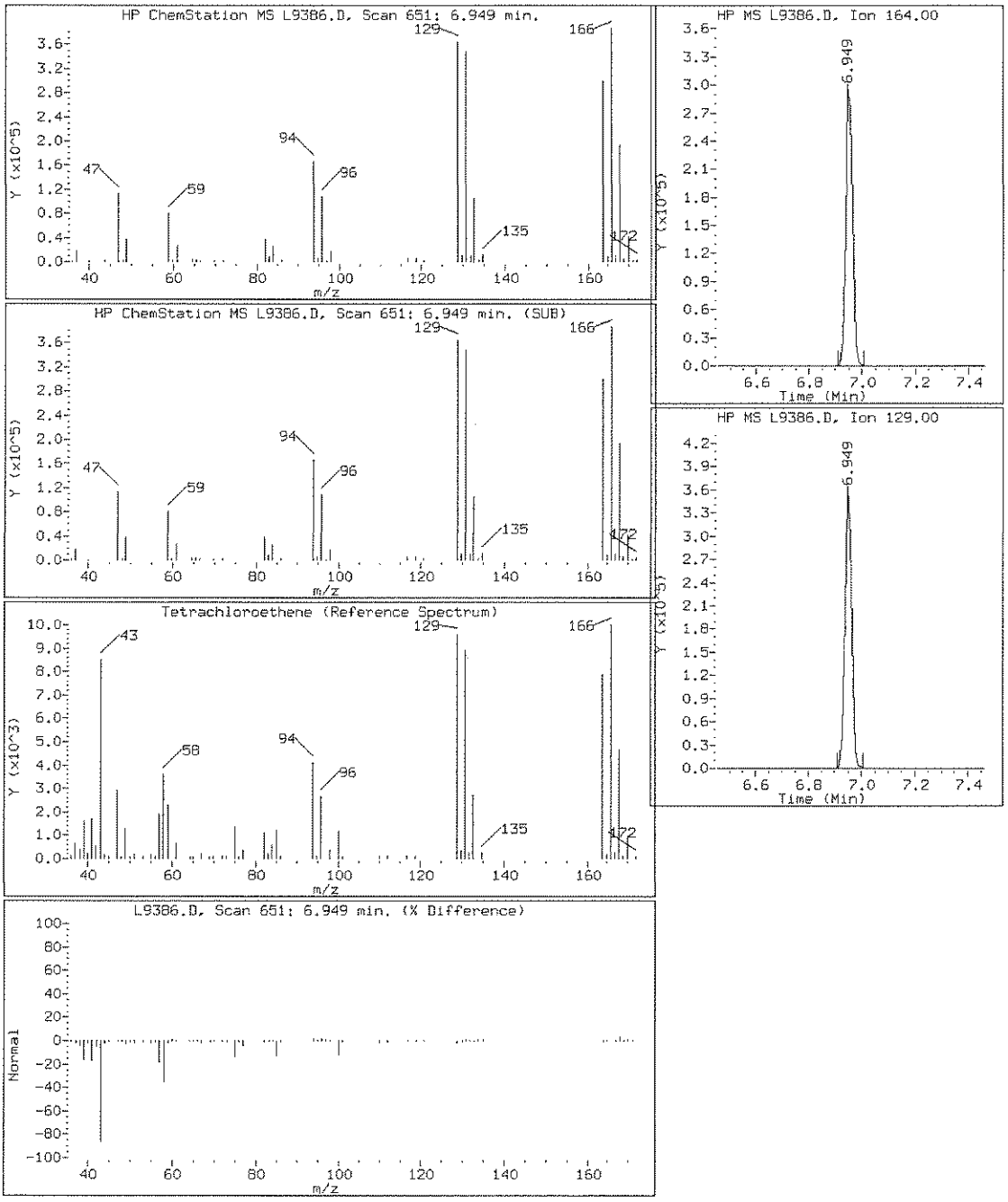
Data File: L9386.D

Date: 02-AUG-2007 16:48

Client ID: A-6 Instrument: msl.i

Sample Info: 220-2277-A-2 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-7 | Lab Sample ID: 220-2277-3

Matrix: Water | Lab File ID: L9330.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/01/2007 04:38

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8307 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	3.8	J	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	26		10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	41		5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079310.b\L9330.D
Lab Smp Id: 220-2277-A-3 Client Smp ID: A-7
Inj Date : 01-AUG-2007 04:38 MS Autotune Date: 26-APR-2004 14:21
Operator : D. GAYDA Inst ID: msl.i
Smp Info : 220-2277-A-3
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079310.b\L8260BNW.m
Meth Date : 31-Jul-2007 21:07 target Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 21
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

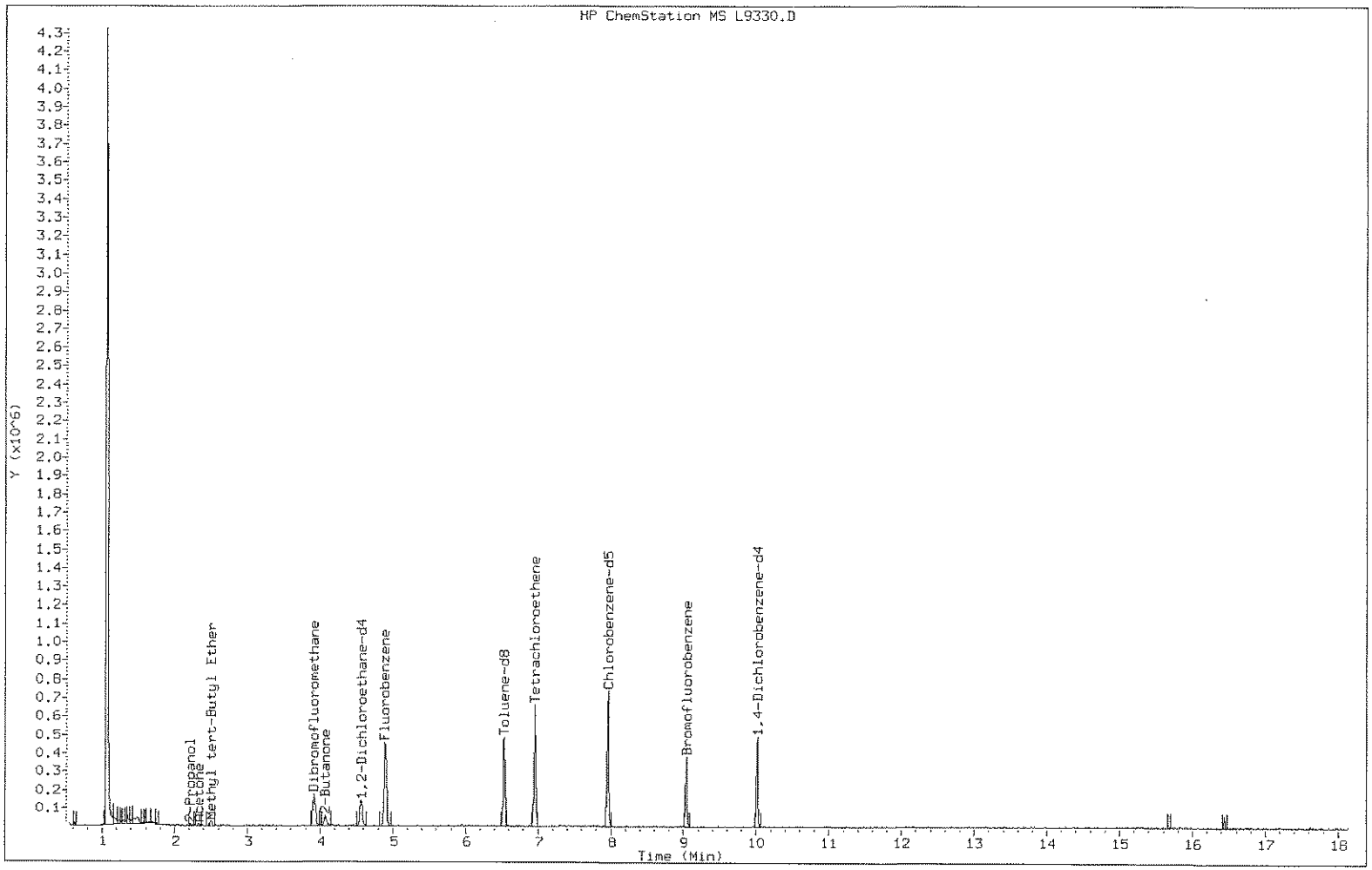
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.900	4.896	(1.000)	451910	25.0000	
	18 2-Propanol	45	2.214	2.219	(0.452)	12881	12.4211	12
	21 Acetone	43	2.322	2.328	(0.474)	10801	3.76874	4
	24 Methyl tert-Butyl Ether	73	2.499	2.495	(0.510)	29245	1.55872	2
$	41 Dibromofluoromethane	111	3.916	3.922	(0.799)	109224	18.7346	19
	45 2-Butanone	43	4.073	4.069	(0.831)	106985	25.7164	26
$	55 1,2-Dichloroethane-d4	65	4.565	4.561	(0.932)	131881	18.5540	18
*	75 Chlorobenzene-d5	117	7.960	7.956	(1.000)	425701	25.0000	
$	77 Toluene-d8	98	6.533	6.529	(0.821)	314755	20.7939	21
	80 Tetrachloroethene	164	6.956	6.952	(0.874)	146801	41.0434	41
*	95 1,4-Dichlorobenzene-d4	152	10.016	10.012	(1.000)	119044	25.0000	
$	125 Bromofluorobenzene	95	9.032	9.038	(0.902)	125454	28.0908	28

Data File: L9330.D

Date: 01-AUG-2007 04:38

Client ID: A-7 Instrument: msl.i

Sample Info: 220-2277-A-3 Operator: D. GAYDA



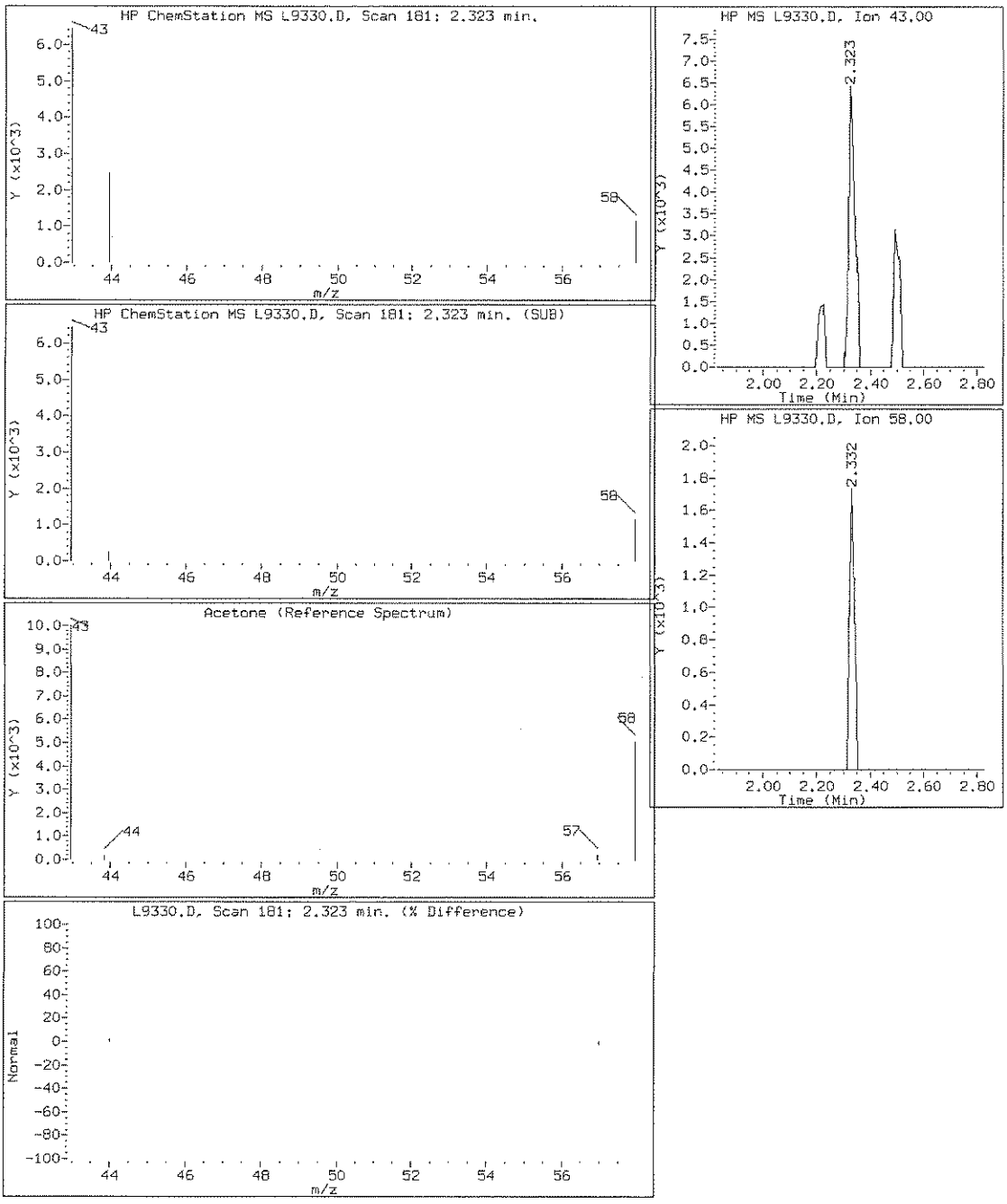
Data File: L9330.D

Date: 01-AUG-2007 04:38

Client ID: A-7 Instrument: msl.i

Sample Info: 220-2277-A-3 Operator: D. GAYDA

21 Acetone



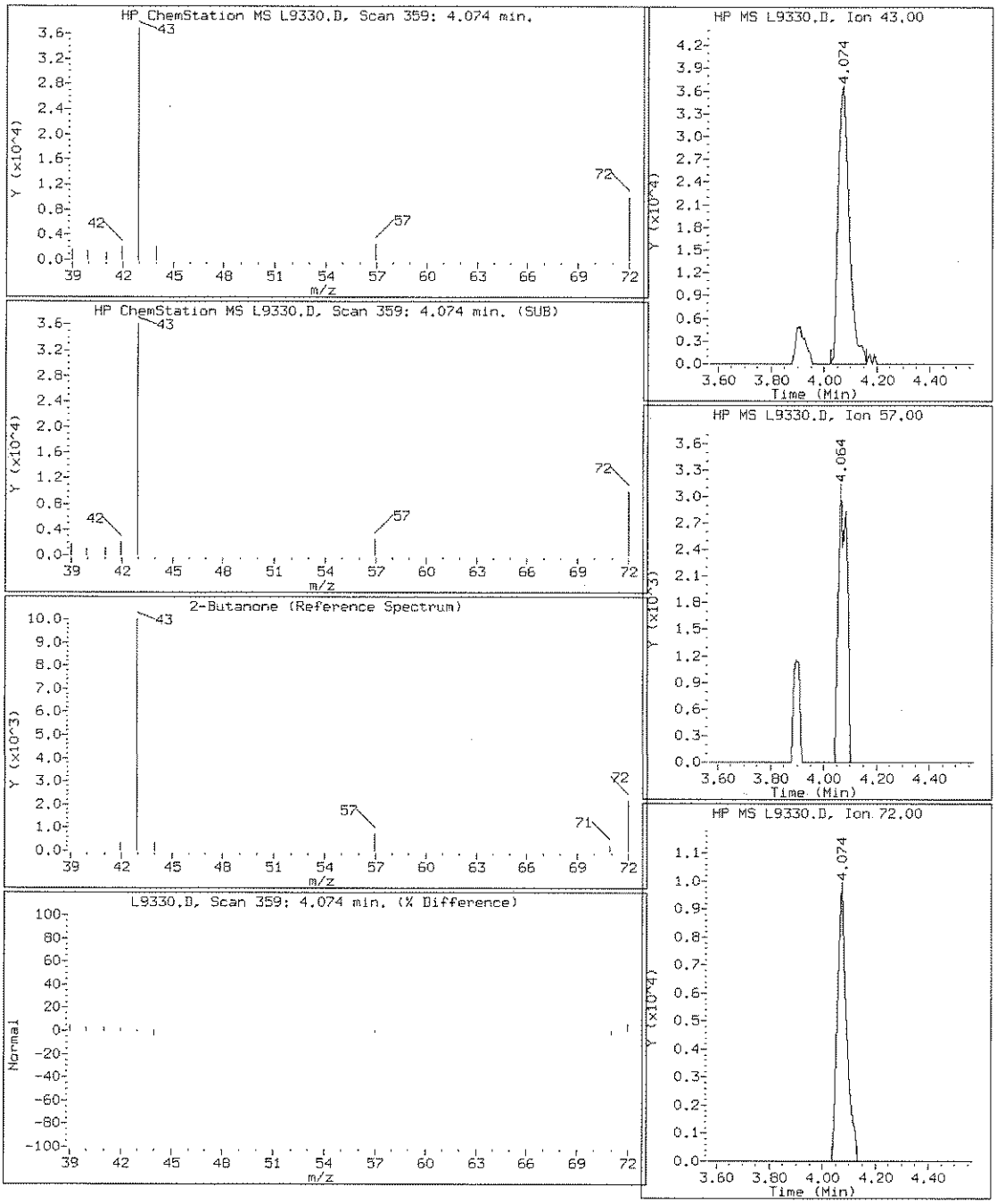
Data File: L9330.D

Date: 01-AUG-2007 04:38

Client ID: A-7 Instrument: msl.i

Sample Info: 220-2277-A-3 Operator: D. GAYDA

45 2-Butanone



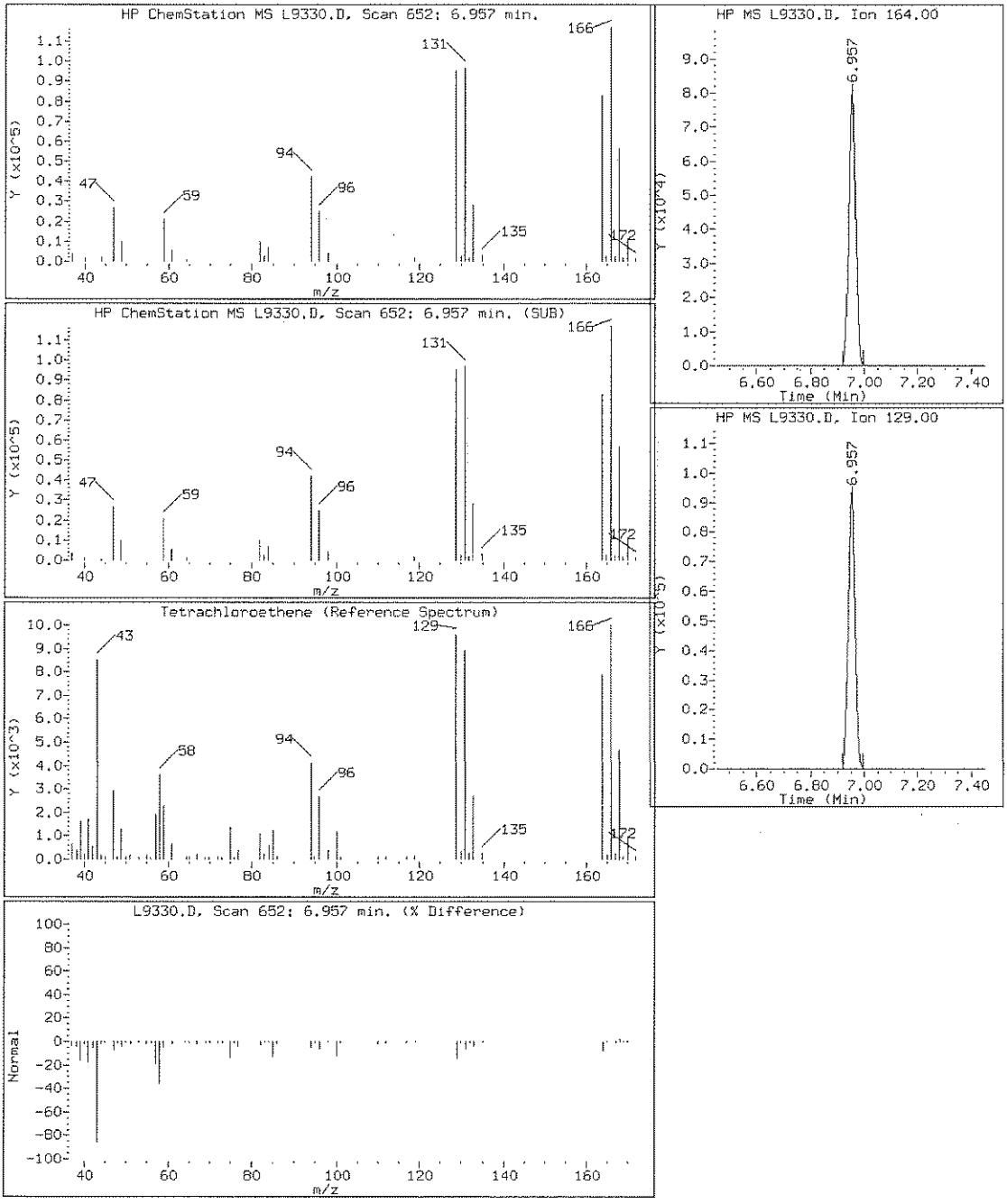
Data File: L9330.D

Date: 01-AUG-2007 04:38

Client ID: A-7 Instrument: msl.i

Sample Info: 220-2277-A-3 Operator: D. GAYDA

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-8 Lab Sample ID: 220-2277-4

Matrix: Water Lab File ID: L9382.D

Analysis Method: 8260B Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) Date Analyzed: 08/02/2007 15:08

Level: (low/med) Low Dilution Factor: 50

GC Column/ID: RTX-VMS 0.18 (mm) Soil Aliquot:

Soil Extract Vol.: % Moisture:

Analy. Batch No.: 8356 Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	97	J	500	70
71-43-2	Benzene	250	U	250	20
75-27-4	Bromodichloromethane	250	U	250	20
75-25-2	Bromoform	250	U	250	40
74-83-9	Bromomethane	250	U	250	60
78-93-3	2-Butanone (MEK)	500	U	500	60
75-15-0	Carbon disulfide	250	U	250	45
56-23-5	Carbon tetrachloride	250	U	250	50
108-90-7	Chlorobenzene	250	U	250	20
75-00-3	Chloroethane	250	U	250	40
67-66-3	Chloroform	250	U	250	35
74-87-3	Chloromethane	250	U	250	25
124-48-1	Dibromochloromethane	250	U	250	25
75-34-3	1,1-Dichloroethane	250	U	250	30
107-06-2	1,2-Dichloroethane	250	U	250	30
75-35-4	1,1-Dichloroethene	250	U	250	35
78-87-5	1,2-Dichloropropane	250	U	250	45
10061-01-5	cis-1,3-Dichloropropene	250	U	250	25
10061-02-6	trans-1,3-Dichloropropene	250	U	250	40
100-41-4	Ethylbenzene	250	U	250	50
591-78-6	2-Hexanone	500	U	500	40
75-09-2	Methylene Chloride	250	U	250	20
108-10-1	4-Methyl-2-pentanone (MIBK)	500	U	500	35
100-42-5	Styrene	250	U	250	25
79-34-5	1,1,2,2-Tetrachloroethane	250	U	250	20
127-18-4	Tetrachloroethene	4400		250	25
108-88-3	Toluene	250	U	250	15
71-55-6	1,1,1-Trichloroethane	250	U	250	20
79-00-5	1,1,2-Trichloroethane	250	U	250	30
79-01-6	Trichloroethene	250	U	250	35
75-01-4	Vinyl chloride	250	U	250	40
1330-20-7	Xylenes, Total	250	U	250	50
156-59-2	cis-1,2-Dichloroethene	250	U	250	30
156-60-5	trans-1,2-Dichloroethene	250	U	250	25

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079370.b\L9382.D
Lab Smp Id: 220-2277-B-4 Client Smp ID: A-8
Inj Date : 02-AUG-2007 15:08 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-B-4
Misc Info : : ;;; ; 8260 ; 50 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 64
Dil Factor: 50.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	50.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

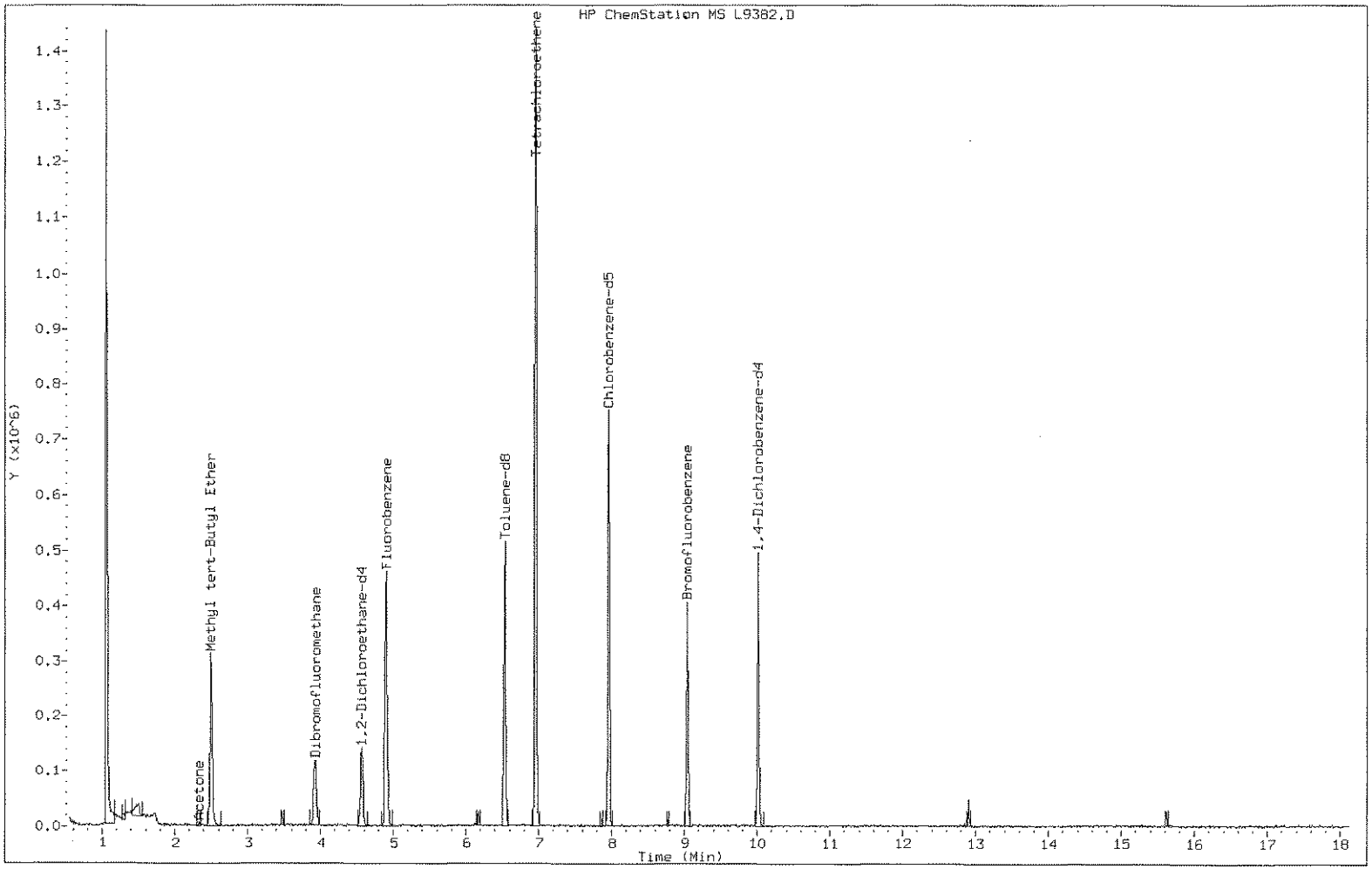
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.901	4.901	(1.000)	452451	25.0000	
	21 Acetone	43	2.333	2.323	(0.476)	5573	1.94223	97
	24 Methyl tert-Butyl Ether	73	2.491	2.491	(0.508)	342850	18.2516	910
$	41 Dibromofluoromethane	111	3.917	3.927	(0.799)	109090	18.6892	19
$	55 1,2-Dichloroethane-d4	65	4.567	4.567	(0.932)	132348	18.5974	18
*	75 Chlorobenzene-d5	117	7.961	7.961	(1.000)	428736	25.0000	
$	77 Toluene-d8	98	6.534	6.535	(0.821)	320424	21.0185	21
	80 Tetrachloroethene	164	6.958	6.958	(0.874)	316947	87.9865	4400
*	95 1,4-Dichlorobenzene-d4	152	10.017	10.018	(1.000)	121030	25.0000	
$	125 Bromofluorobenzene	95	9.043	9.043	(0.903)	132624	29.2090	29

Data File: L9382.D

Date: 02-AUG-2007 15:08

Client ID: A-8 Instrument: msl.i

Sample Info: 220-2277-B-4 Operator: D. HUMBERT



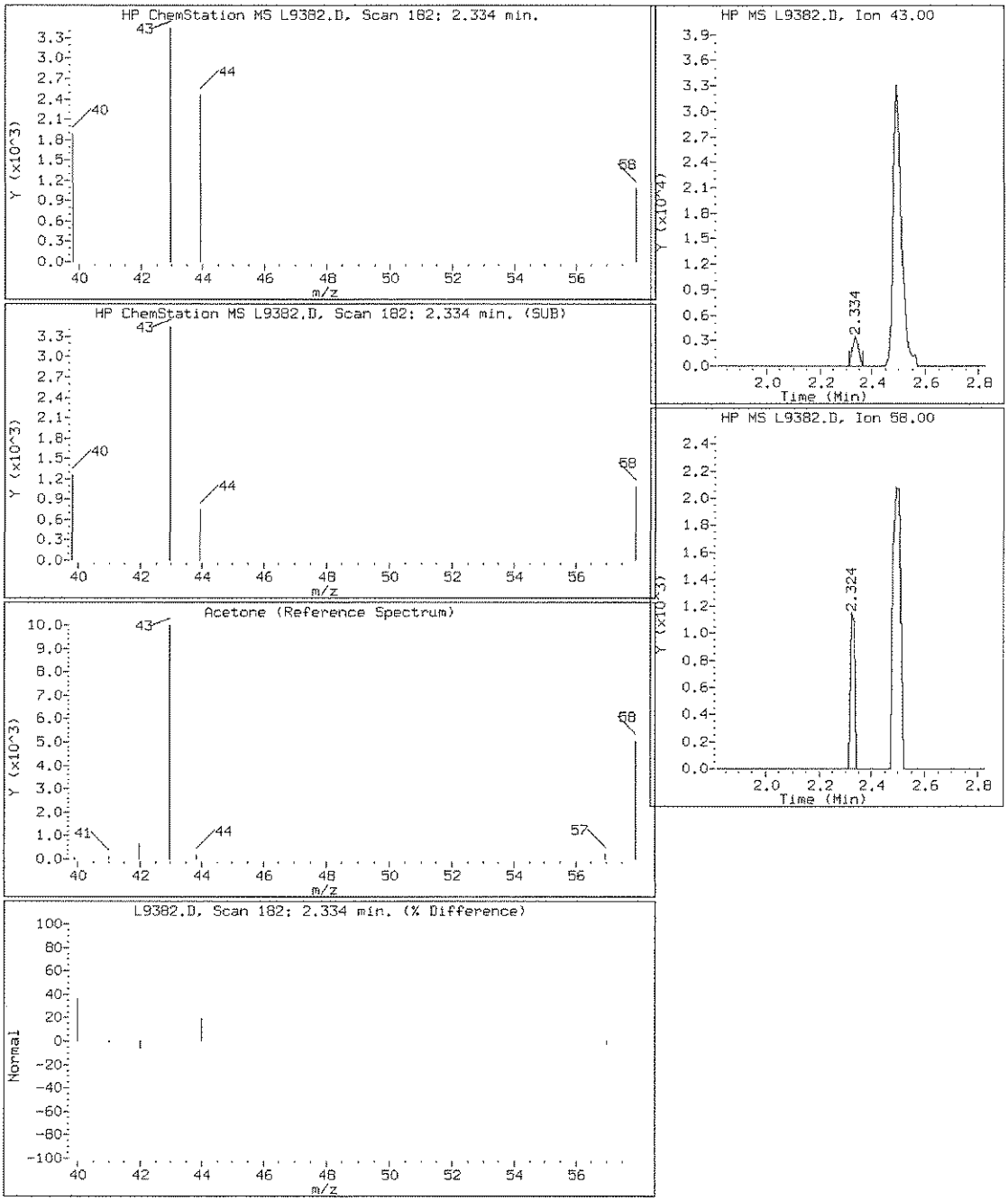
Data File: L9382.D

Date: 02-AUG-2007 15:08

Client ID: A-8 Instrument: msl.i

Sample Info: 220-2277-B-4 Operator: D. HUMBERT

21 Acetone



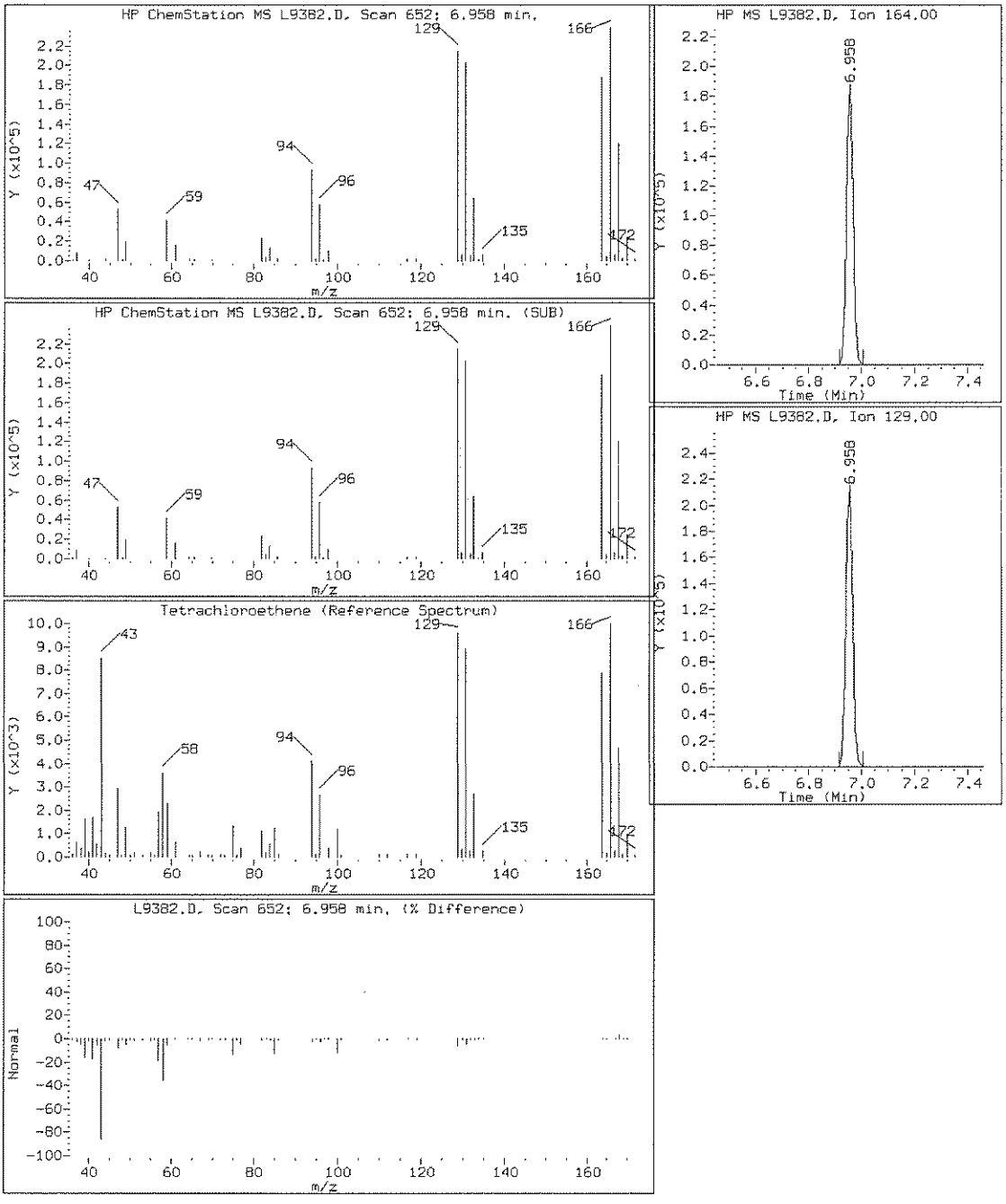
Data File: L9382.D

Date: 02-AUG-2007 15:08

Client ID: A-8 Instrument: msl.i

Sample Info: 220-2277-B-4 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-3 | Lab Sample ID: 220-2277-5

Matrix: Water | Lab File ID: L9376.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/02/2007 12:32

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8356 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	1.4	J	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	14		5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079370.b\L9376.D
Lab Smp Id: 220-2277-A-5 Client Smp ID: A-3
Inj Date : 02-AUG-2007 12:32 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-A-5
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 58
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

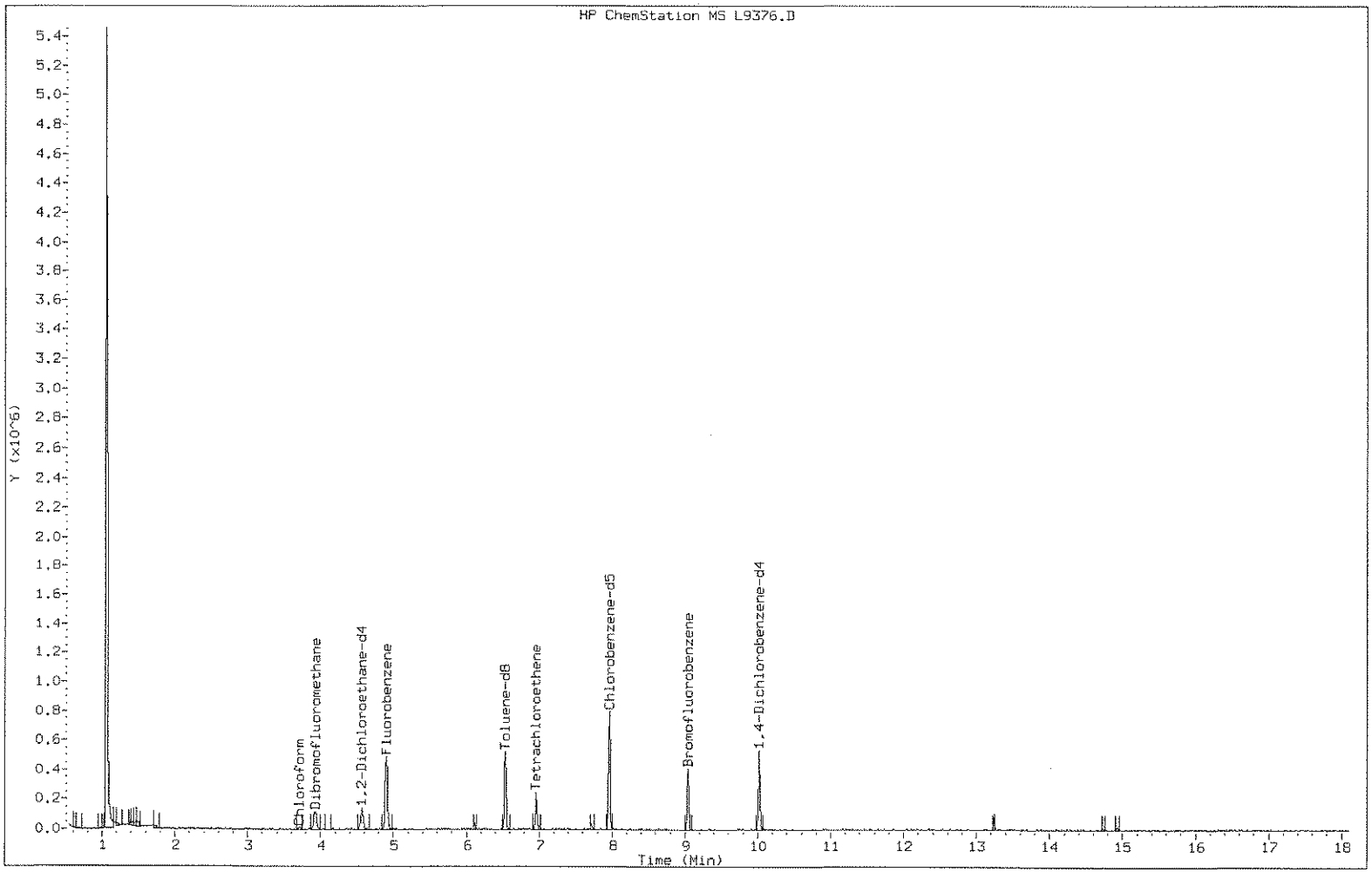
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.901	4.901	(1.000)	462027	25.0000	
	38 Chloroform	83	3.711	3.701	(0.757)	13416	1.38965	1
$	41 Dibromofluoromethane	111	3.927	3.927	(0.801)	107656	18.0613	18
$	55 1,2-Dichloroethane-d4	65	4.567	4.567	(0.932)	129675	17.8441	18
*	75 Chlorobenzene-d5	117	7.961	7.961	(1.000)	454599	25.0000	
$	77 Toluene-d8	98	6.534	6.535	(0.821)	314630	19.4643	19
	80 Tetrachloroethene	164	6.958	6.958	(0.874)	51606	13.5111	14
*	95 1,4-Dichlorobenzene-d4	152	10.017	10.018	(1.000)	124058	25.0000	
$	125 Bromofluorobenzene	95	9.043	9.043	(0.903)	130765	28.0966	28

Data File: L9376.D

Date: 02-AUG-2007 12:32

Client ID: A-3 Instrument: msl.i

Sample Info: 220-2277-A-5 Operator: D. HUMBERT



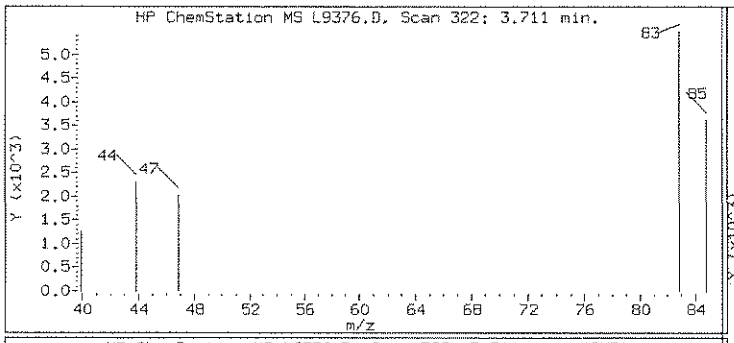
Data File: L9376.D

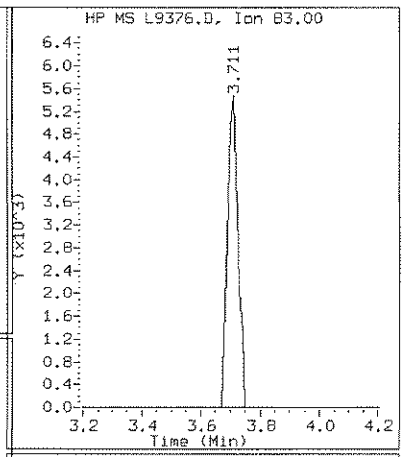
Date: 02-AUG-2007 12:32

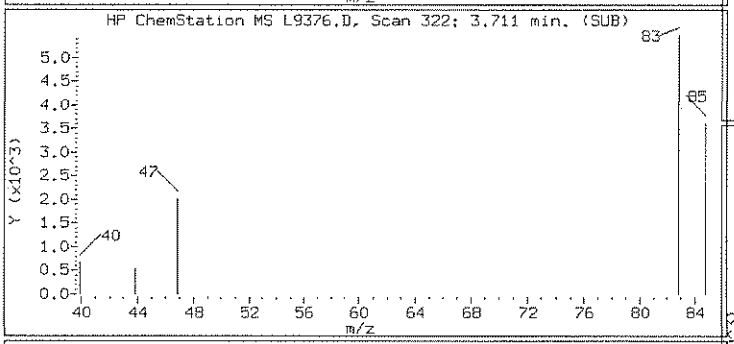
Client ID: A-3 Instrument: msl.i

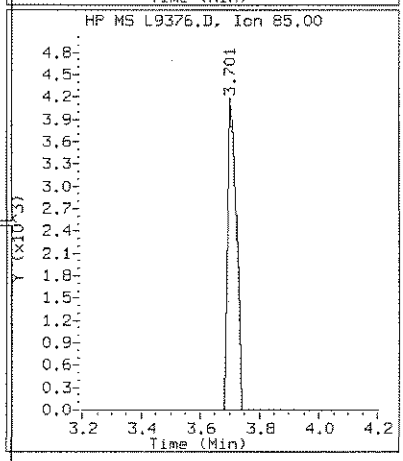
Sample Info: 220-2277-A-5 Operator: D. HUMBERT

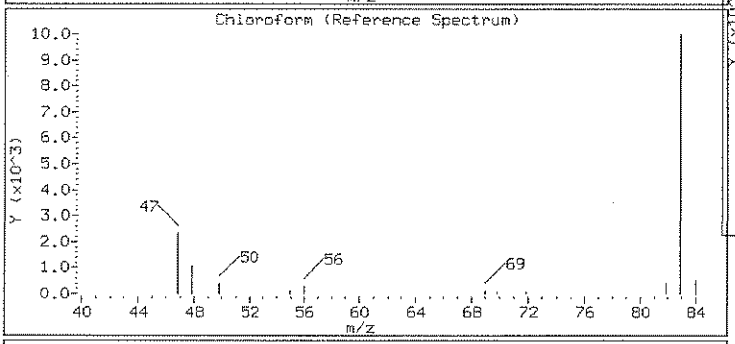
38 Chloroform

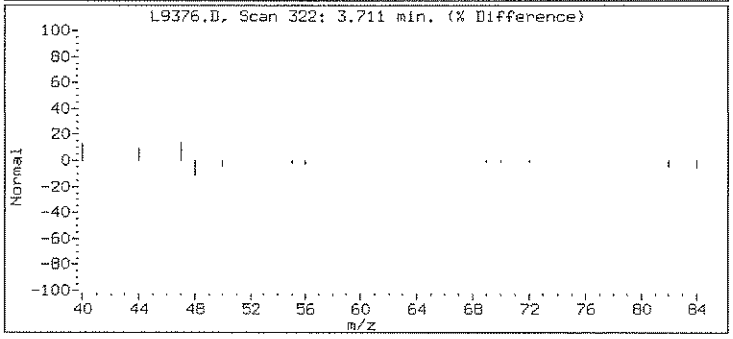












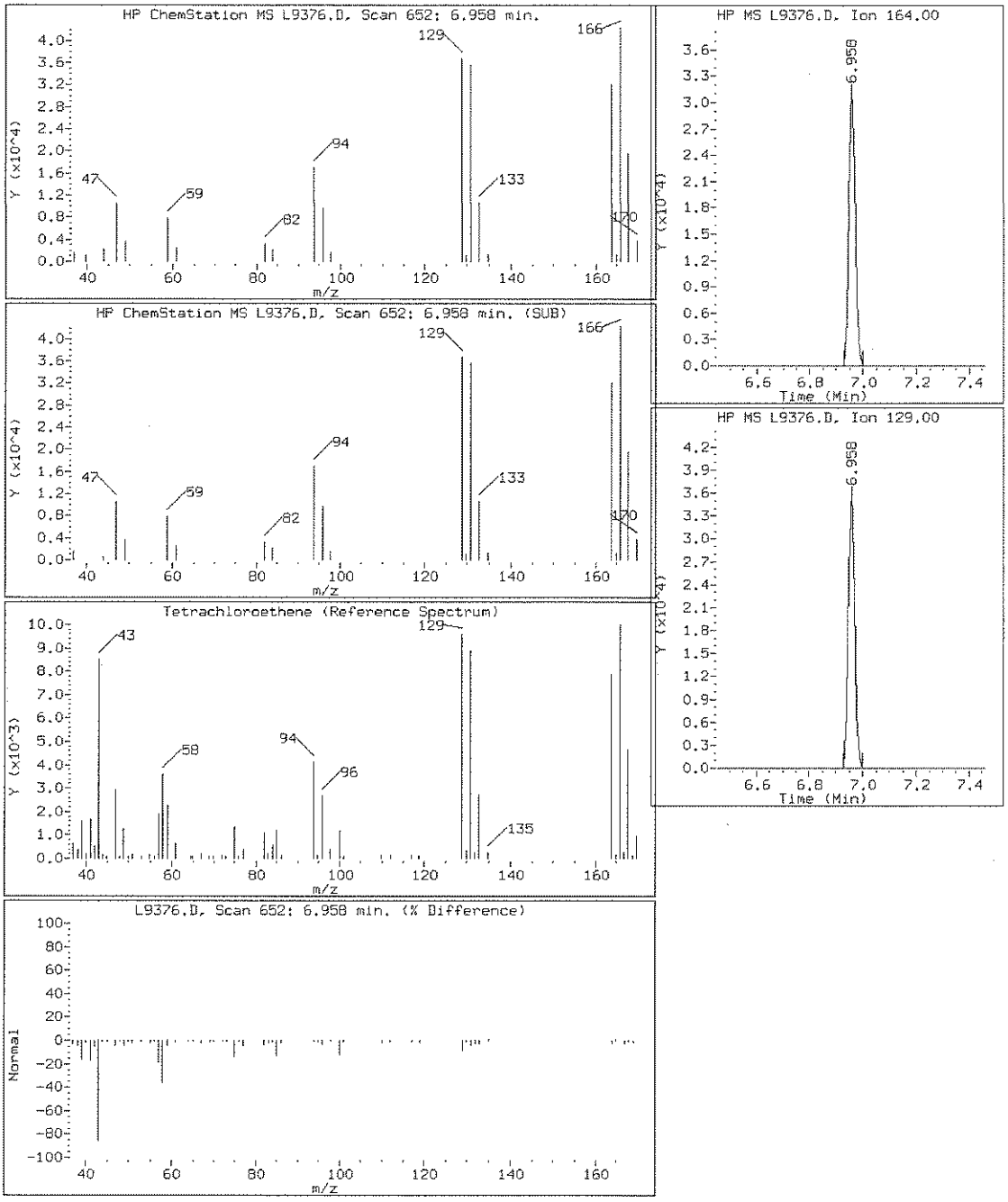
Data File: L9376.D

Date: 02-AUG-2007 12:32

Client ID: A-3 Instrument: msl.i

Sample Info: 220-2277-A-5 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-11 | Lab Sample ID: 220-2277-6

Matrix: Water | Lab File ID: L9377.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/02/2007 12:57

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8356 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	8.5	J	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	0.78	J	5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079370.b\L9377.D
Lab Smp Id: 220-2277-B-6 Client Smp ID: A-11
Inj Date : 02-AUG-2007 12:57 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-B-6
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 59
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

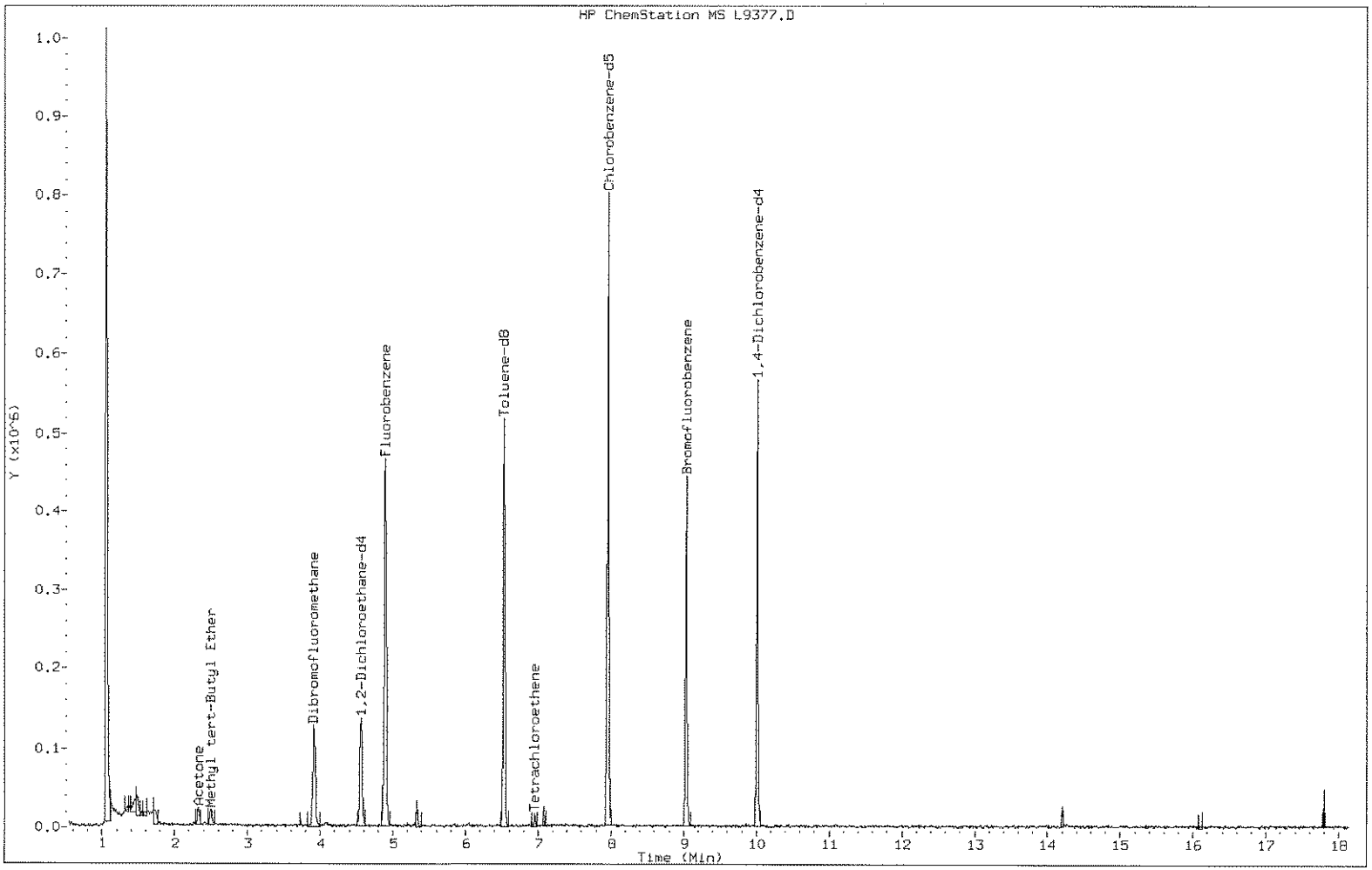
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.896	4.901	(1.000)	447349	25.0000	
	21 Acetone	43	2.328	2.323	(0.476)	24236	8.54276	8
	24 Methyl tert-Butyl Ether	73	2.495	2.491	(0.510)	24581	1.32349	1
$	41 Dibromofluoromethane	111	3.912	3.927	(0.799)	110202	19.0950	19
$	55 1,2-Dichloroethane-d4	65	4.561	4.567	(0.932)	127976	18.1881	18
*	75 Chlorobenzene-d5	117	7.956	7.961	(1.000)	437111	25.0000	
$	77 Toluene-d8	98	6.529	6.535	(0.821)	314931	20.2624	20
	80 Tetrachloroethene	164	6.952	6.958	(0.874)	2868	0.78092	0.8
*	95 1,4-Dichlorobenzene-d4	152	10.012	10.018	(1.000)	122354	25.0000	
$	125 Bromofluorobenzene	95	9.038	9.043	(0.903)	132858	28.9439	29

Data File: L9377.D

Date: 02-AUG-2007 12:57

Client ID: A-11 Instrument: msl.i

Sample Info: 220-2277-B-6 Operator: D. HUMBERT



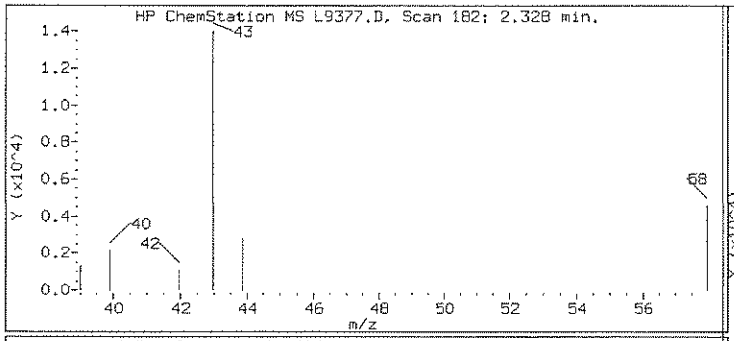
Data File: L9377.D

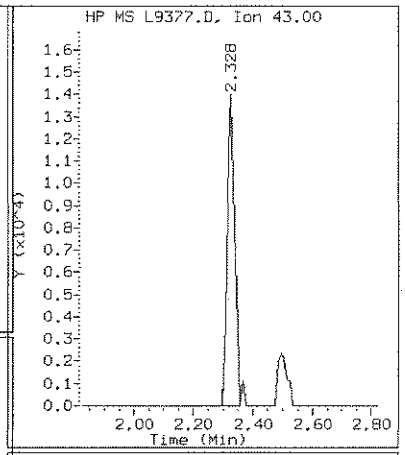
Date: 02-AUG-2007 12:57

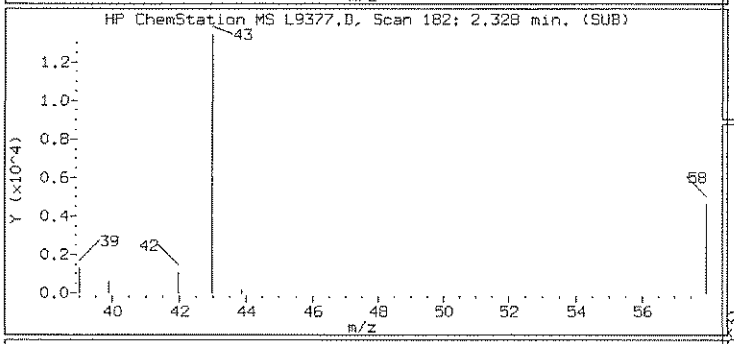
Client ID: A-11 Instrument: msl.i

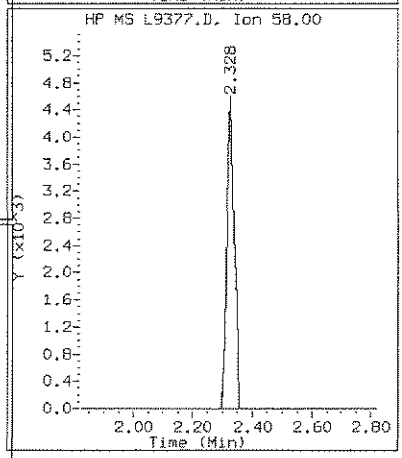
Sample Info: 220-2277-B-6 Operator: D. HUMBERT

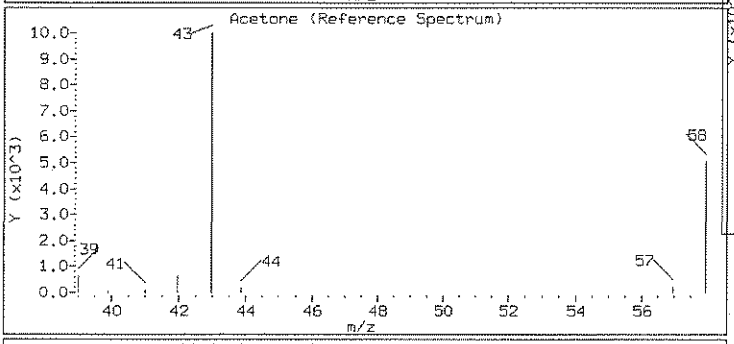
21 Acetone

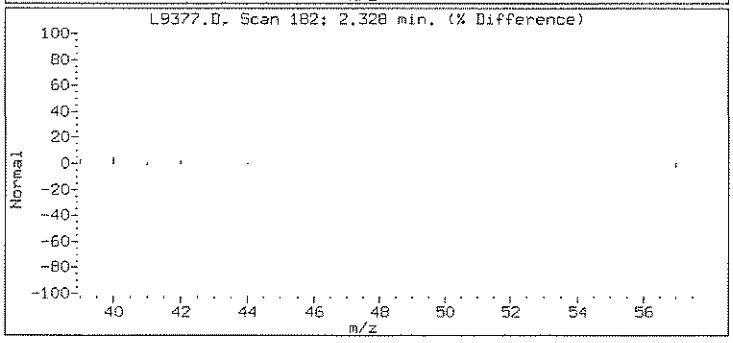












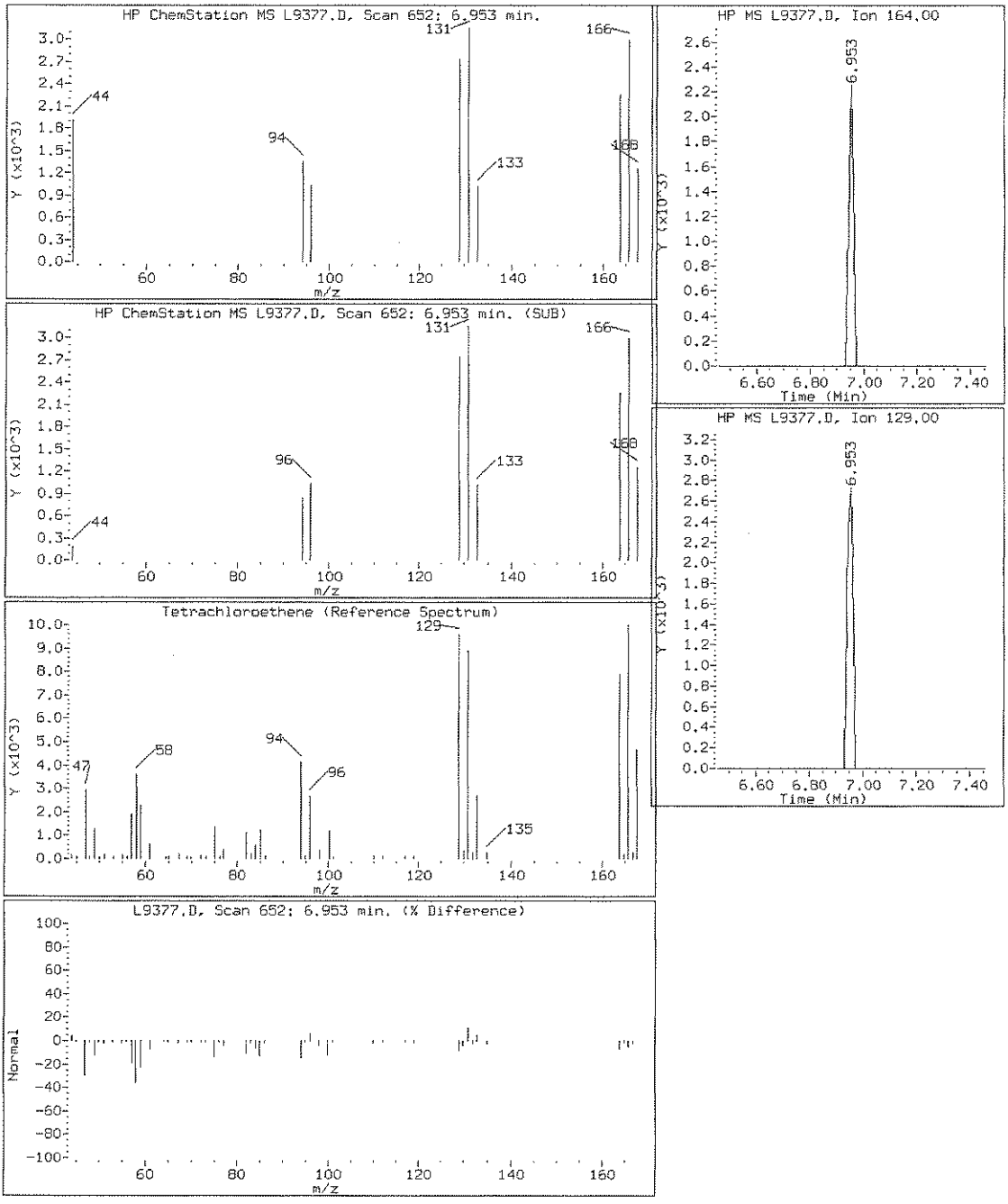
Data File: L9377.D

Date: 02-AUG-2007 12:57

Client ID: A-11 Instrument: msl.i

Sample Info: 220-2277-B-6 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-12 Lab Sample ID: 220-2277-7

Matrix: Water Lab File ID: L9383.D

Analysis Method: 8260B Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) Date Analyzed: 08/02/2007 15:33

Level: (low/med) Low Dilution Factor: 20

GC Column/ID: RTX-VMS 0.18 (mm) Soil Aliquot:

Soil Extract Vol.: % Moisture:

Analy. Batch No.: 8356 Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	200	U	200	28
71-43-2	Benzene	100	U	100	8.0
75-27-4	Bromodichloromethane	100	U	100	8.0
75-25-2	Bromoform	100	U	100	16
74-83-9	Bromomethane	100	U	100	24
78-93-3	2-Butanone (MEK)	200	U	200	24
75-15-0	Carbon disulfide	100	U	100	18
56-23-5	Carbon tetrachloride	100	U	100	20
108-90-7	Chlorobenzene	100	U	100	8.0
75-00-3	Chloroethane	100	U	100	16
67-66-3	Chloroform	100	U	100	14
74-87-3	Chloromethane	100	U	100	10
124-48-1	Dibromochloromethane	100	U	100	10
75-34-3	1,1-Dichloroethane	100	U	100	12
107-06-2	1,2-Dichloroethane	100	U	100	12
75-35-4	1,1-Dichloroethene	100	U	100	14
78-87-5	1,2-Dichloropropane	100	U	100	18
10061-01-5	cis-1,3-Dichloropropene	100	U	100	10
10061-02-6	trans-1,3-Dichloropropene	100	U	100	16
100-41-4	Ethylbenzene	100	U	100	20
591-78-6	2-Hexanone	200	U	200	16
75-09-2	Methylene Chloride	100	U	100	8.0
108-10-1	4-Methyl-2-pentanone (MIBK)	200	U	200	14
100-42-5	Styrene	100	U	100	10
79-34-5	1,1,2,2-Tetrachloroethane	100	U	100	8.0
127-18-4	Tetrachloroethene	1800		100	10
108-88-3	Toluene	100	U	100	6.0
71-55-6	1,1,1-Trichloroethane	100	U	100	8.0
79-00-5	1,1,2-Trichloroethane	100	U	100	12
79-01-6	Trichloroethene	100	U	100	14
75-01-4	Vinyl chloride	100	U	100	16
1330-20-7	Xylenes, Total	100	U	100	20
156-59-2	cis-1,2-Dichloroethene	100	U	100	12
156-60-5	trans-1,2-Dichloroethene	100	U	100	10

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079370.b\L9383.D
Lab Smp Id: 220-2277-B-7 Client Smp ID: A-12
Inj Date : 02-AUG-2007 15:33 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-B-7
Misc Info : : ;;; ; 8260 ; 20 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 65
Dil Factor: 20.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	20.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

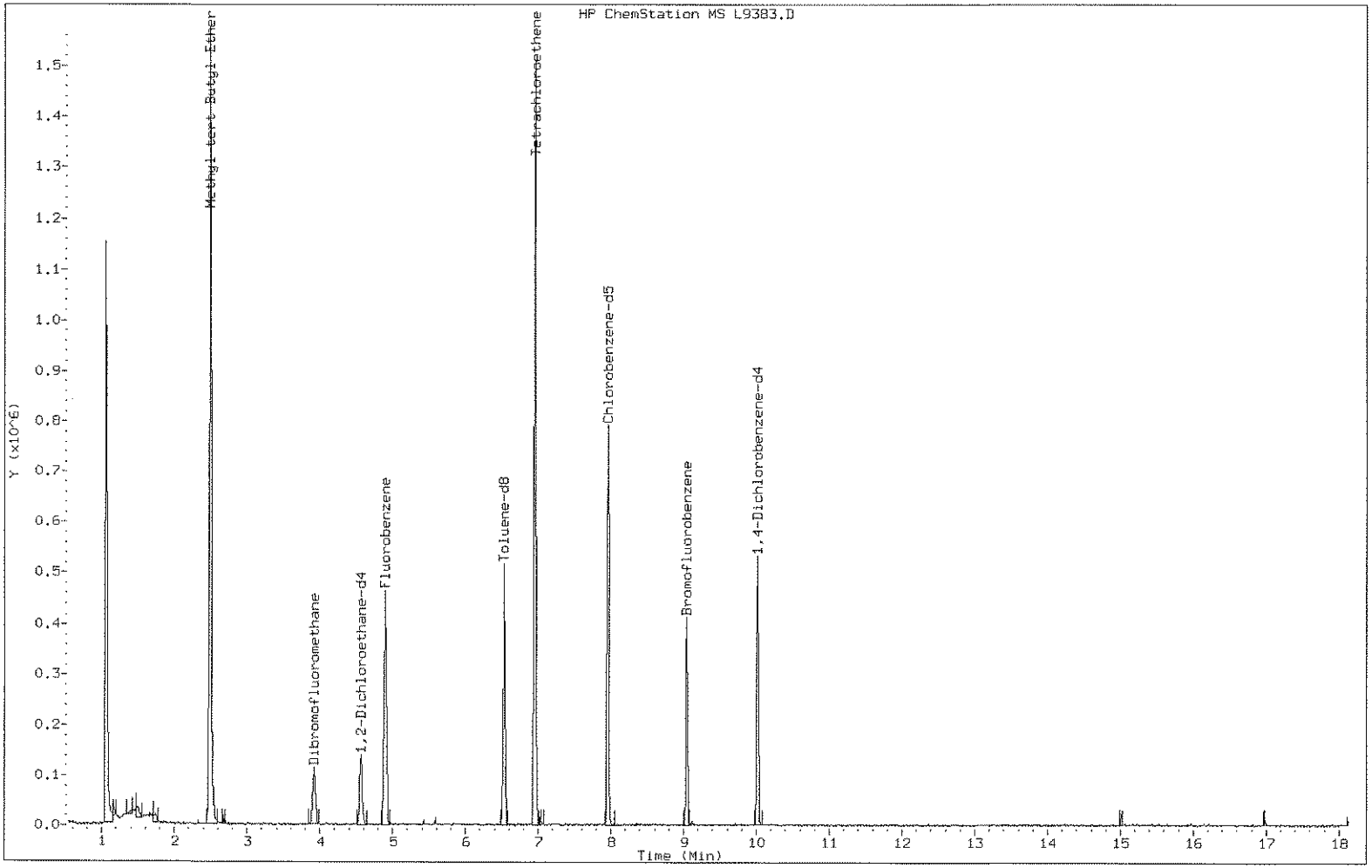
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.897	4.901	(1.000)	452541	25.0000	
	24 Methyl tert-Butyl Ether	73	2.486	2.491	(0.508)	1596201	84.9569	1700
$	41 Dibromofluoromethane	111	3.913	3.927	(0.799)	104807	17.9519	18
$	55 1,2-Dichloroethane-d4	65	4.562	4.567	(0.932)	132894	18.6704	19
*	75 Chlorobenzene-d5	117	7.957	7.961	(1.000)	436646	25.0000	
$	77 Toluene-d8	98	6.530	6.535	(0.821)	314256	20.2405	20
	80 Tetrachloroethene	164	6.953	6.958	(0.874)	328725	89.6030	1800
*	95 1,4-Dichlorobenzene-d4	152	10.013	10.018	(1.000)	124274	25.0000	
$	125 Bromofluorobenzene	95	9.039	9.043	(0.903)	130239	27.9350	28

Data File: L9383.D

Date: 02-AUG-2007 15:33

Client ID: A-12 Instrument: msl.i

Sample Info: 220-2277-B-7 Operator: D. HUMBERT



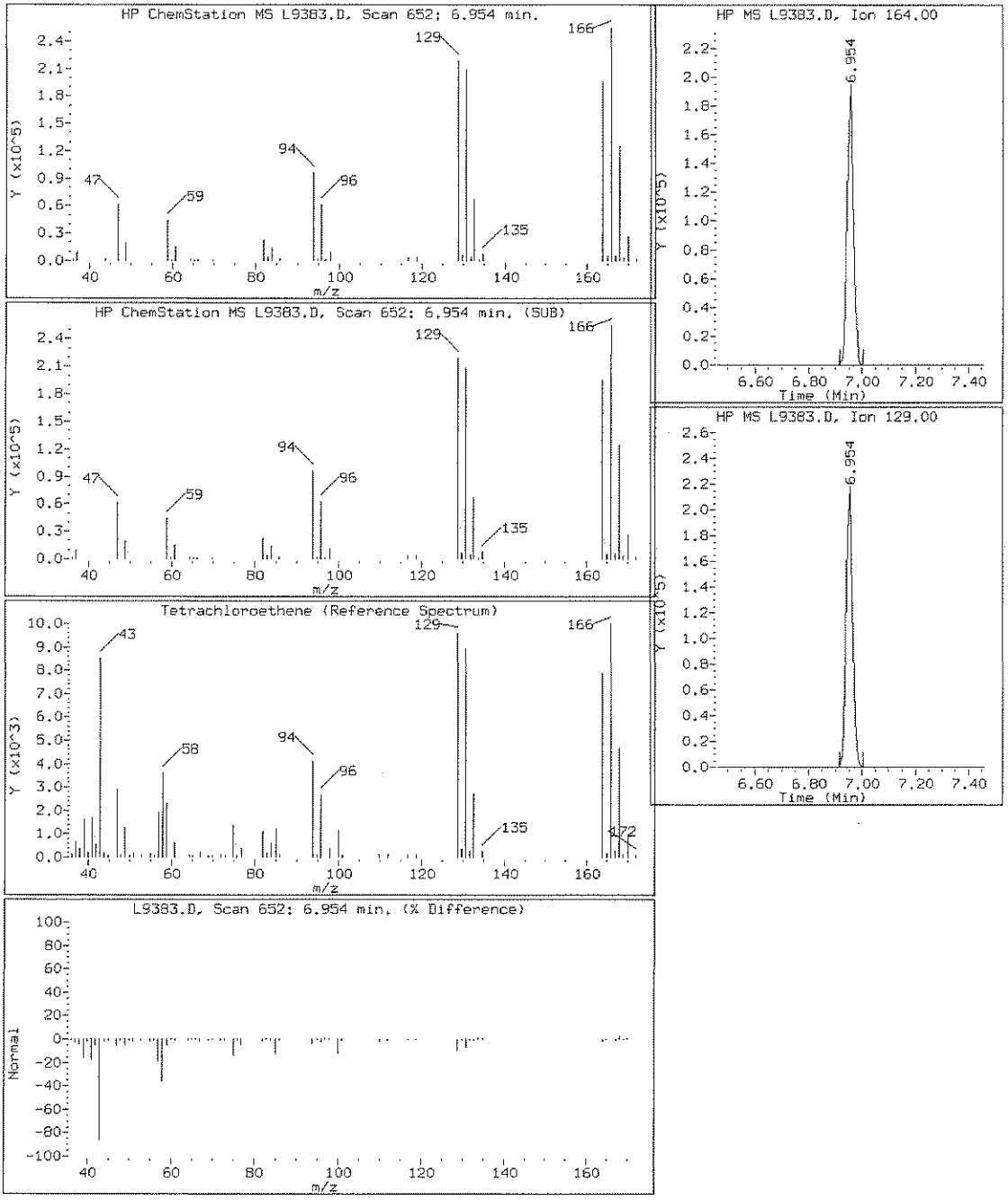
Data File: L9383.D

Date: 02-AUG-2007 15:33

Client ID: A-12 Instrument: msl.i

Sample Info: 220-2277-B-7 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: ME-7 | Lab Sample ID: 220-2277-8

Matrix: Water | Lab File ID: L9378.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/02/2007 13:22

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8356 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	0.88	J	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	41		5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\Files\chem\VOA\msl.i\L079370.b\L9378.D
Lab Smp Id:
Inj Date : 02-AUG-2007 13:22 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-B-8
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\target1_ct\Files\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 60
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONVOAMGR1

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

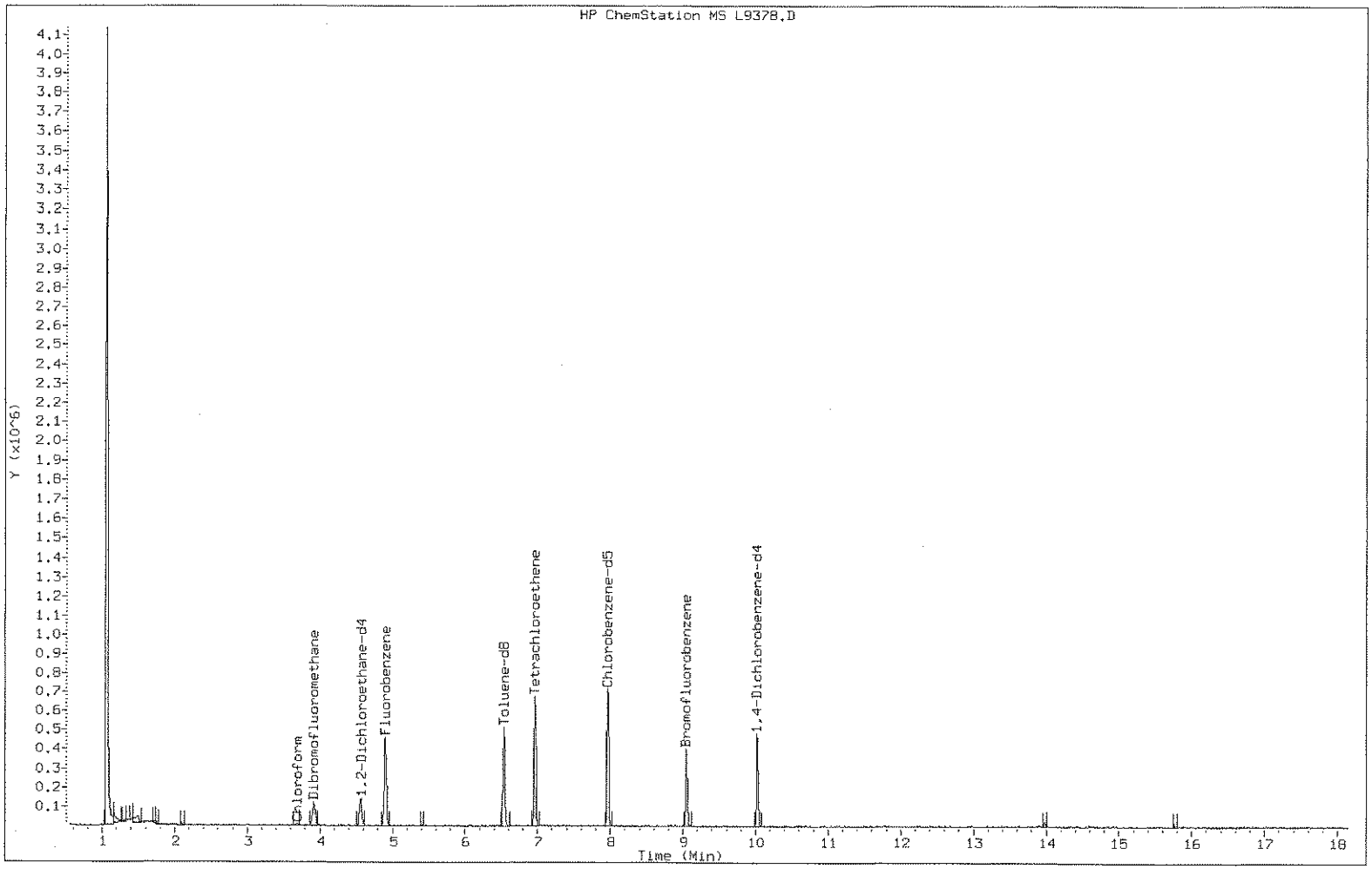
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.901	4.901	(1.000)	444846	25.0000	
	38 Chloroform	83	3.700	3.701	(0.755)	8195	0.88163	0.9
$	41 Dibromofluoromethane	111	3.917	3.927	(0.799)	107365	18.7081	19
$	55 1,2-Dichloroethane-d4	65	4.556	4.567	(0.930)	134676	19.2481	19
*	75 Chlorobenzene-d5	117	7.970	7.961	(1.000)	434210	25.0000	
$	77 Toluene-d8	98	6.534	6.535	(0.820)	317459	20.5615	20
	80 Tetrachloroethene	164	6.957	6.958	(0.873)	147844	40.5250	40
*	95 1,4-Dichlorobenzene-d4	152	10.027	10.018	(1.000)	119729	25.0000	
$	125 Bromofluorobenzene	95	9.043	9.043	(0.902)	129711	28.8779	29

Data File: L9378.D

Date: 02-AUG-2007 13:22

Client ID: ME-7 Instrument: msl.i

Sample Info: 220-2277-B-8 Operator: D. HUMBERT



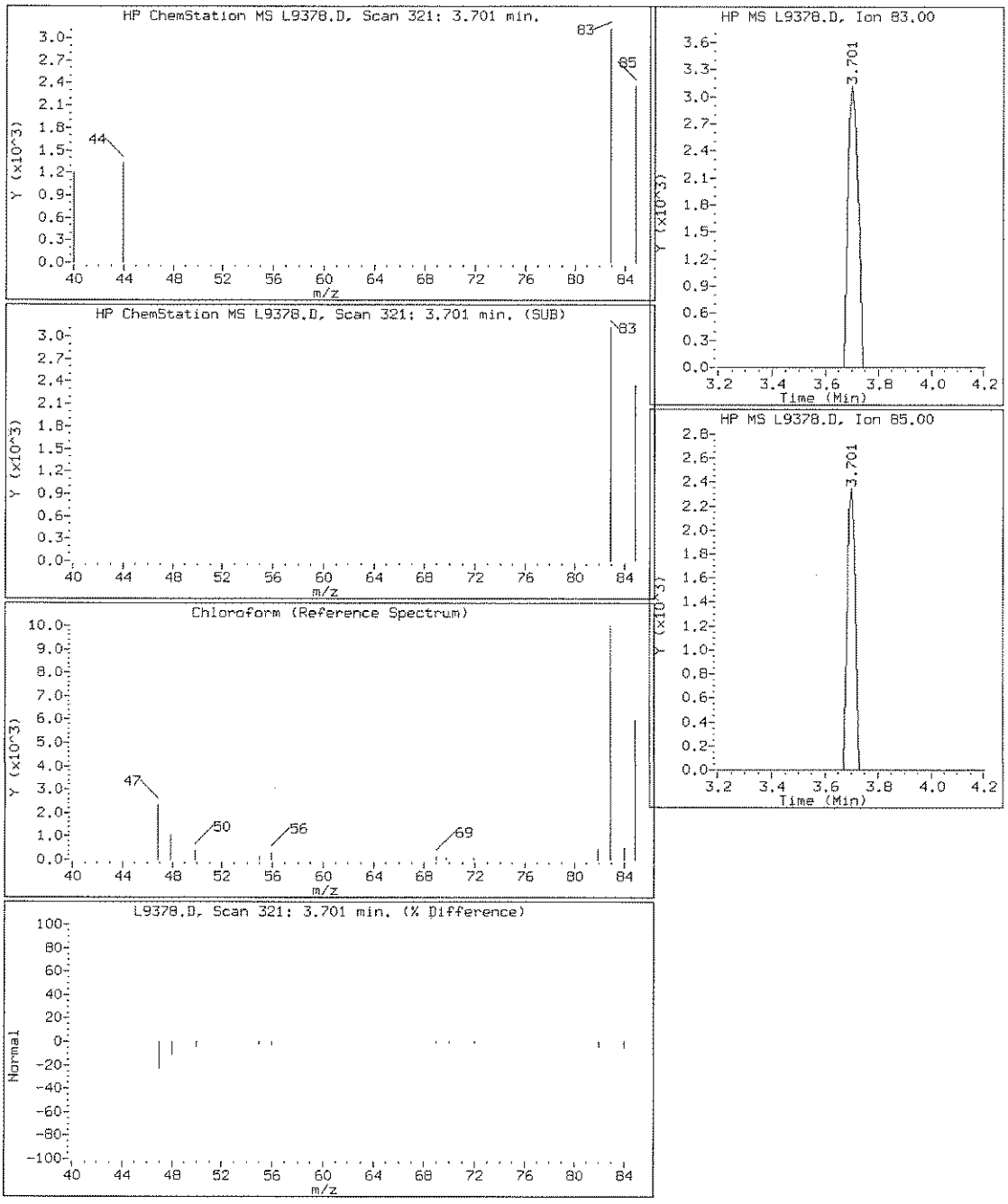
Data File: L9378.D

Date: 02-AUG-2007 13:22

Client ID: ME-7 Instrument: msl.i

Sample Info: 220-2277-B-8 Operator: D. HUMBERT

38 Chloroform



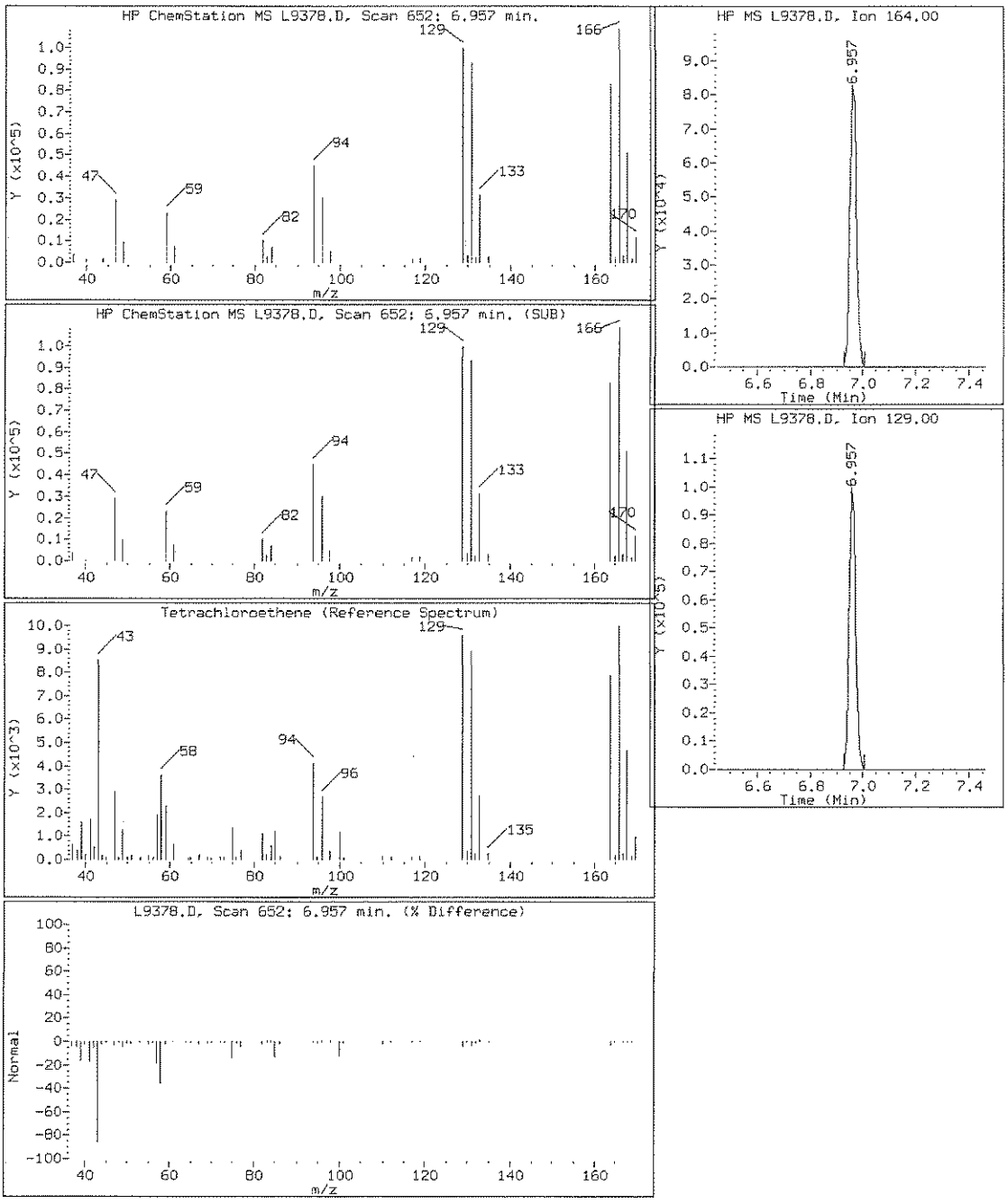
Data File: L9378.D

Date: 02-AUG-2007 13:22

Client ID: ME-7 Instrument: msl.i

Sample Info: 220-2277-B-8 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: ME-11 Lab Sample ID: 220-2277-9

Matrix: Water Lab File ID: L9387.D

Analysis Method: 8260B Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) Date Analyzed: 08/02/2007 17:13

Level: (low/med) Low Dilution Factor: 2

GC Column/ID: RTX-VMS 0.18 (mm) Soil Aliquot:

Soil Extract Vol.: % Moisture:

Analy. Batch No.: 8356 Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	20	U	20	2.8
71-43-2	Benzene	10	U	10	0.80
75-27-4	Bromodichloromethane	10	U	10	0.80
75-25-2	Bromoform	10	U	10	1.6
74-83-9	Bromomethane	10	U	10	2.4
78-93-3	2-Butanone (MEK)	20	U	20	2.4
75-15-0	Carbon disulfide	10	U	10	1.8
56-23-5	Carbon tetrachloride	10	U	10	2.0
108-90-7	Chlorobenzene	10	U	10	0.80
75-00-3	Chloroethane	10	U	10	1.6
67-66-3	Chloroform	10	U	10	1.4
74-87-3	Chloromethane	10	U	10	1.0
124-48-1	Dibromochloromethane	10	U	10	1.0
75-34-3	1,1-Dichloroethane	10	U	10	1.2
107-06-2	1,2-Dichloroethane	10	U	10	1.2
75-35-4	1,1-Dichloroethene	10	U	10	1.4
78-87-5	1,2-Dichloropropane	10	U	10	1.8
10061-01-5	cis-1,3-Dichloropropene	10	U	10	1.0
10061-02-6	trans-1,3-Dichloropropene	10	U	10	1.6
100-41-4	Ethylbenzene	10	U	10	2.0
591-78-6	2-Hexanone	20	U	20	1.6
75-09-2	Methylene Chloride	10	U	10	0.80
108-10-1	4-Methyl-2-pentanone (MIBK)	20	U	20	1.4
100-42-5	Styrene	10	U	10	1.0
79-34-5	1,1,2,2-Tetrachloroethane	10	U	10	0.80
127-18-4	Tetrachloroethene	290		10	1.0
108-88-3	Toluene	10	U	10	0.60
71-55-6	1,1,1-Trichloroethane	10	U	10	0.80
79-00-5	1,1,2-Trichloroethane	10	U	10	1.2
79-01-6	Trichloroethene	10	U	10	1.4
75-01-4	Vinyl chloride	10	U	10	1.6
1330-20-7	Xylenes, Total	10	U	10	2.0
156-59-2	cis-1,2-Dichloroethene	10	U	10	1.2
156-60-5	trans-1,2-Dichloroethene	10	U	10	1.0

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\Files\chem\VOA\msl.i\L079370.b\L9387.D
Lab Smp Id:
Inj Date : 02-AUG-2007 17:13 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-A-9
Misc Info : : ;;; ; 8260 ; 2 ; LLW
Comment :
Method : \\target1_ct\Files\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 69
Dil Factor: 2.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONVOAMGR1

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	2.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

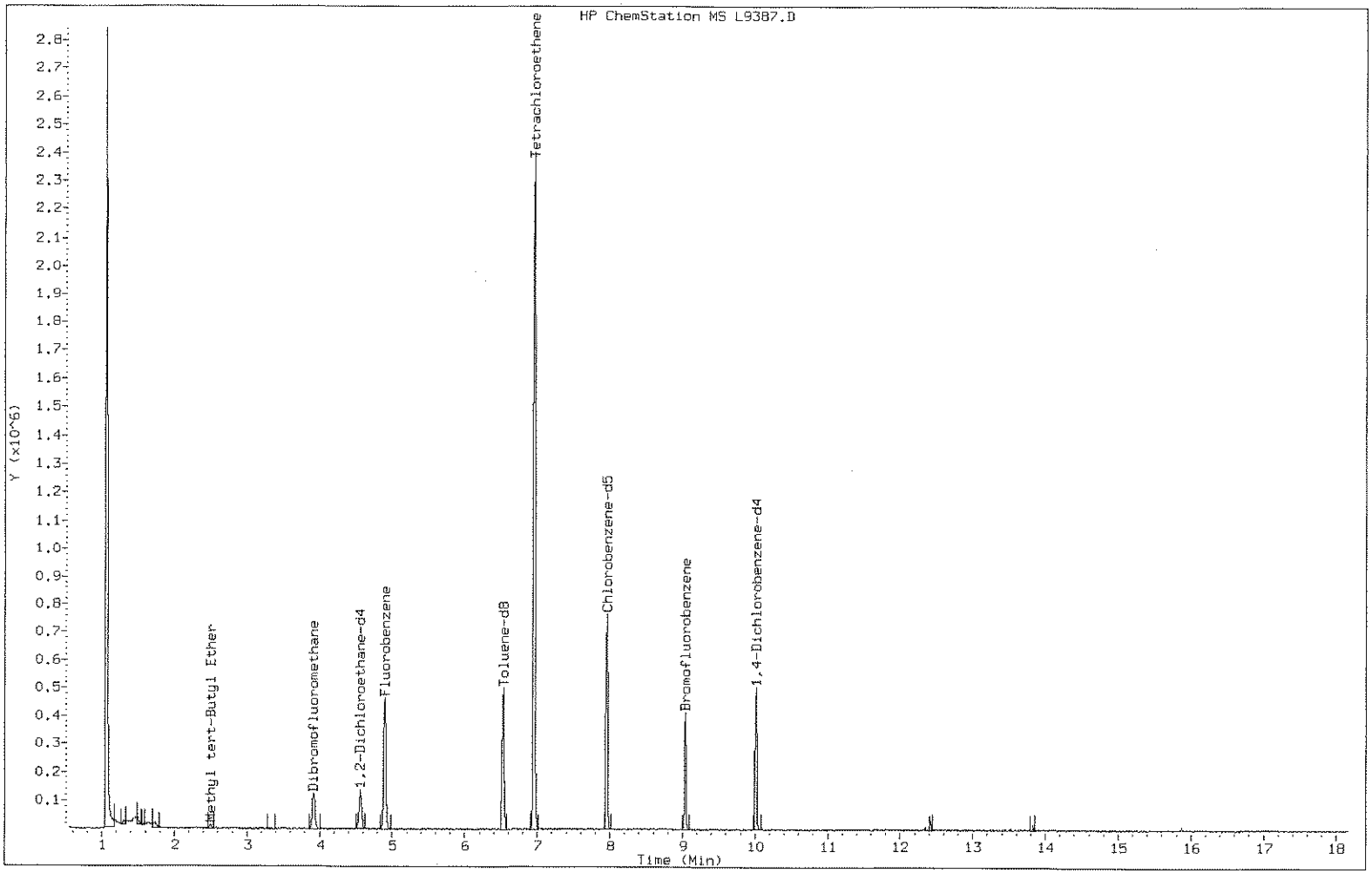
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.898	4.901	(1.000)	458179	25.0000	
	24 Methyl tert-Butyl Ether	73	2.498	2.491	(0.510)	15922	0.83701	2
$	41 Dibromofluoromethane	111	3.924	3.927	(0.801)	111214	18.8149	19
$	55 1,2-Dichloroethane-d4	65	4.564	4.567	(0.932)	131030	18.1820	18
*	75 Chlorobenzene-d5	117	7.958	7.961	(1.000)	441225	25.0000	
$	77 Toluene-d8	98	6.532	6.535	(0.821)	315271	20.0952	20
	80 Tetrachloroethene	164	6.955	6.958	(0.874)	538031	145.133	290
*	95 1,4-Dichlorobenzene-d4	152	10.015	10.018	(1.000)	123236	25.0000	
$	125 Bromofluorobenzene	95	9.041	9.043	(0.903)	128060	27.6990	28

Data File: L9387.D

Date: 02-AUG-2007 17:13

Client ID: ME-11 Instrument: msl.i

Sample Info: 220-2277-A-9 Operator: D. HUMBERT



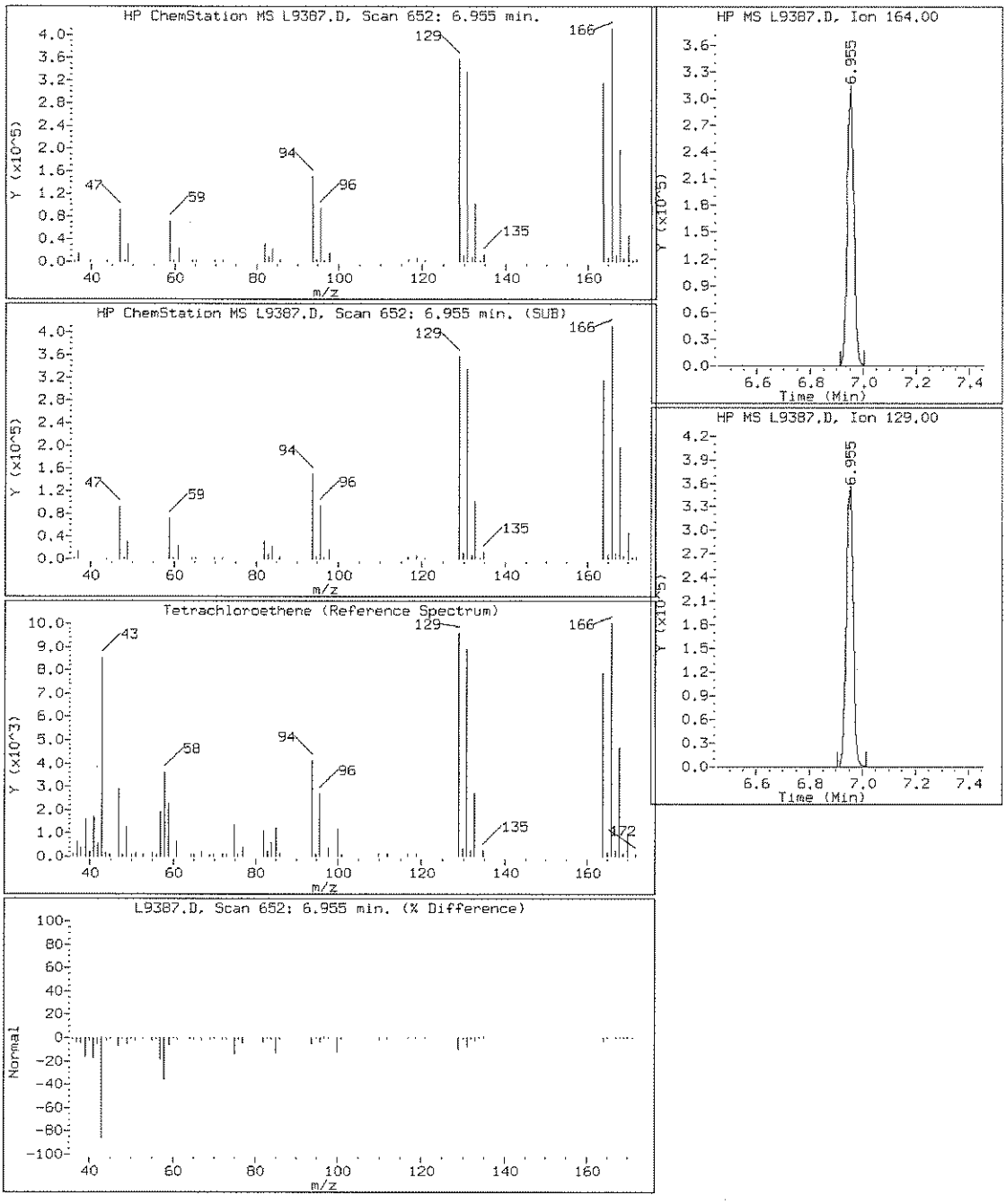
Data File: L9387.D

Date: 02-AUG-2007 17:13

Client ID: ME-11 Instrument: msl.i

Sample Info: 220-2277-A-9 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: FB0723 | Lab Sample ID: 220-2277-10

Matrix: Water | Lab File ID: L9326.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/01/2007 02:59

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8307 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	0.54	J	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	5.0	U	5.0	0.50
108-88-3	Toluene	0.40	J	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079310.b\L9326.D
Lab Smp Id: 220-2277-B-10 Client Smp ID: FB0723
Inj Date : 01-AUG-2007 02:59 MS Autotune Date: 26-APR-2004 14:21
Operator : D. GAYDA Inst ID: msl.i
Smp Info : 220-2277-B-10
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079310.b\L8260BNW.m
Meth Date : 31-Jul-2007 21:07 target Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 17
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

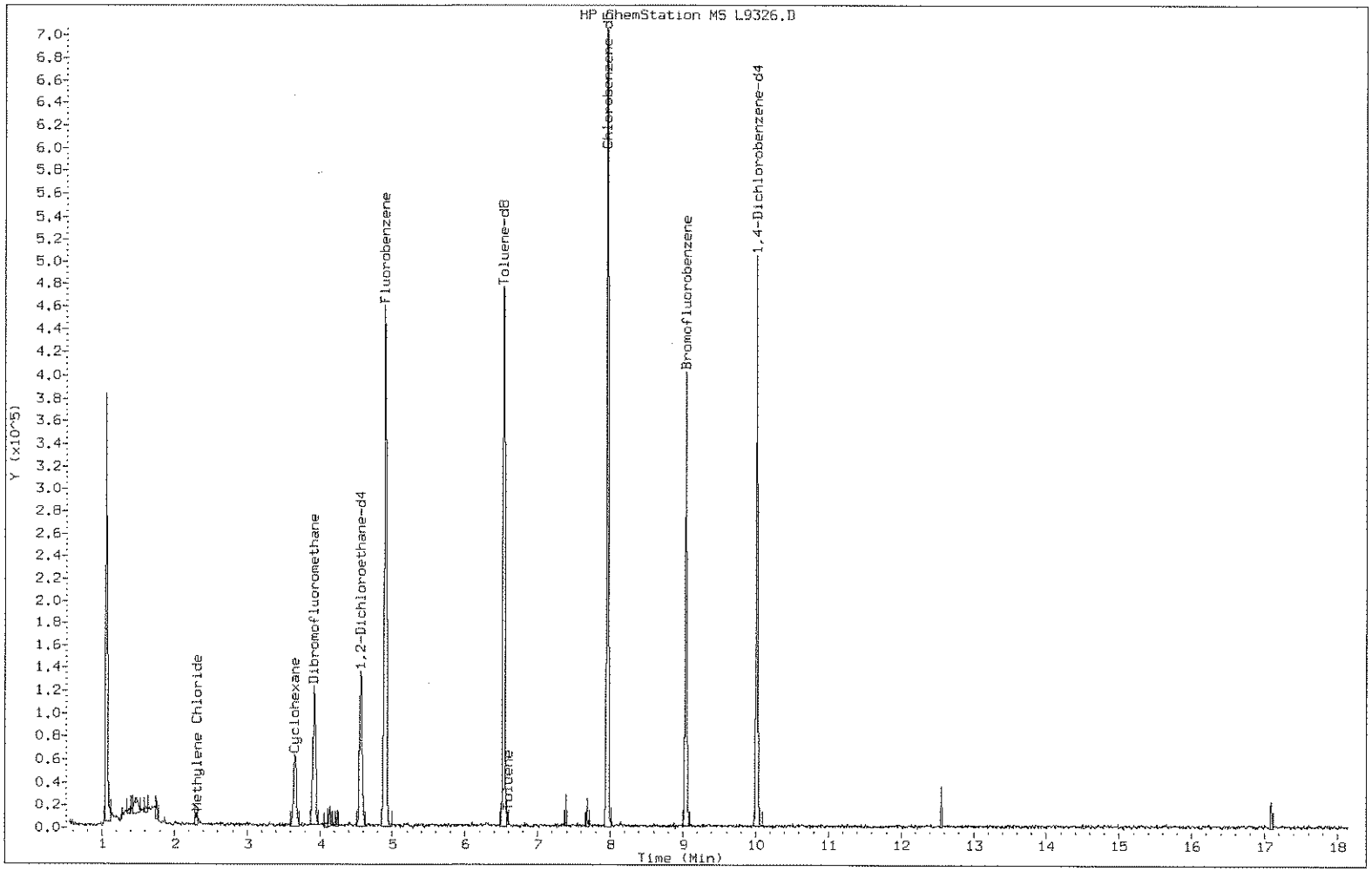
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.891	4.896	(1.000)	451044	25.0000	
	20 Methylene Chloride	84	2.304	2.298	(0.471)	2877	0.54443	0.5
	37 Cyclohexane	84	3.661	3.656	(0.749)	26326	3.99589	4
$	41 Dibromofluoromethane	111	3.917	3.922	(0.801)	109623	18.8391	19
$	55 1,2-Dichloroethane-d4	65	4.557	4.561	(0.932)	125141	17.6395	18
*	75 Chlorobenzene-d5	117	7.961	7.956	(1.000)	437455	25.0000	
	76 Toluene	91	6.584	6.578	(0.827)	7494	0.40232	0.4
$	77 Toluene-d8	98	6.525	6.529	(0.820)	310334	19.9509	20
*	95 1,4-Dichlorobenzene-d4	152	10.017	10.012	(1.000)	124592	25.0000	
$	125 Bromofluorobenzene	95	9.034	9.038	(0.902)	128375	27.4649	27

Data File: L9326.D

Date: 01-AUG-2007 02:59

Client ID: FB0723 Instrument: msl.i

Sample Info: 220-2277-B-10 Operator: D. GAYDA



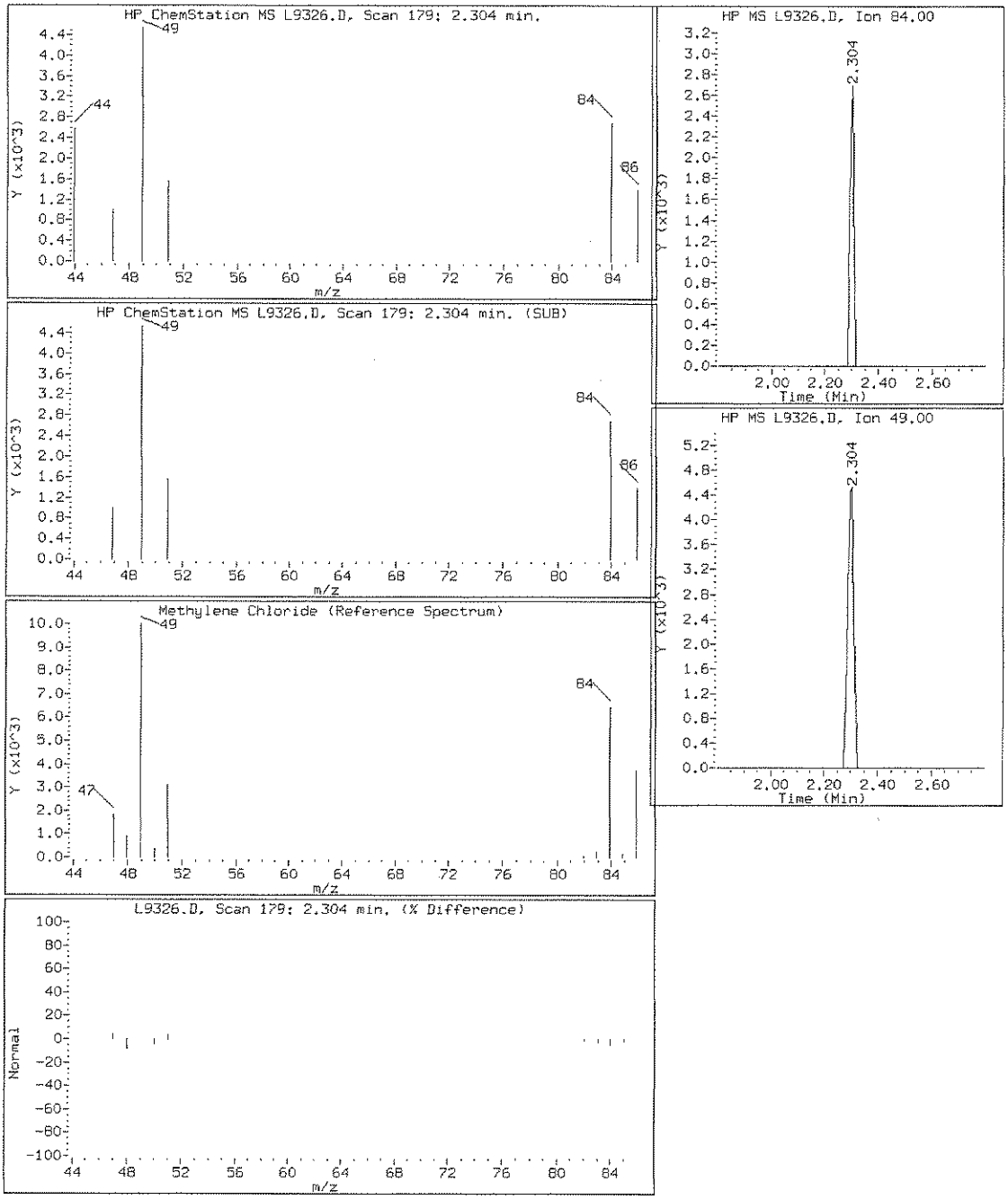
Data File: L9326.D

Date: 01-AUG-2007 02:59

Client ID: FB0723 Instrument: msl.i

Sample Info: 220-2277-B-10 Operator: D. GAYDA

20 Methylene Chloride



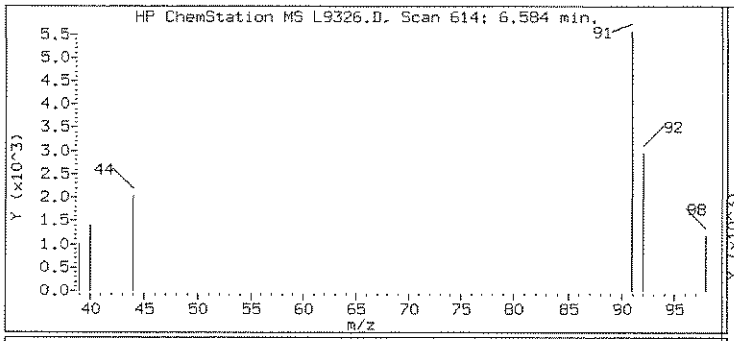
Data File: L9326.D

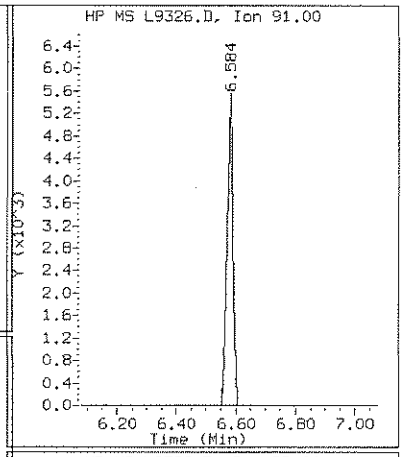
Date: 01-AUG-2007 02:59

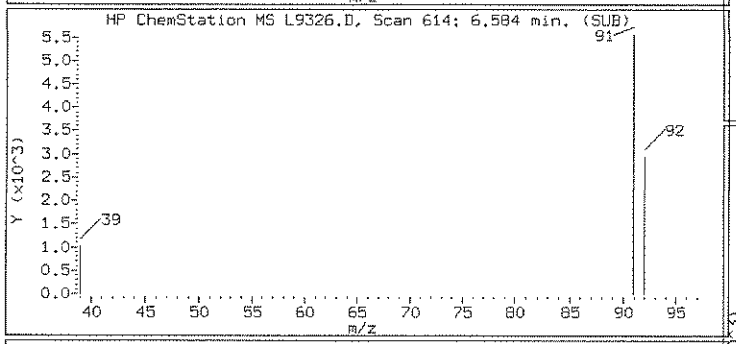
Client ID: FB0723 Instrument: msl.i

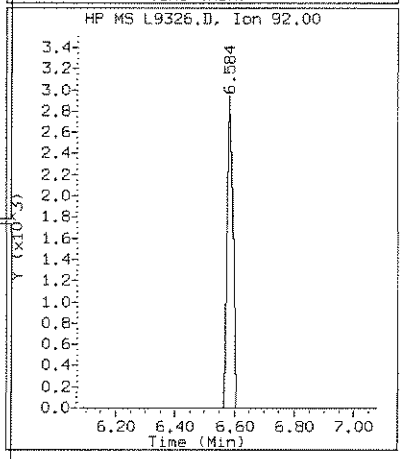
Sample Info: 220-2277-B-10 Operator: D. GAYDA

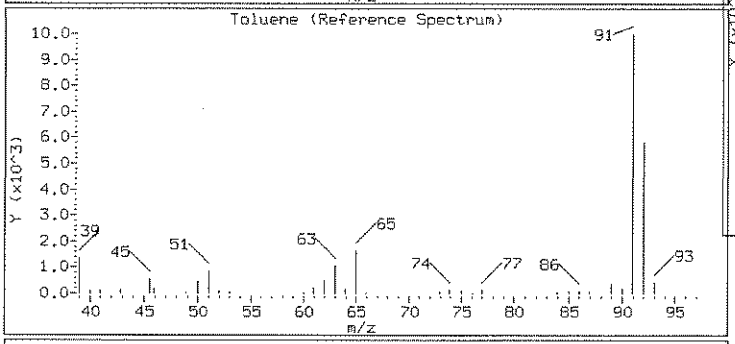
76 Toluene

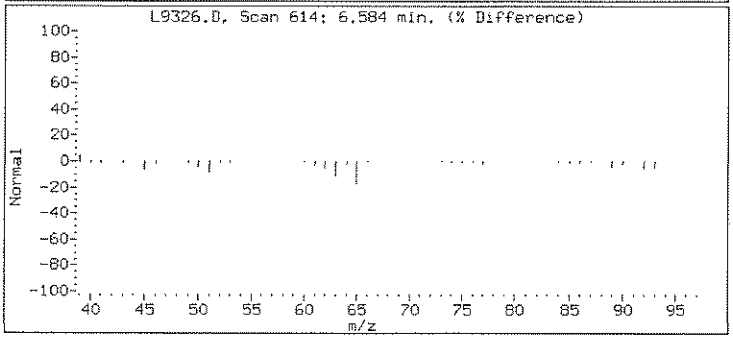












FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: TRIP BLANK | Lab Sample ID: 220-2277-11

Matrix: Water | Lab File ID: L9327.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/01/2007 03:23

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8307 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	3.1	J	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	3.6	J	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	2.8	J	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	5.0	U	5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079310.b\L9327.D
Lab Smp Id: 220-2277-B-11 Client Smp ID: TRIP BLANK
Inj Date : 01-AUG-2007 03:23 MS Autotune Date: 26-APR-2004 14:21
Operator : D. GAYDA Inst ID: msl.i
Smp Info : 220-2277-B-11
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079310.b\L8260BNW.m
Meth Date : 31-Jul-2007 21:07 target Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 18
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.896	4.896	(1.000)	446792	25.0000	
	20 Methylene Chloride	84	2.299	2.298	(0.470)	14454	2.76122	3
	21 Acetone	43	2.328	2.328	(0.476)	8755	3.08983	3
$	41 Dibromofluoromethane	111	3.922	3.922	(0.801)	107151	18.5895	18
	45 2-Butanone	43	4.090	4.069	(0.835)	14791	3.59609	4
$	55 1,2-Dichloroethane-d4	65	4.562	4.561	(0.932)	135038	19.2157	19
*	75 Chlorobenzene-d5	117	7.956	7.956	(1.000)	417702	25.0000	
$	77 Toluene-d8	98	6.530	6.529	(0.821)	320783	21.5979	22
*	95 1,4-Dichlorobenzene-d4	152	10.013	10.012	(1.000)	117531	25.0000	
$	125 Bromofluorobenzene	95	9.039	9.038	(0.903)	133119	30.1908	30

Data File: L9327.D

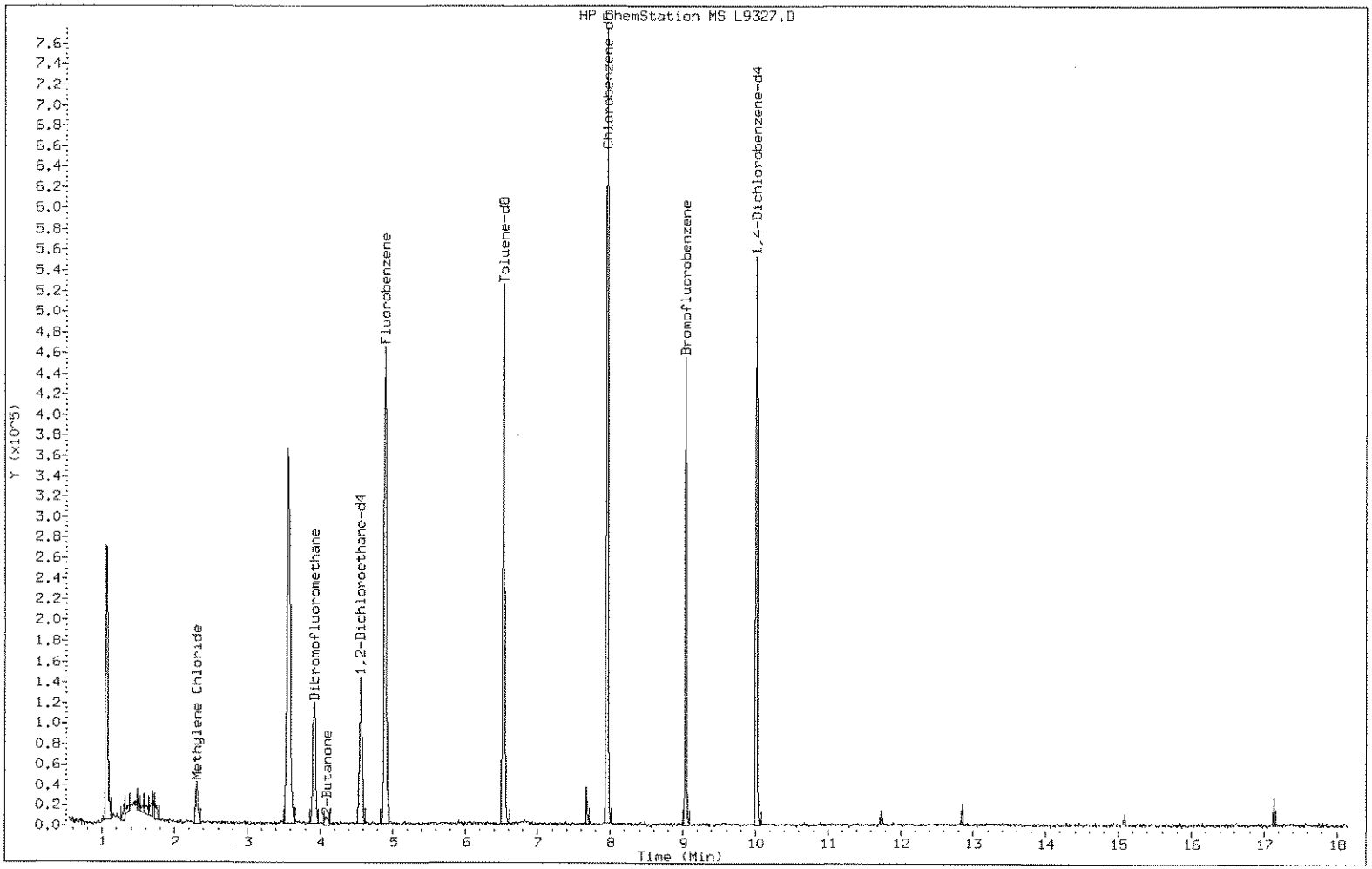
Date: 01-AUG-2007 03:23

Client ID: TRIP BLANK

Instrument: msl.i

Sample Info: 220-2277-B-11

Operator: D. GAYDA



Data File: L9327.D

Date: 01-AUG-2007 03:23

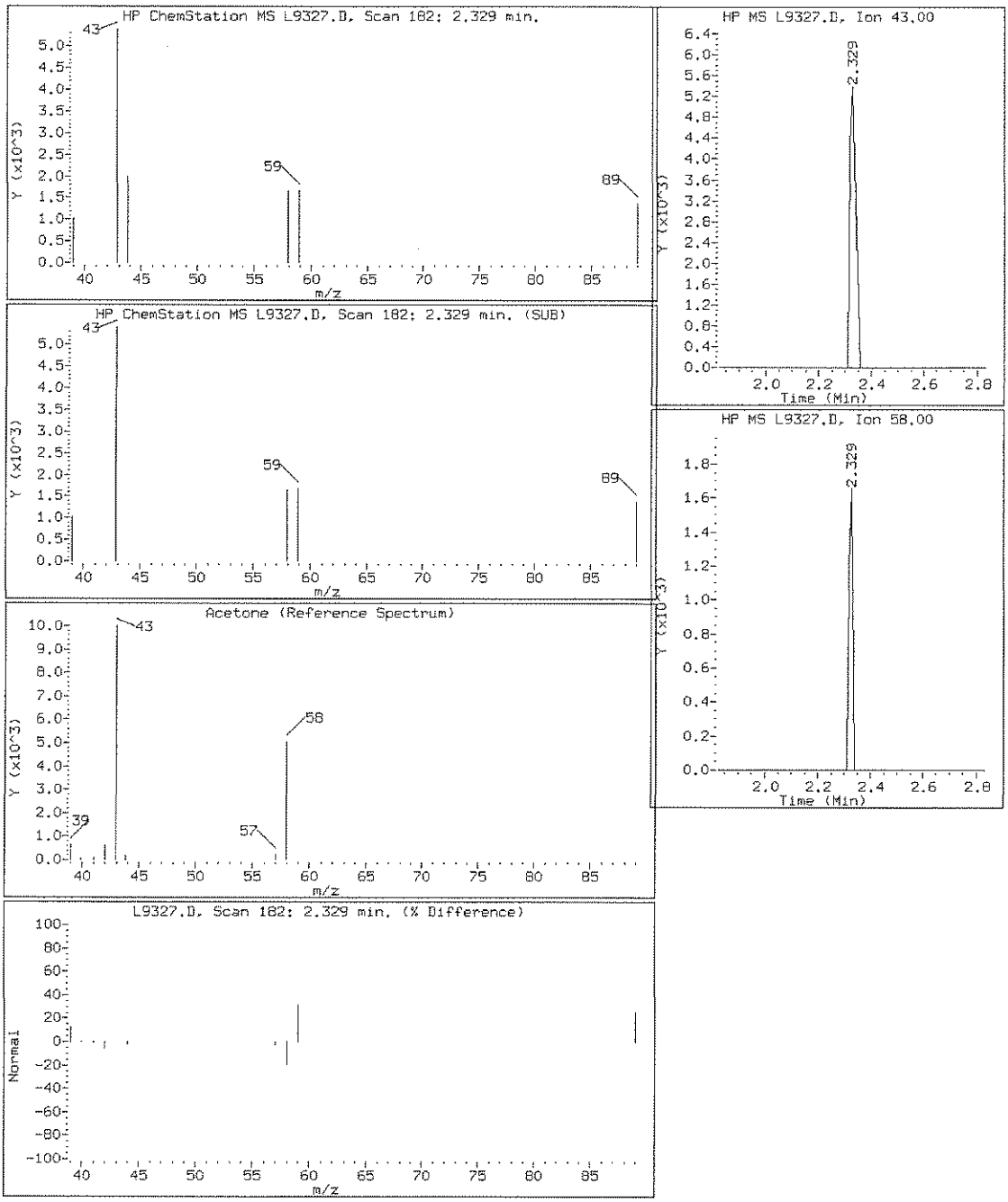
Client ID: TRIP BLANK

Instrument: ms1.i

Sample Info: 220-2277-B-11

Operator: D. GAYDA

21 Acetone



Data File: L9327.D

Date: 01-AUG-2007 03:23

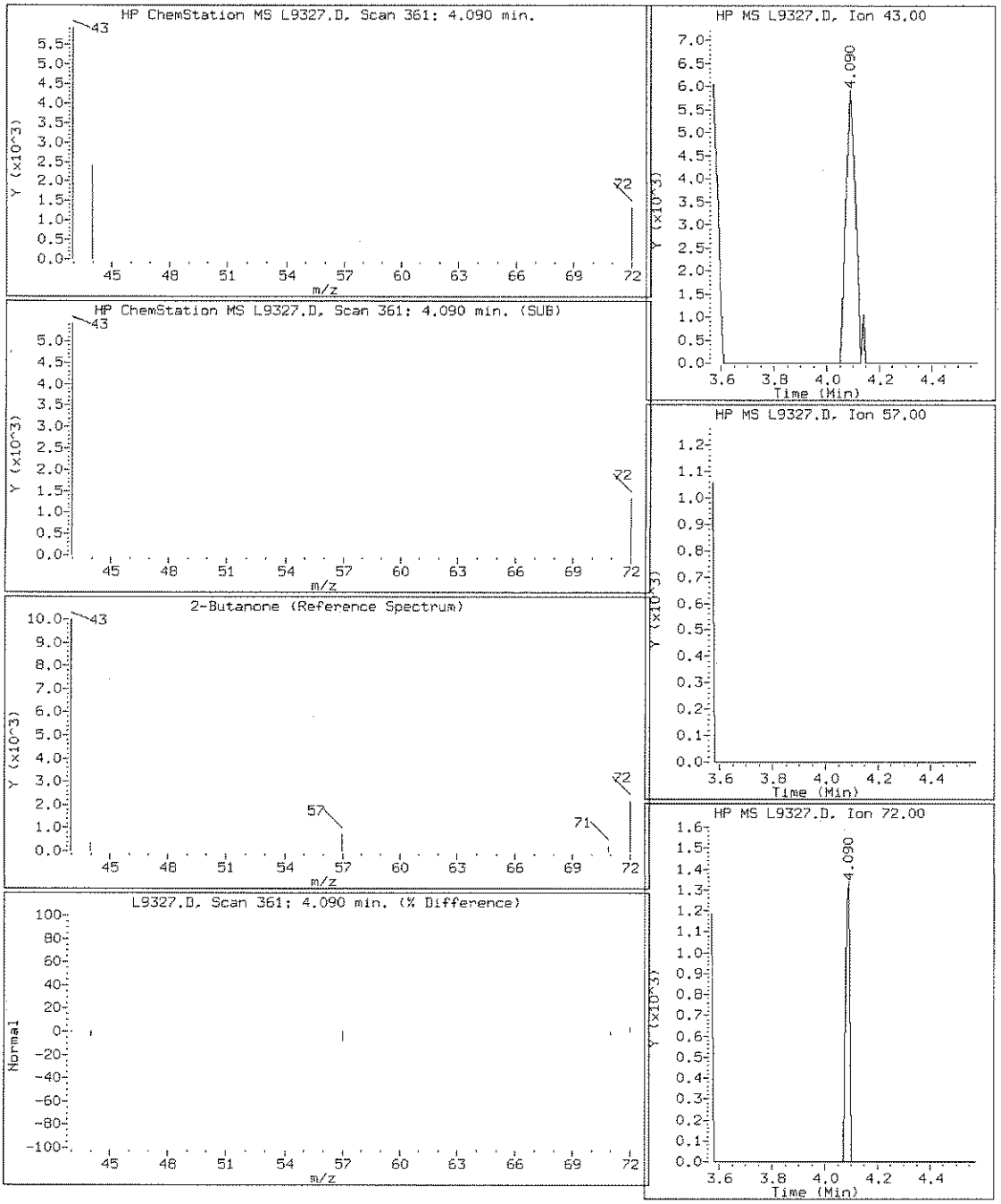
Client ID: TRIP BLANK

Instrument: msl.i

Sample Info: 220-2277-B-11

Operator: D. GAYDA

45 2-Butanone



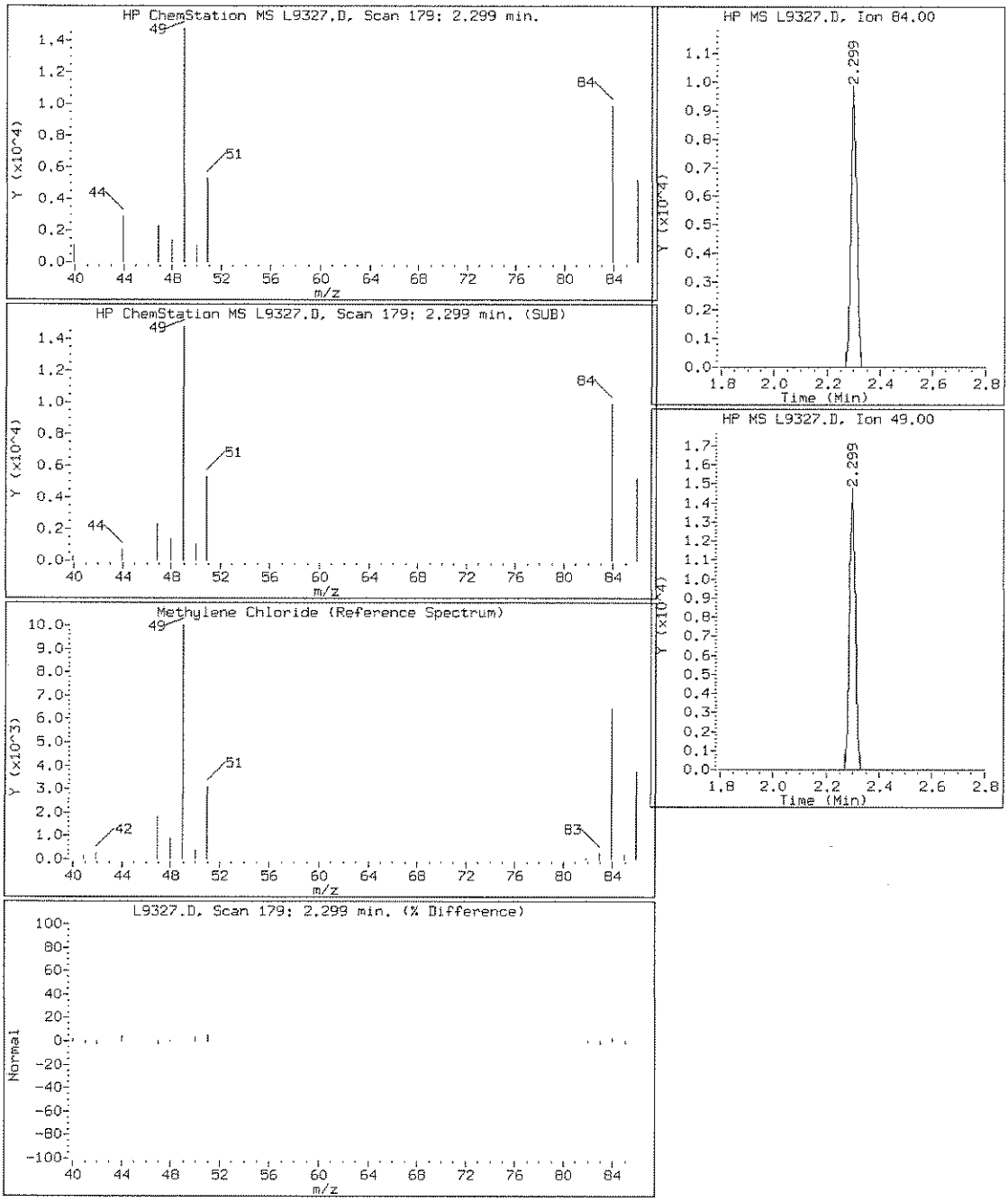
Data File: L9327.D

Date: 01-AUG-2007 03:23

Client ID: TRIP BLANK Instrument: msl.i

Sample Info: 220-2277-B-11 Operator: D. GAYDA

20 Methylene Chloride



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-16 | Lab Sample ID: 220-2277-12

Matrix: Water | Lab File ID: L9379.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/02/2007 13:47

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8356 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	5.0	U	5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079370.b\L9379.D
Lab Smp Id: 220-2277-A-12 Client Smp ID: A-16
Inj Date : 02-AUG-2007 13:47 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-A-12
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 61
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

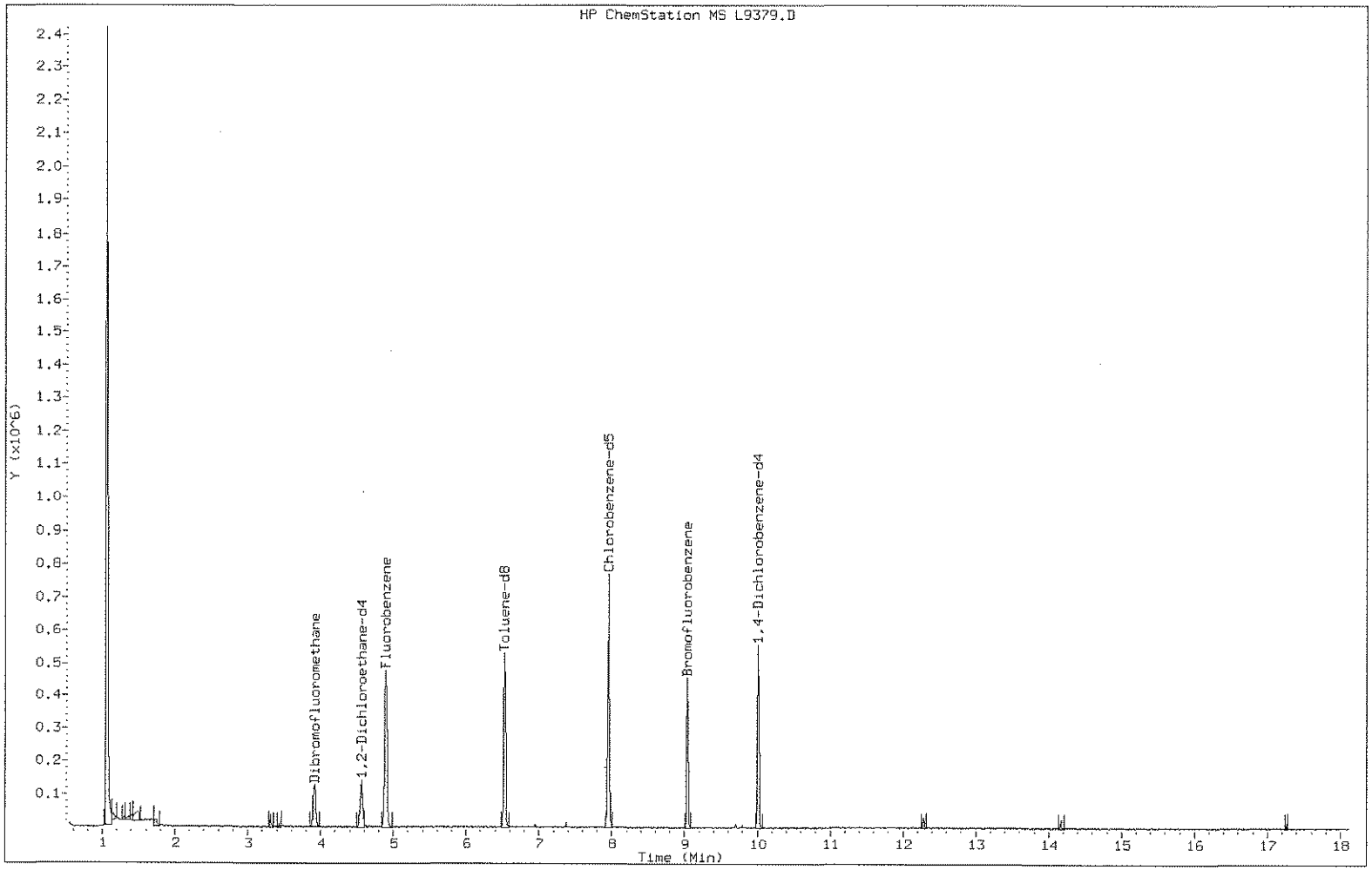
Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1 Fluorobenzene	96	4.892	4.901	(1.000)	452863	25.0000	
$ 41 Dibromofluoromethane	111	3.918	3.927	(0.801)	113968	19.5071	20
$ 55 1,2-Dichloroethane-d4	65	4.557	4.567	(0.932)	134626	18.9003	19
* 75 Chlorobenzene-d5	117	7.952	7.961	(1.000)	434522	25.0000	
$ 77 Toluene-d8	98	6.525	6.535	(0.821)	323944	20.9665	21
80 Tetrachloroethene	164	6.958	6.958	(0.875)	1436	0.39333	0.4
* 95 1,4-Dichlorobenzene-d4	152	10.008	10.018	(1.000)	122772	25.0000	
$ 125 Bromofluorobenzene	95	9.034	9.043	(0.903)	132876	28.8493	29

Data File: L9379.D

Date: 02-AUG-2007 13:47

Client ID: A-16 Instrument: msl.i

Sample Info: 220-2277-A-12 Operator: D. HUMBERT



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-17 | Lab Sample ID: 220-2277-13

Matrix: Water | Lab File ID: L9351.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/01/2007 14:45

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8321 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	0.79	J	5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079341.b\L9351.D
Lab Smp Id: 220-2277-A-13 Client Smp ID: A-17
Inj Date : 01-AUG-2007 14:45 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-A-13
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079341.b\L8260BNW.m
Meth Date : 02-Aug-2007 09:12 target Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 38
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

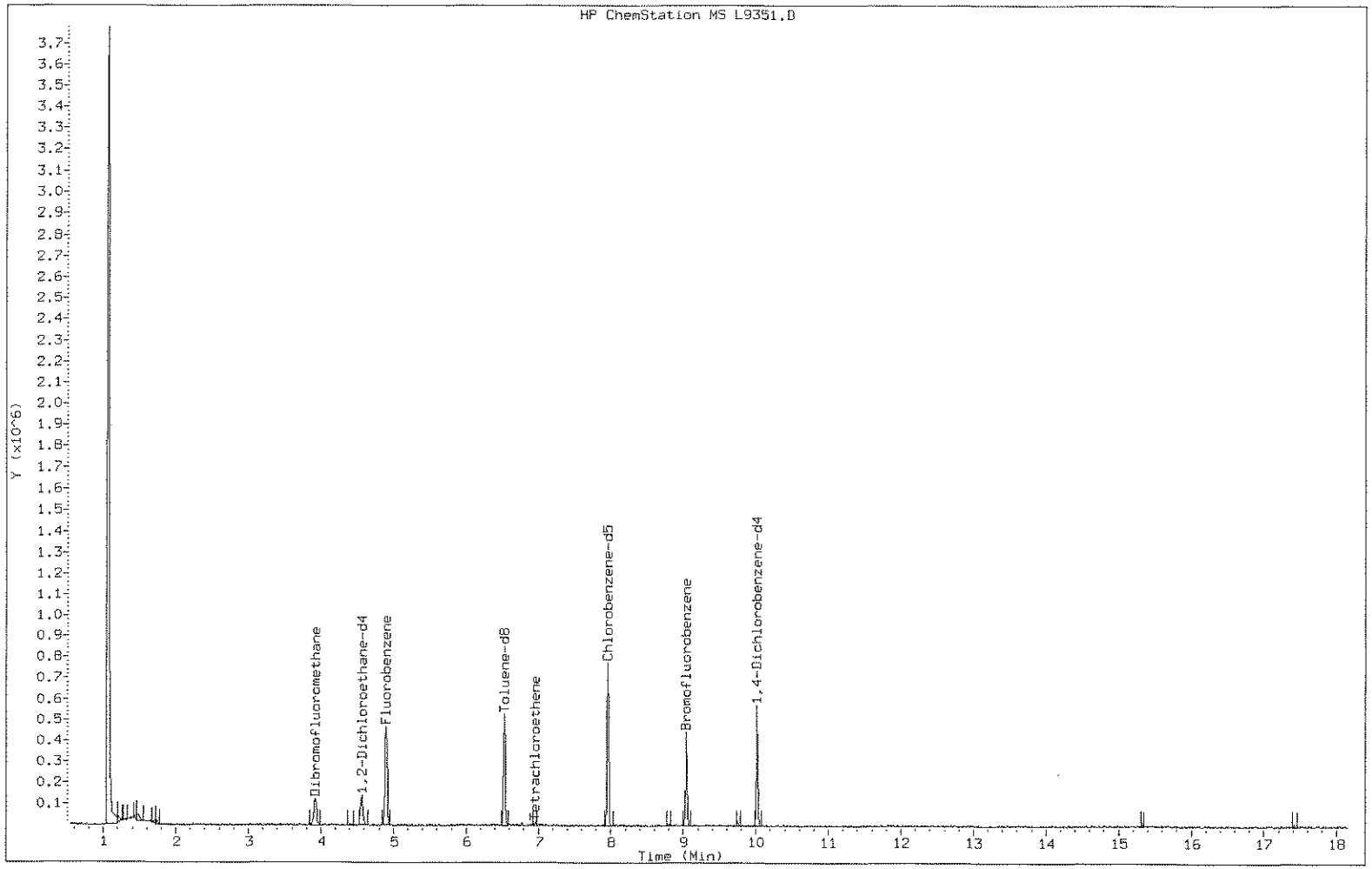
Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1 Fluorobenzene	96	4.897	4.908	(1.000)	448065	25.0000	
$ 41 Dibromofluoromethane	111	3.923	3.924	(0.801)	107045	18.5184	18
$ 55 1,2-Dichloroethane-d4	65	4.562	4.563	(0.932)	127328	18.0671	18
* 75 Chlorobenzene-d5	117	7.957	7.968	(1.000)	425019	25.0000	
$ 77 Toluene-d8	98	6.530	6.541	(0.821)	314344	20.8000	21
80 Tetrachloroethene	164	6.953	6.964	(0.874)	2806	0.78578	0.8
* 95 1,4-Dichlorobenzene-d4	152	10.013	10.014	(1.000)	126762	25.0000	
$ 125 Bromofluorobenzene	95	9.039	9.040	(0.903)	126685	26.6394	27

Data File: L9351.D

Date: 01-AUG-2007 14:45

Client ID: A-17 Instrument: msl.i

Sample Info: 220-2277-A-13 Operator: D. HUMBERT



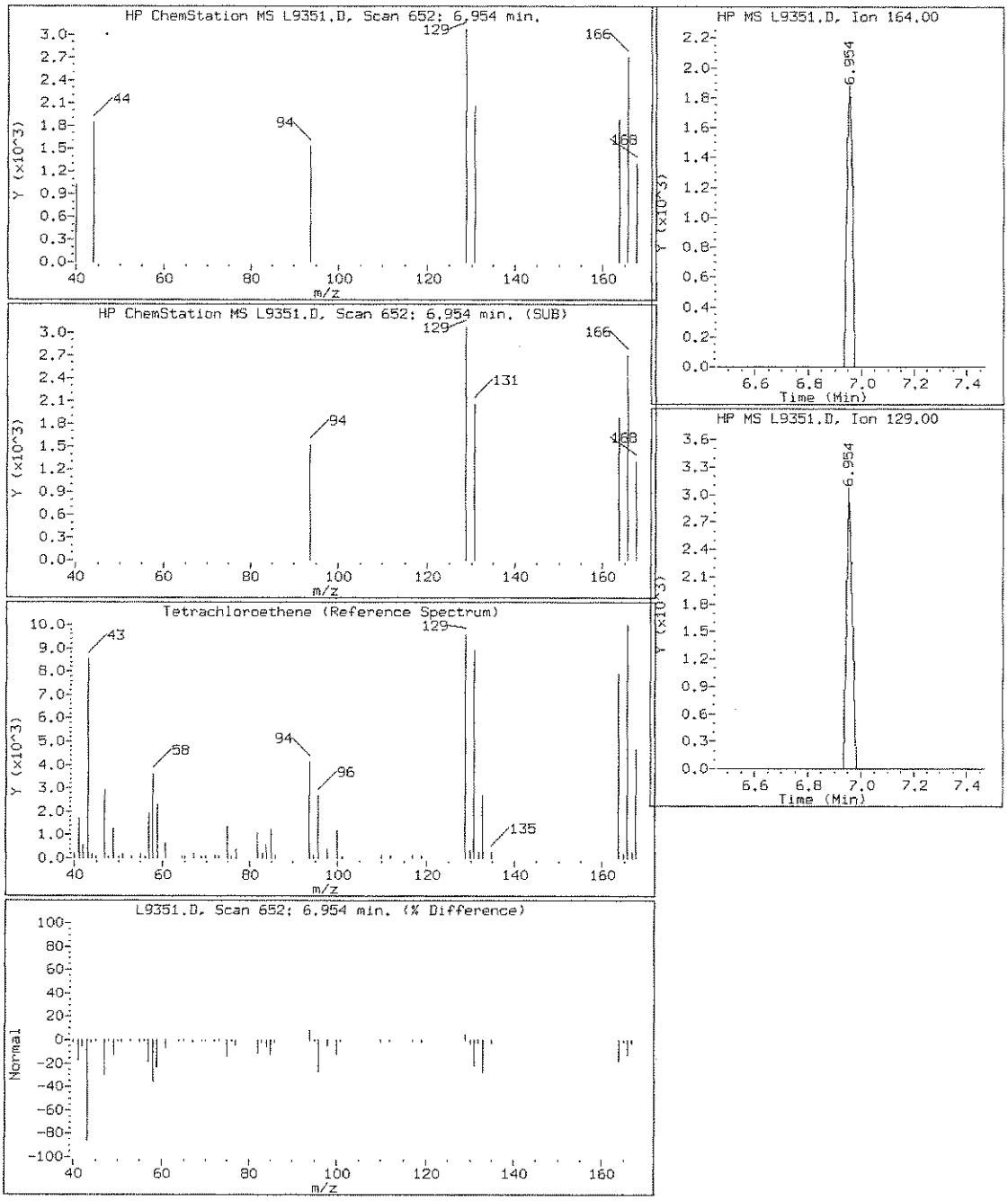
Data File: L9351.D

Date: 01-AUG-2007 14:45

Client ID: A-17 Instrument: msl.i

Sample Info: 220-2277-A-13 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: MW-109 | Lab Sample ID: 220-2277-14

Matrix: Water | Lab File ID: L9352.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/01/2007 15:10

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8321 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	2.6	J	5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079341.b\L9352.D
Lab Smp Id: 220-2277-A-14 Client Smp ID: MW-109
Inj Date : 01-AUG-2007 15:10 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-A-14
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079341.b\L8260BNW.m
Meth Date : 02-Aug-2007 09:12 target Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 39
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

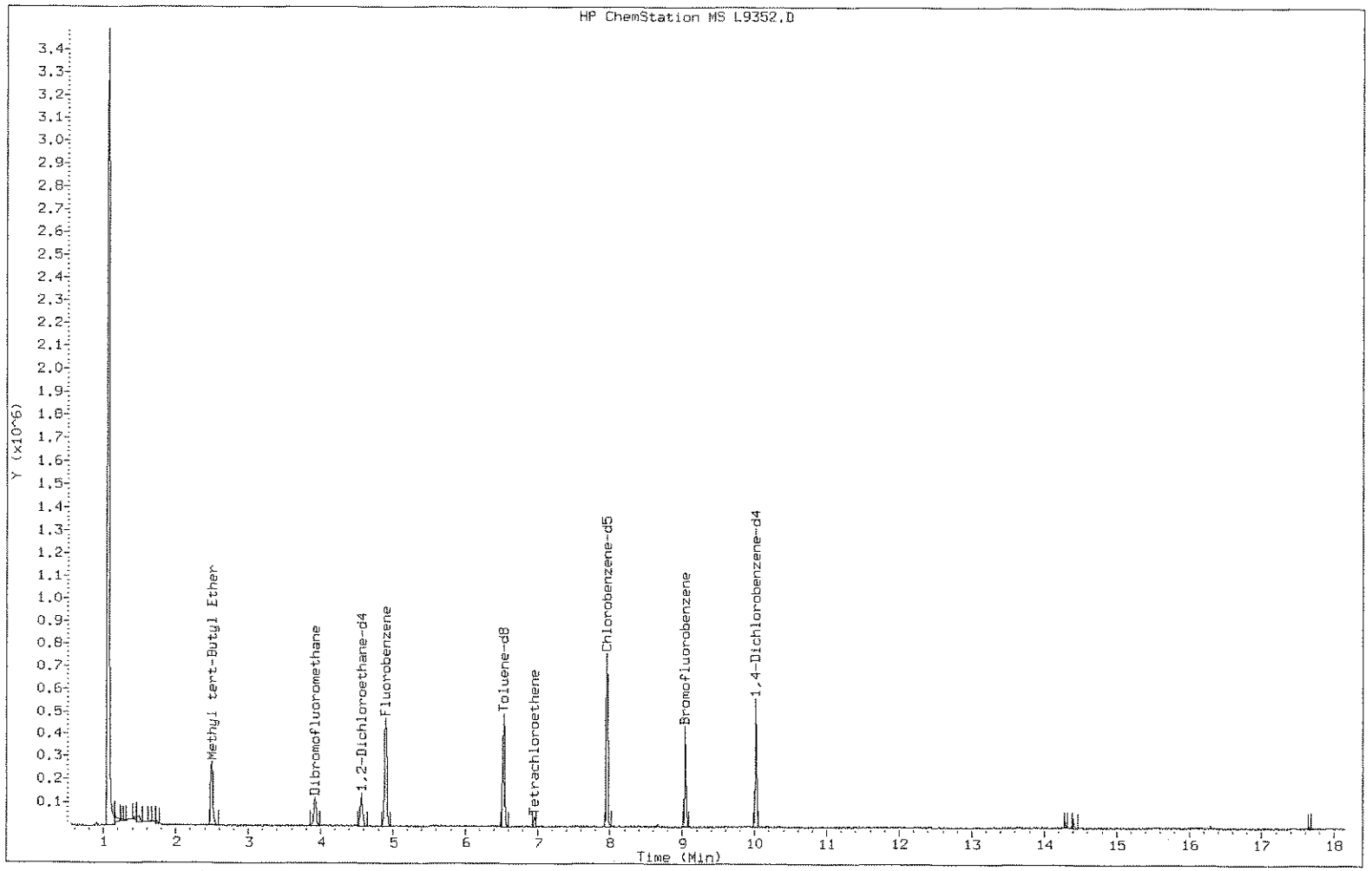
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.897	4.908	(1.000)	451047	25.0000	
	24 Methyl tert-Butyl Ether	73	2.496	2.497	(0.510)	308079	16.4516	16
$	41 Dibromofluoromethane	111	3.923	3.924	(0.801)	104029	17.8776	18
$	55 1,2-Dichloroethane-d4	65	4.562	4.563	(0.932)	125864	17.7413	18
*	75 Chlorobenzene-d5	117	7.957	7.968	(1.000)	422636	25.0000	
$	77 Toluene-d8	98	6.530	6.541	(0.821)	306050	20.3654	20
	80 Tetrachloroethene	164	6.953	6.964	(0.874)	9174	2.58352	2
*	95 1,4-Dichlorobenzene-d4	152	10.013	10.014	(1.000)	121626	25.0000	
$	125 Bromofluorobenzene	95	9.039	9.040	(0.903)	129893	28.4673	28

Data File: L9352.D

Date: 01-AUG-2007 15:10

Client ID: MW-109 Instrument: msl.i

Sample Info: 220-2277-A-14 Operator: D. HUMBERT



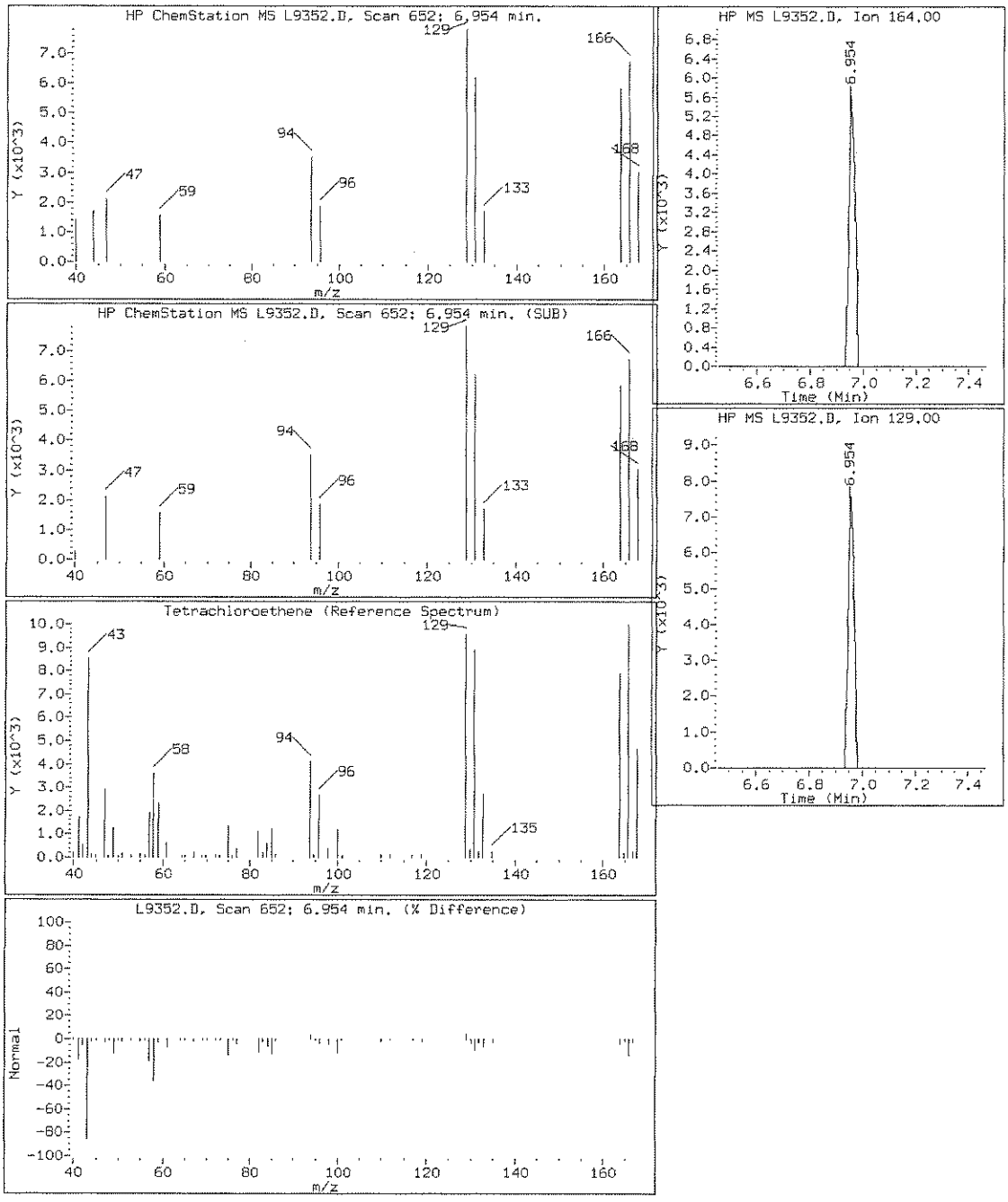
Data File: L9352.D

Date: 01-AUG-2007 15:10

Client ID: MW-109 Instrument: msl.i

Sample Info: 220-2277-A-14 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: MW-110 | Lab Sample ID: 220-2277-15

Matrix: Water | Lab File ID: L9353.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/01/2007 15:35

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8321 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	5.0	U	5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079341.b\L9353.D
Lab Smp Id: 220-2277-B-15 Client Smp ID: MW-110
Inj Date : 01-AUG-2007 15:35 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-B-15
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079341.b\L8260BNW.m
Meth Date : 02-Aug-2007 09:12 target Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 40
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

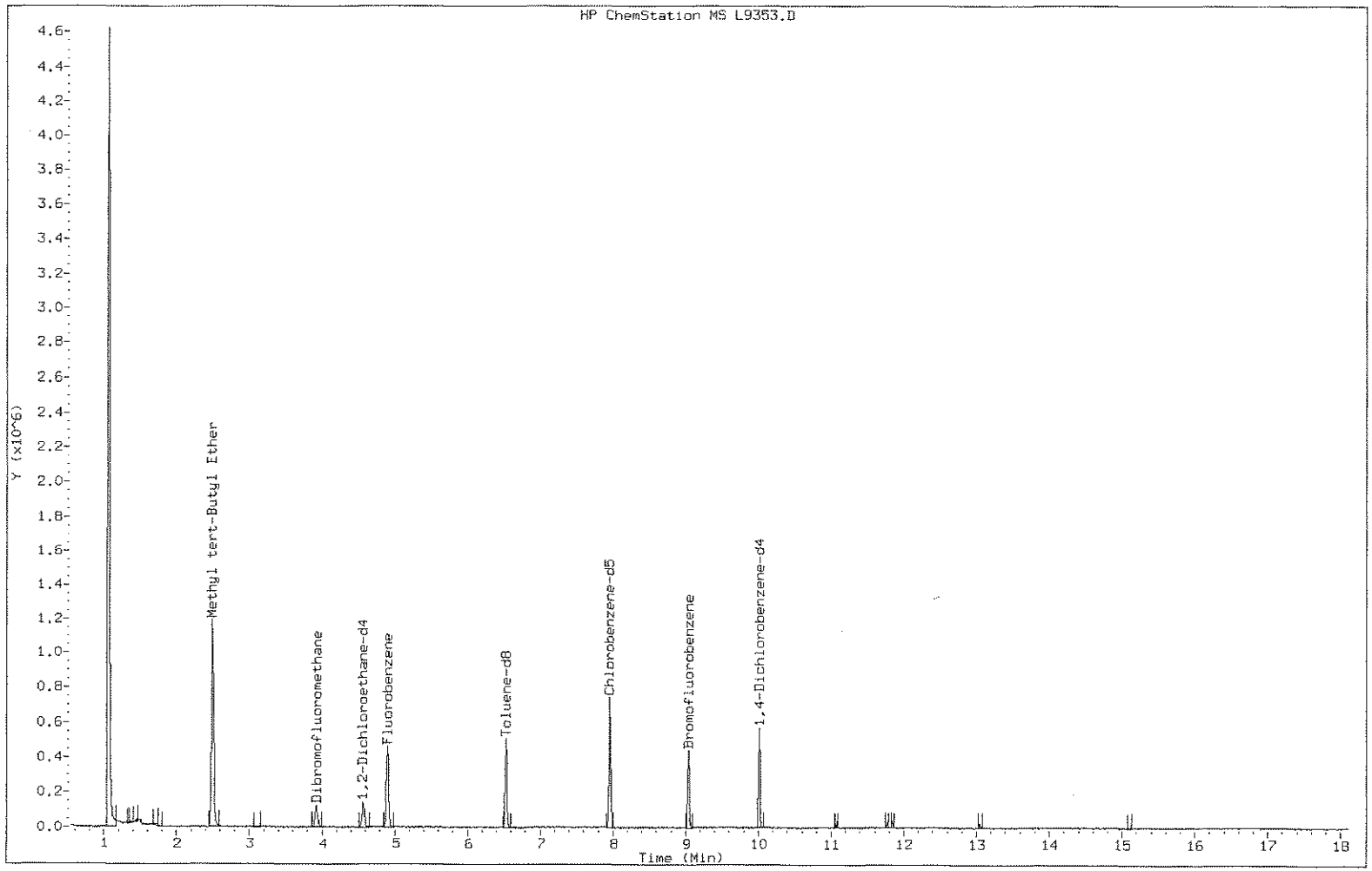
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.896	4.908	(1.000)	432833	25.0000	
	24 Methyl tert-Butyl Ether	73	2.486	2.497	(0.508)	1273466	70.8657	71
$	41 Dibromofluoromethane	111	3.922	3.924	(0.801)	109066	19.5320	20
$	55 1,2-Dichloroethane-d4	65	4.562	4.563	(0.932)	127582	18.7403	19
*	75 Chlorobenzene-d5	117	7.956	7.968	(1.000)	411260	25.0000	
$	77 Toluene-d8	98	6.530	6.541	(0.821)	309564	21.1691	21
*	95 1,4-Dichlorobenzene-d4	152	10.013	10.014	(1.000)	123577	25.0000	
$	125 Bromofluorobenzene	95	9.039	9.040	(0.903)	129679	27.9718	28

Data File: L9353.D

Date: 01-AUG-2007 15:35

Client ID: MW-110 Instrument: msl.i

Sample Info: 220-2277-B-15 Operator: D. HUMBERT



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-18 | Lab Sample ID: 220-2277-16

Matrix: Water | Lab File ID: L9354.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/01/2007 16:00

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8321 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	5.0	U	5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\Files\chem\VOA\msl.i\L079341.b\L9354.D
Lab Smp Id:
Inj Date : 01-AUG-2007 16:00 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-A-16
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\target1_ct\Files\chem\VOA\msl.i\L079341.b\L8260BNW.m
Meth Date : 01-Aug-2007 10:17 ctvoa Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 41
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONVOAMGR1

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

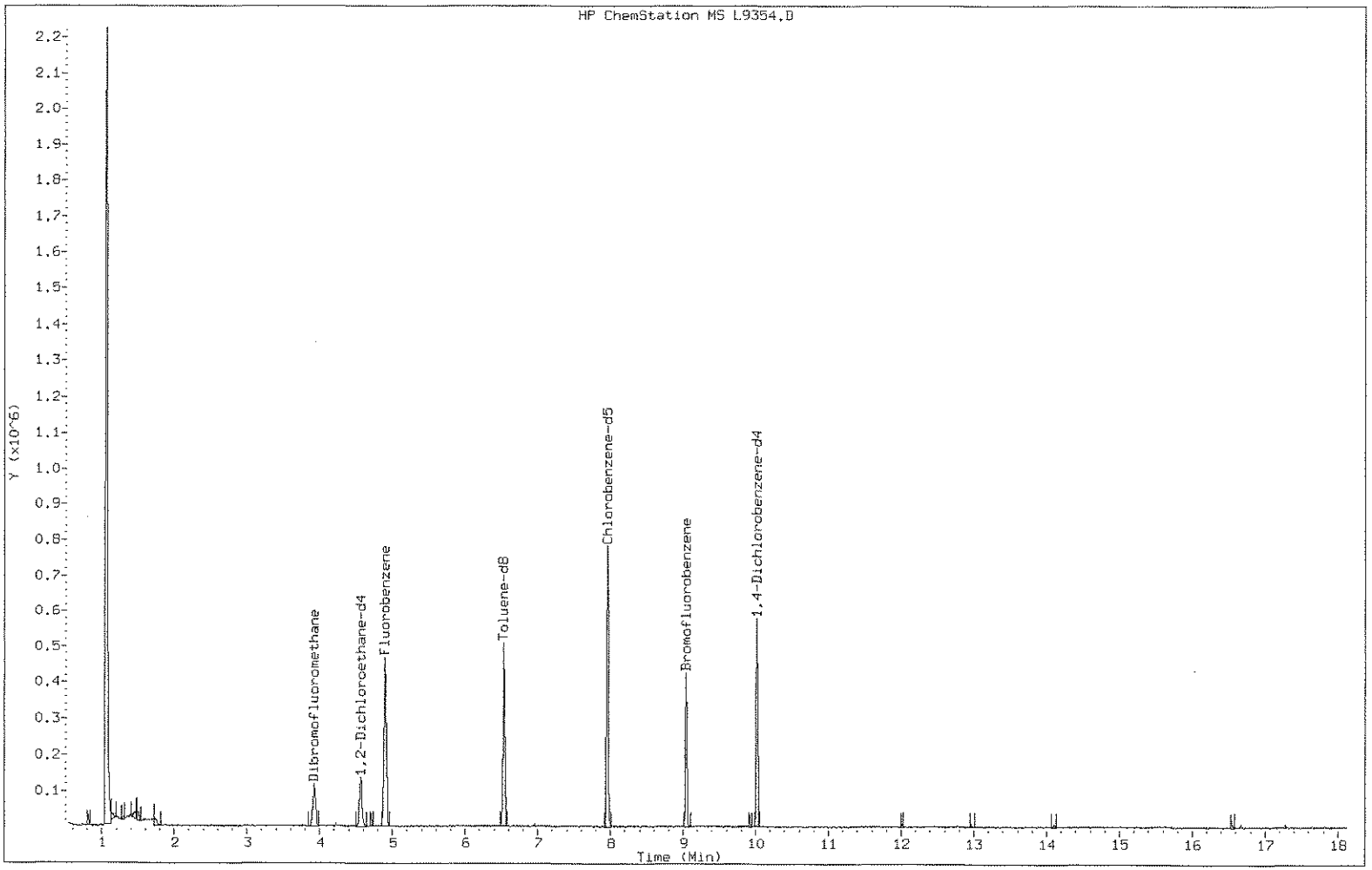
Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1 Fluorobenzene	96	4.896	4.908	(1.000)	443177	25.0000	
$ 41 Dibromofluoromethane	111	3.922	3.924	(0.801)	104159	18.2178	18
$ 55 1,2-Dichloroethane-d4	65	4.562	4.563	(0.932)	123167	17.6695	18
* 75 Chlorobenzene-d5	117	7.956	7.968	(1.000)	435439	25.0000	
$ 77 Toluene-d8	98	6.530	6.541	(0.821)	302263	19.5220	20
* 95 1,4-Dichlorobenzene-d4	152	10.013	10.014	(1.000)	124354	25.0000	
$ 125 Bromofluorobenzene	95	9.039	9.040	(0.903)	124807	26.7527	27

Data File: L9354.D

Date: 01-AUG-2007 16:00

Client ID: A-18 Instrument: msl.i

Sample Info: 220-2277-A-16 Operator: D. HUMBERT



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-14 | Lab Sample ID: 220-2277-17

Matrix: Water | Lab File ID: L9385.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/02/2007 16:23

Level: (low/med) Low | Dilution Factor: 4

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8356 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	40	U	40	5.6
71-43-2	Benzene	20	U	20	1.6
75-27-4	Bromodichloromethane	20	U	20	1.6
75-25-2	Bromoform	20	U	20	3.2
74-83-9	Bromomethane	20	U	20	4.8
78-93-3	2-Butanone (MEK)	40	U	40	4.8
75-15-0	Carbon disulfide	20	U	20	3.6
56-23-5	Carbon tetrachloride	20	U	20	4.0
108-90-7	Chlorobenzene	20	U	20	1.6
75-00-3	Chloroethane	20	U	20	3.2
67-66-3	Chloroform	20	U	20	2.8
74-87-3	Chloromethane	20	U	20	2.0
124-48-1	Dibromochloromethane	20	U	20	2.0
75-34-3	1,1-Dichloroethane	20	U	20	2.4
107-06-2	1,2-Dichloroethane	20	U	20	2.4
75-35-4	1,1-Dichloroethene	20	U	20	2.8
78-87-5	1,2-Dichloropropane	20	U	20	3.6
10061-01-5	cis-1,3-Dichloropropene	20	U	20	2.0
10061-02-6	trans-1,3-Dichloropropene	20	U	20	3.2
100-41-4	Ethylbenzene	20	U	20	4.0
591-78-6	2-Hexanone	40	U	40	3.2
75-09-2	Methylene Chloride	20	U	20	1.6
108-10-1	4-Methyl-2-pentanone (MIBK)	40	U	40	2.8
100-42-5	Styrene	20	U	20	2.0
79-34-5	1,1,2,2-Tetrachloroethane	20	U	20	1.6
127-18-4	Tetrachloroethene	470		20	2.0
108-88-3	Toluene	20	U	20	1.2
71-55-6	1,1,1-Trichloroethane	20	U	20	1.6
79-00-5	1,1,2-Trichloroethane	20	U	20	2.4
79-01-6	Trichloroethene	20	U	20	2.8
75-01-4	Vinyl chloride	20	U	20	3.2
1330-20-7	Xylenes, Total	20	U	20	4.0
156-59-2	cis-1,2-Dichloroethene	20	U	20	2.4
156-60-5	trans-1,2-Dichloroethene	20	U	20	2.0

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079370.b\L9385.D
Lab Smp Id: 220-2277-B-17 Client Smp ID: A-14
Inj Date : 02-AUG-2007 16:23 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-B-17
Misc Info : : ;;; ; 8260 ; 4 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 67
Dil Factor: 4.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	4.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

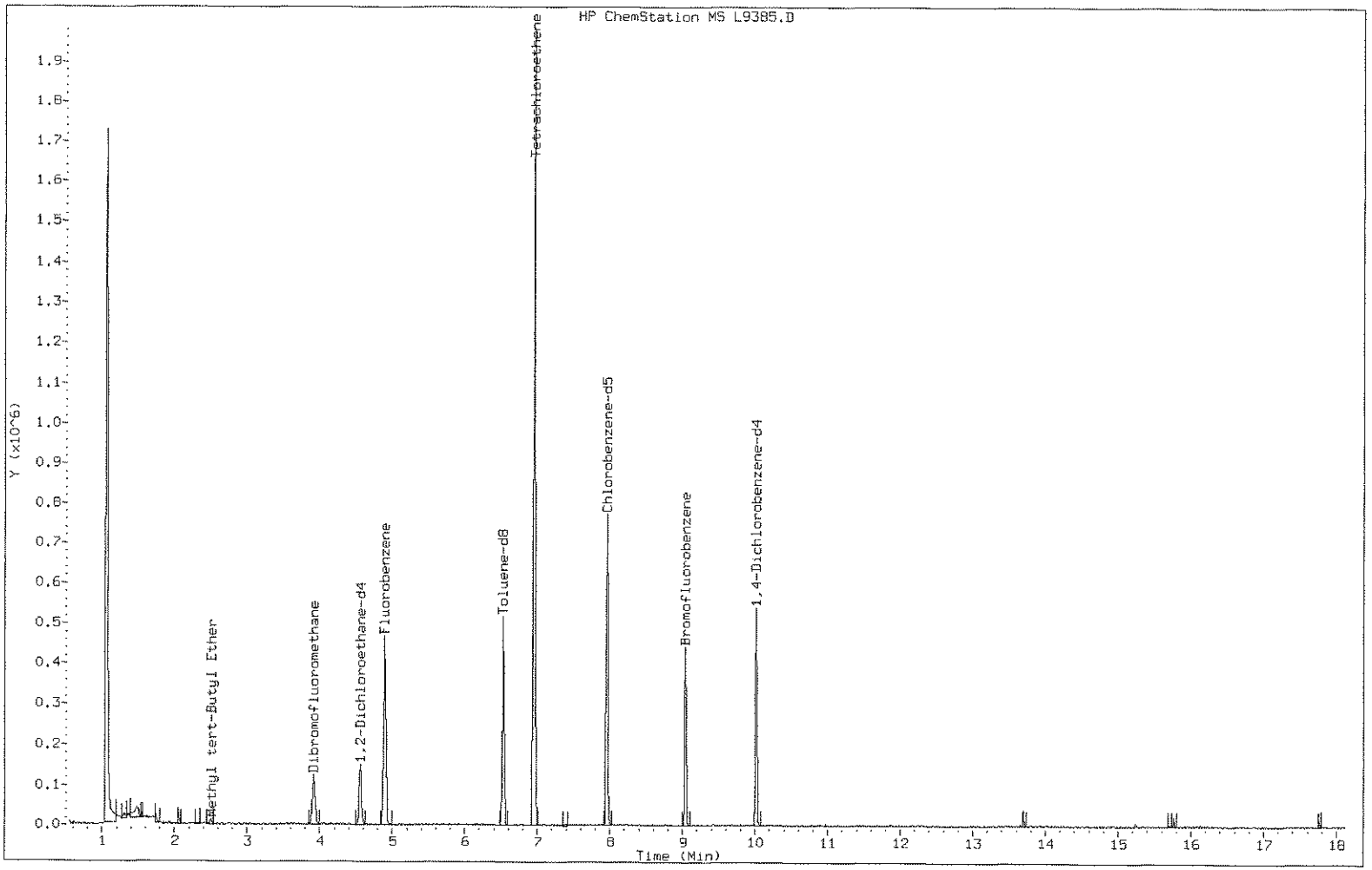
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.895	4.901	(1.000)	447266	25.0000	
	24 Methyl tert-Butyl Ether	73	2.504	2.491	(0.512)	14697	0.79147	3
$	41 Dibromofluoromethane	111	3.921	3.927	(0.801)	109523	18.9809	19
$	55 1,2-Dichloroethane-d4	65	4.561	4.567	(0.932)	134051	19.0551	19
*	75 Chlorobenzene-d5	117	7.955	7.961	(1.000)	427654	25.0000	
$	77 Toluene-d8	98	6.528	6.535	(0.821)	315488	20.7471	21
	80 Tetrachloroethene	164	6.952	6.958	(0.874)	417757	116.265	460
*	95 1,4-Dichlorobenzene-d4	152	10.011	10.018	(1.000)	119467	25.0000	
$	125 Bromofluorobenzene	95	9.037	9.043	(0.903)	128701	28.7158	29

Data File: L9385.D

Date: 02-AUG-2007 16:23

Client ID: A-14 Instrument: msl.i

Sample Info: 220-2277-B-17 Operator: D. HUMBERT



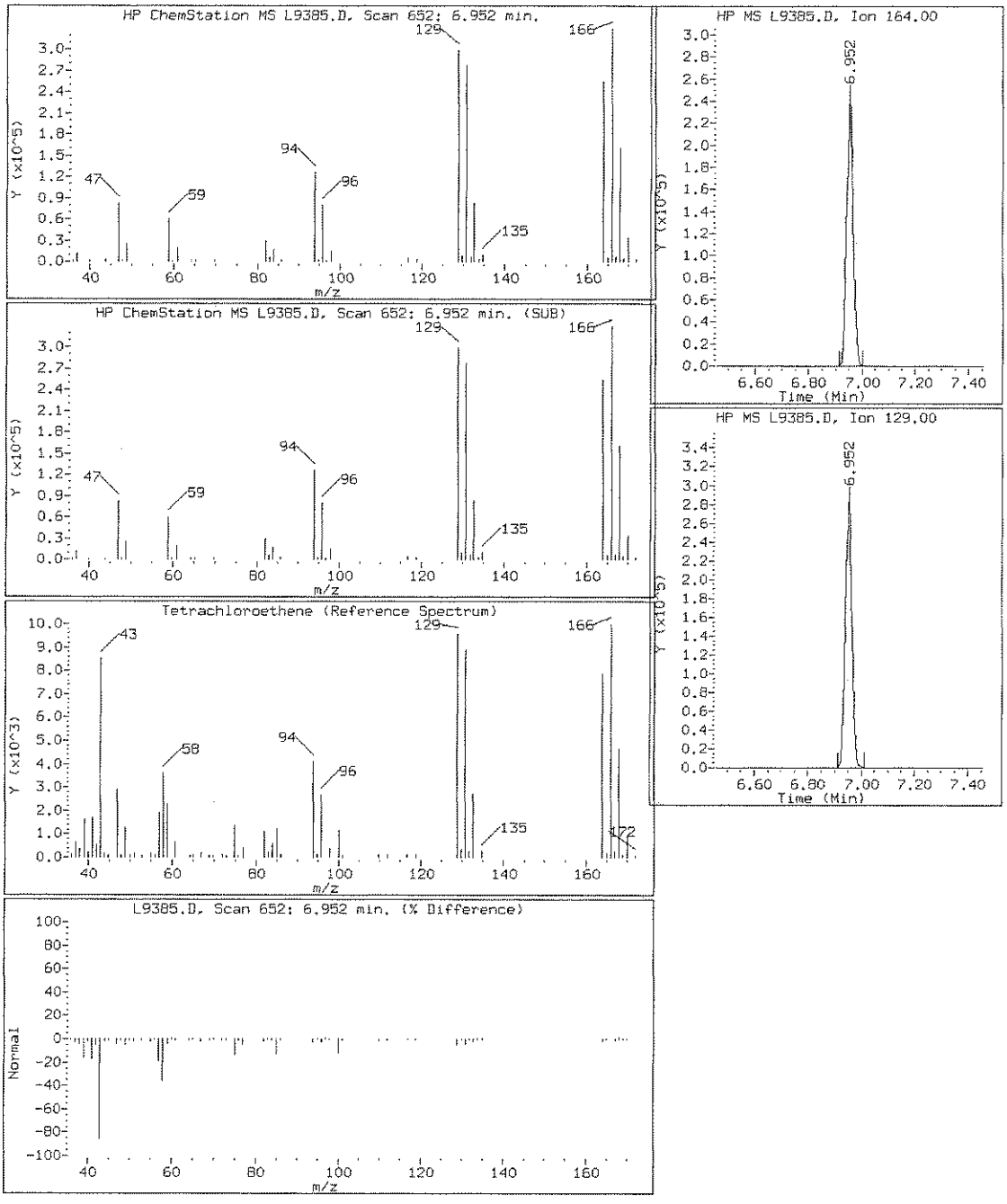
Data File: L9385.D

Date: 02-AUG-2007 16:23

Client ID: A-14 Instrument: msl.i

Sample Info: 220-2277-B-17 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-15 | Lab Sample ID: 220-2277-18

Matrix: Water | Lab File ID: L9380.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/02/2007 14:12

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8356 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	10		5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\Files\chem\VOA\msl.i\L079370.b\L9380.D
Lab Smp Id:
Inj Date : 02-AUG-2007 14:12 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-B-18
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\target1_ct\Files\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 62
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONVOAMGR1

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

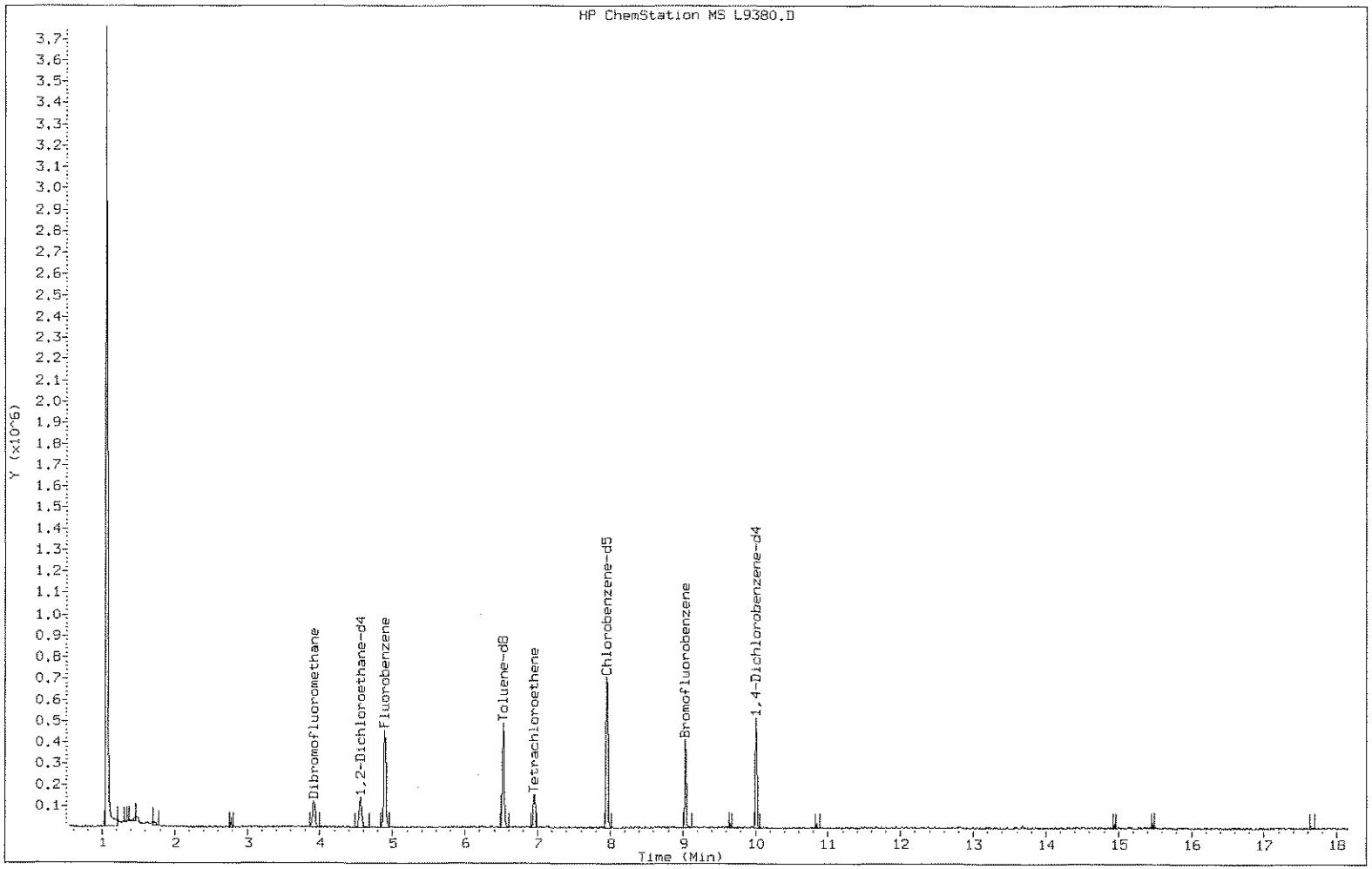
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1	Fluorobenzene	96	4.891	4.901	(1.000)	441024	25.0000	
$ 41	Dibromofluoromethane	111	3.917	3.927	(0.801)	106509	18.7198	19
$ 55	1,2-Dichloroethane-d4	65	4.556	4.567	(0.932)	130244	18.7760	19
* 75	Chlorobenzene-d5	117	7.951	7.961	(1.000)	425077	25.0000	
$ 77	Toluene-d8	98	6.524	6.535	(0.821)	303160	20.0573	20
80	Tetrachloroethene	164	6.947	6.958	(0.874)	36227	10.1434	10
* 95	1,4-Dichlorobenzene-d4	152	10.007	10.018	(1.000)	119186	25.0000	
$ 125	Bromofluorobenzene	95	9.033	9.043	(0.903)	125880	28.1526	28

Data File: L9380.D

Date: 02-AUG-2007 14:12

Client ID: A-15 Instrument: msl.i

Sample Info: 220-2277-B-18 Operator: D. HUMBERT



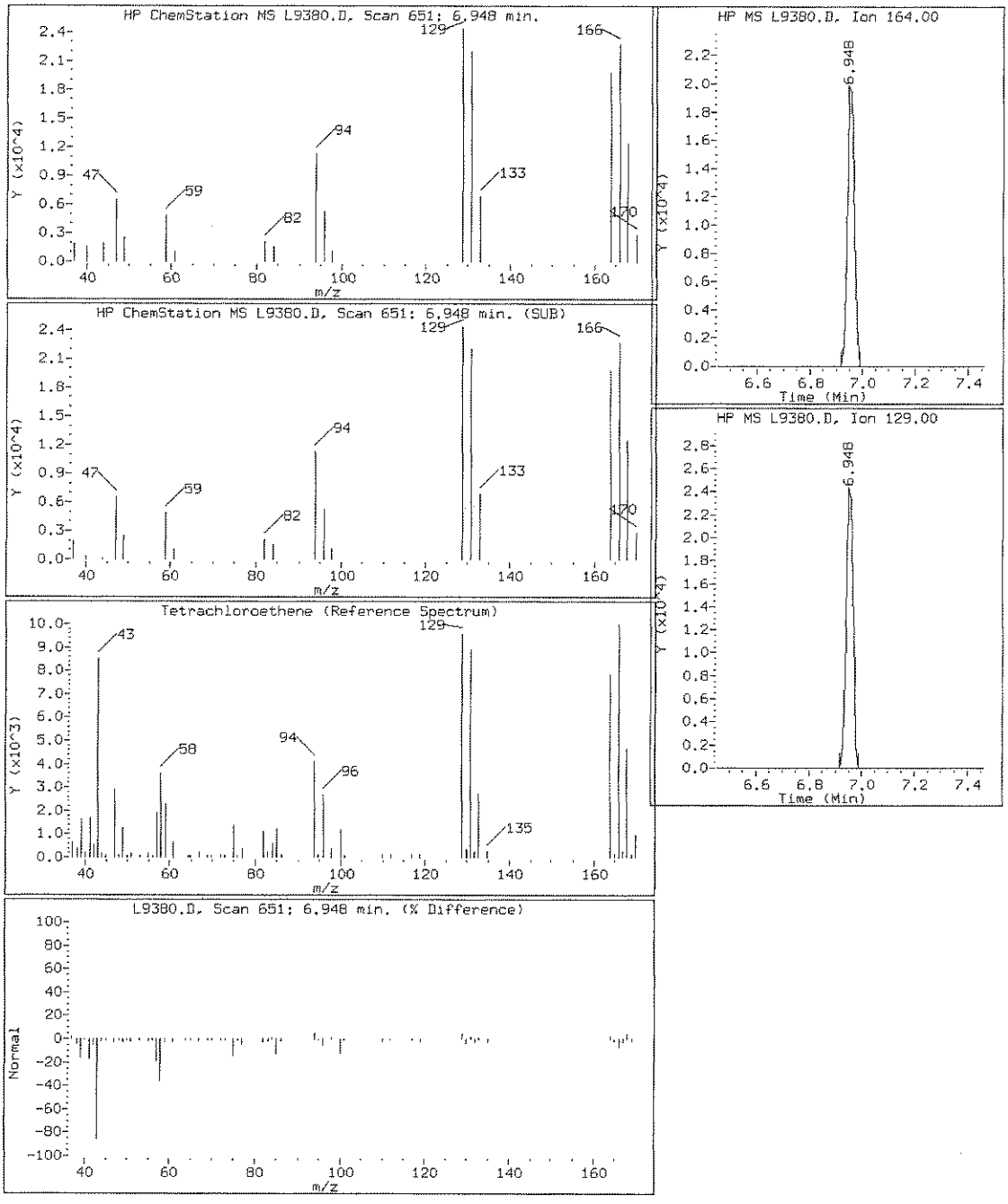
Data File: L9380.D

Date: 02-AUG-2007 14:12

Client ID: A-15 Instrument: msl.i

Sample Info: 220-2277-B-18 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-9 | Lab Sample ID: 220-2277-19

Matrix: Water | Lab File ID: L9357.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/01/2007 17:15

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8321 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	23		5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\Files\chem\VOA\msl.i\L079341.b\L9357.D
Lab Smp Id:
Inj Date : 01-AUG-2007 17:15 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-A-19
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\target1_ct\Files\chem\VOA\msl.i\L079341.b\L8260BNW.m
Meth Date : 01-Aug-2007 10:17 ctvoa Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 44
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONVOAMGR1

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

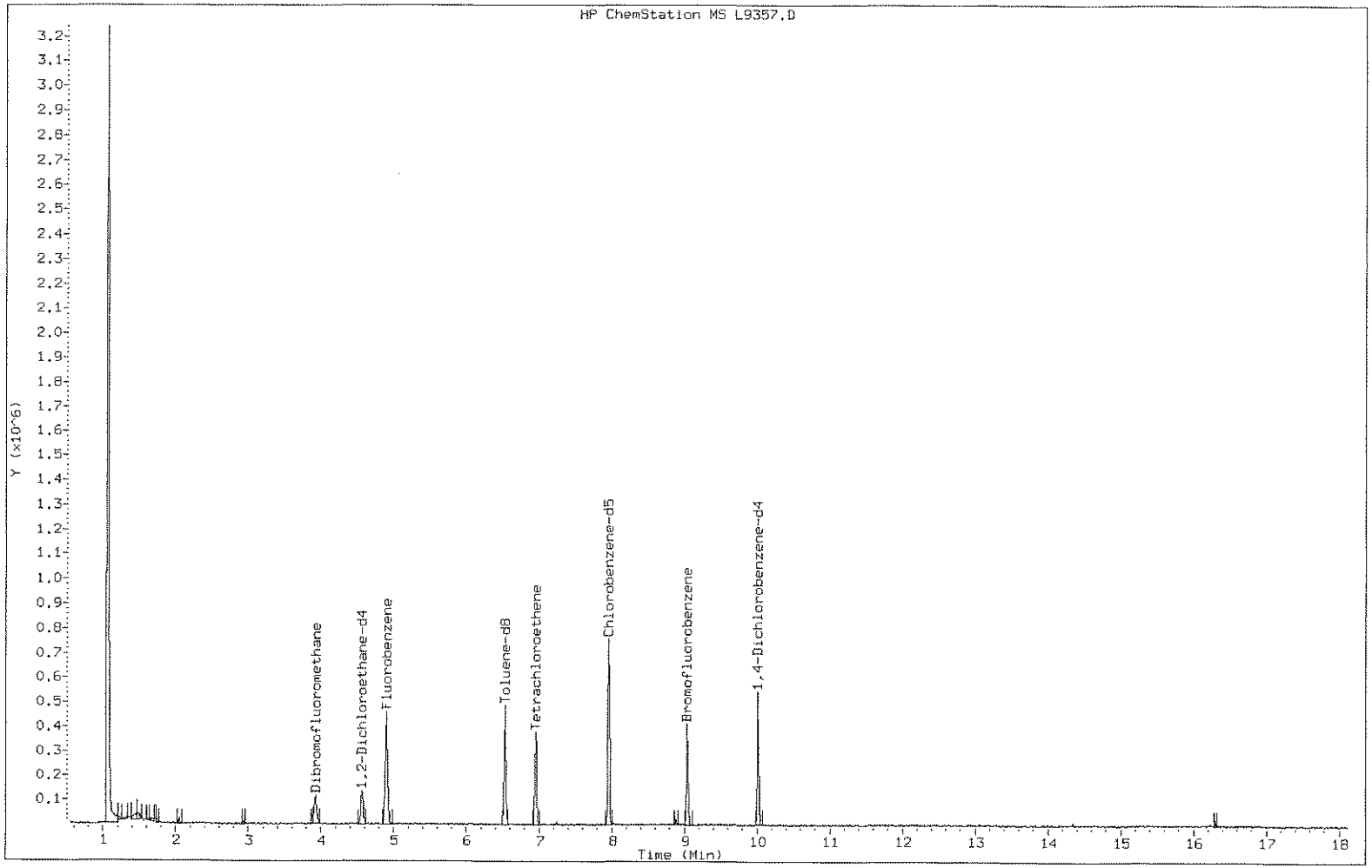
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.896	4.908	(1.000)	443072	25.0000	
	38 Chloroform	83	3.705	3.707	(0.757)	2898	0.31302	0.3
$	41 Dibromofluoromethane	111	3.921	3.924	(0.801)	104221	18.2330	18
$	55 1,2-Dichloroethane-d4	65	4.561	4.563	(0.932)	125527	18.0123	18
*	75 Chlorobenzene-d5	117	7.955	7.968	(1.000)	420067	25.0000	
$	77 Toluene-d8	98	6.529	6.541	(0.821)	300143	20.0945	20
	80 Tetrachloroethene	164	6.952	6.964	(0.874)	80461	22.7974	23
*	95 1,4-Dichlorobenzene-d4	152	10.012	10.014	(1.000)	117011	25.0000	
$	125 Bromofluorobenzene	95	9.038	9.040	(0.903)	120603	27.4738	27

Data File: L9357.D

Date: 01-AUG-2007 17:15

Client ID: A-9 Instrument: msl.i

Sample Info: 220-2277-A-19 Operator: D. HUMBERT



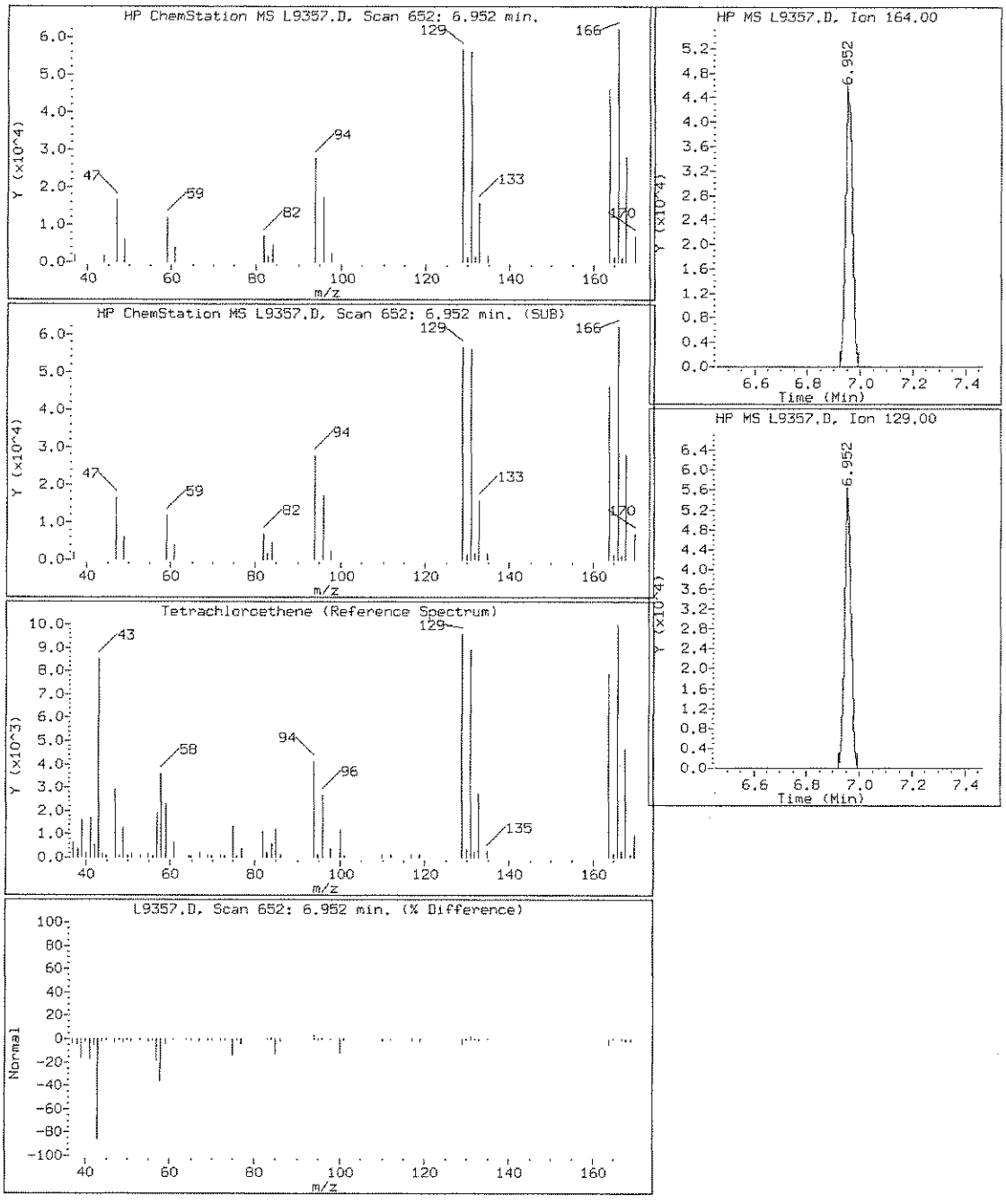
Data File: L9357.D

Date: 01-AUG-2007 17:15

Client ID: A-9 Instrument: msl.i

Sample Info: 220-2277-A-19 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-10 | Lab Sample ID: 220-2277-20

Matrix: Water | Lab File ID: L9358.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/01/2007 17:40

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8321 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	10		5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079341.b\L9358.D
Lab Smp Id: 220-2277-A-20 Client Smp ID: A-10
Inj Date : 01-AUG-2007 17:40 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-A-20
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079341.b\L8260BNW.m
Meth Date : 02-Aug-2007 09:12 target Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 45
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

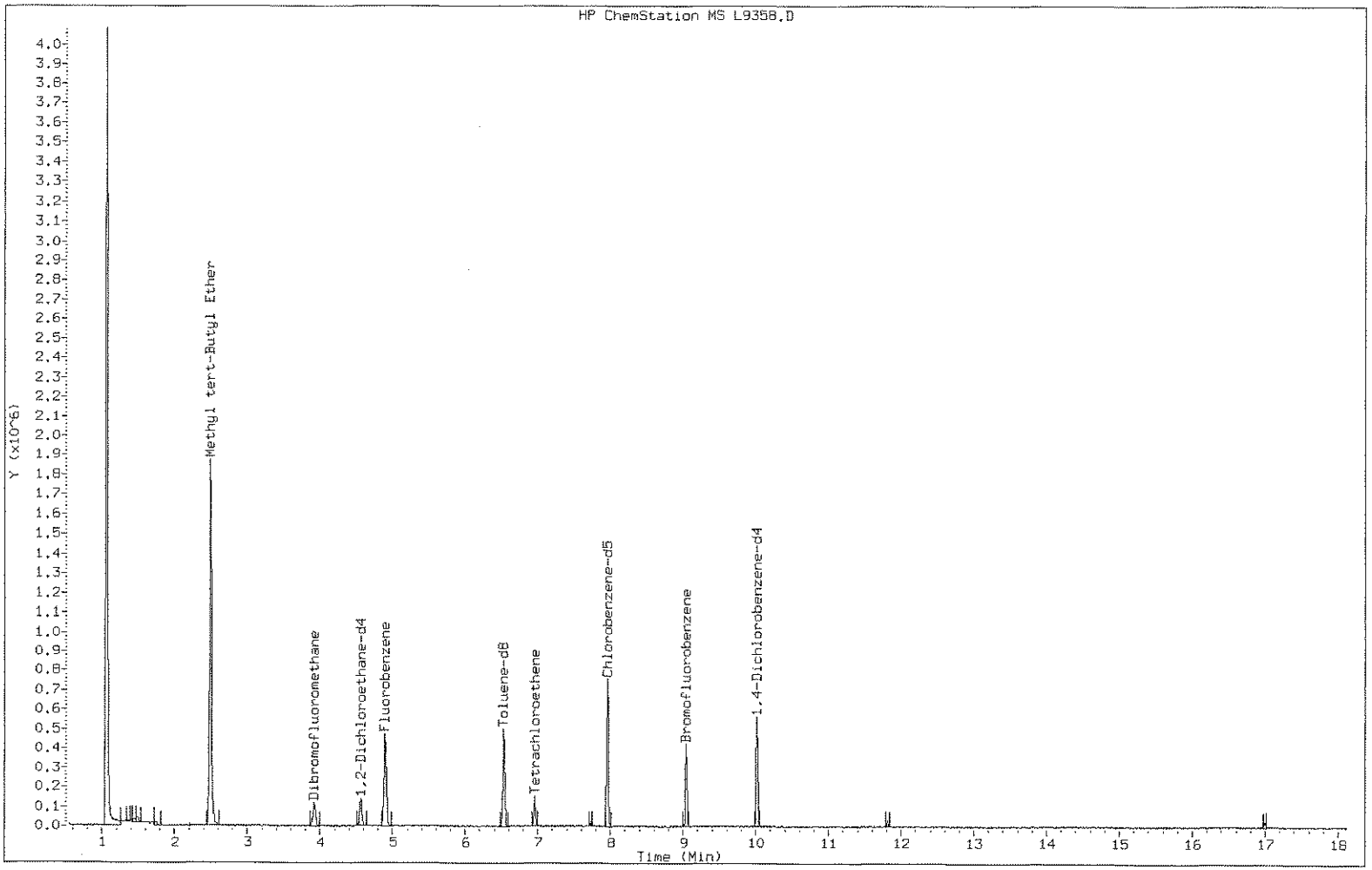
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1	Fluorobenzene	96	4.895	4.908	(1.000)	444007	25.0000	
24	Methyl tert-Butyl Ether	73	2.485	2.497	(0.508)	1960279	106.340	110
$ 41	Dibromofluoromethane	111	3.921	3.924	(0.801)	105732	18.4584	18
$ 55	1,2-Dichloroethane-d4	65	4.561	4.563	(0.932)	127827	18.3037	18
* 75	Chlorobenzene-d5	117	7.955	7.968	(1.000)	428130	25.0000	
$ 77	Toluene-d8	98	6.528	6.541	(0.821)	303985	19.9684	20
80	Tetrachloroethene	164	6.951	6.964	(0.874)	35853	9.96710	10
* 95	1,4-Dichlorobenzene-d4	152	10.011	10.014	(1.000)	123039	25.0000	
$ 125	Bromofluorobenzene	95	9.037	9.040	(0.903)	125674	27.2264	27

Data File: L9358.D

Date: 01-AUG-2007 17:40

Client ID: A-10 Instrument: msl.i

Sample Info: 220-2277-A-20 Operator: D. HUMBERT



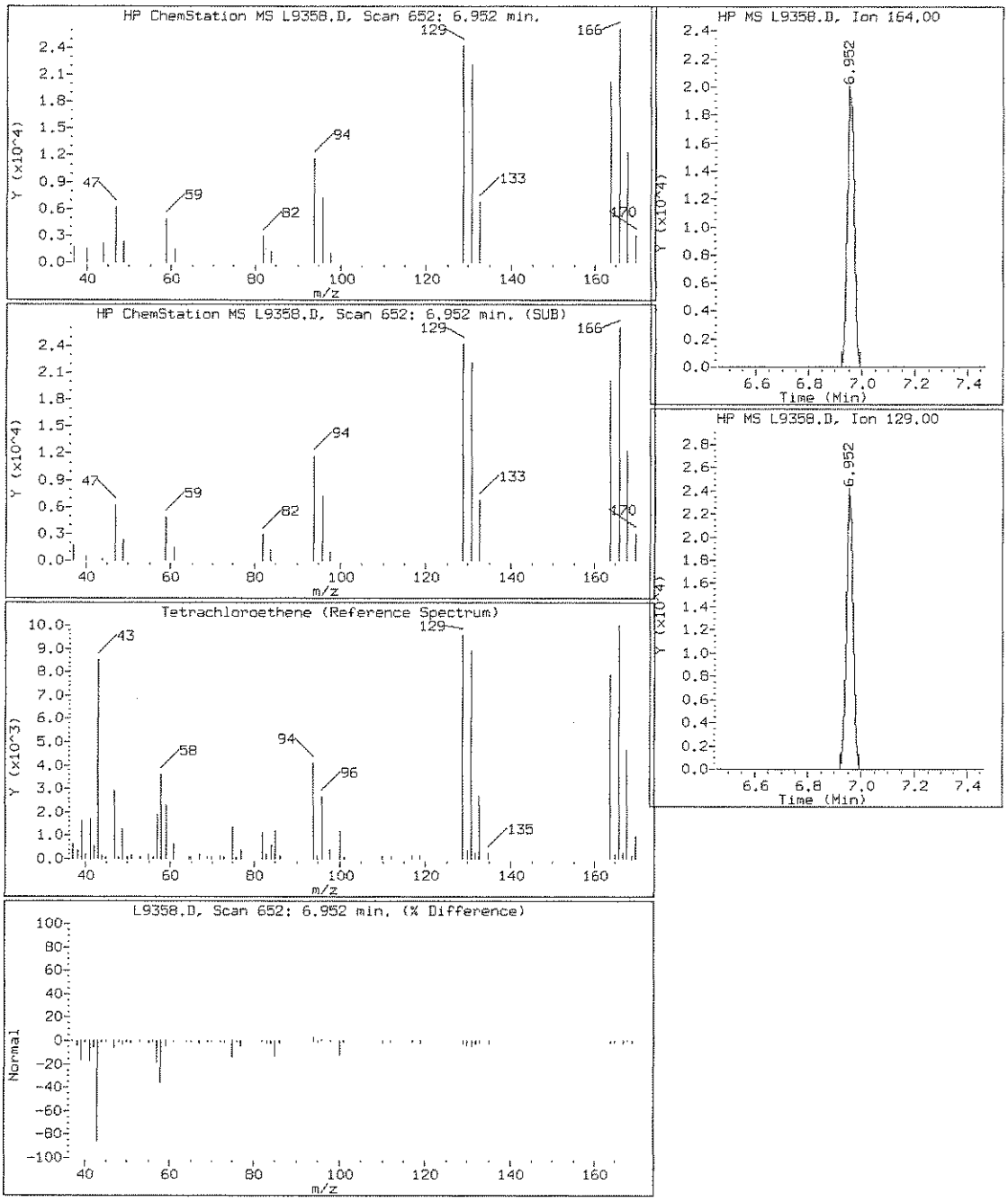
Data File: L9358.D

Date: 01-AUG-2007 17:40

Client ID: A-10 Instrument: msl.i

Sample Info: 220-2277-A-20 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: MW-113 | Lab Sample ID: 220-2277-21

Matrix: Water | Lab File ID: L9359.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/01/2007 18:04

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8321 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	3.4	J B	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	8.7		5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079341.b\L9359.D
Lab Smp Id: 220-2277-A-21 Client Smp ID: MW-113
Inj Date : 01-AUG-2007 18:04 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-A-21
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079341.b\L8260BNW.m
Meth Date : 02-Aug-2007 09:12 target Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 46
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

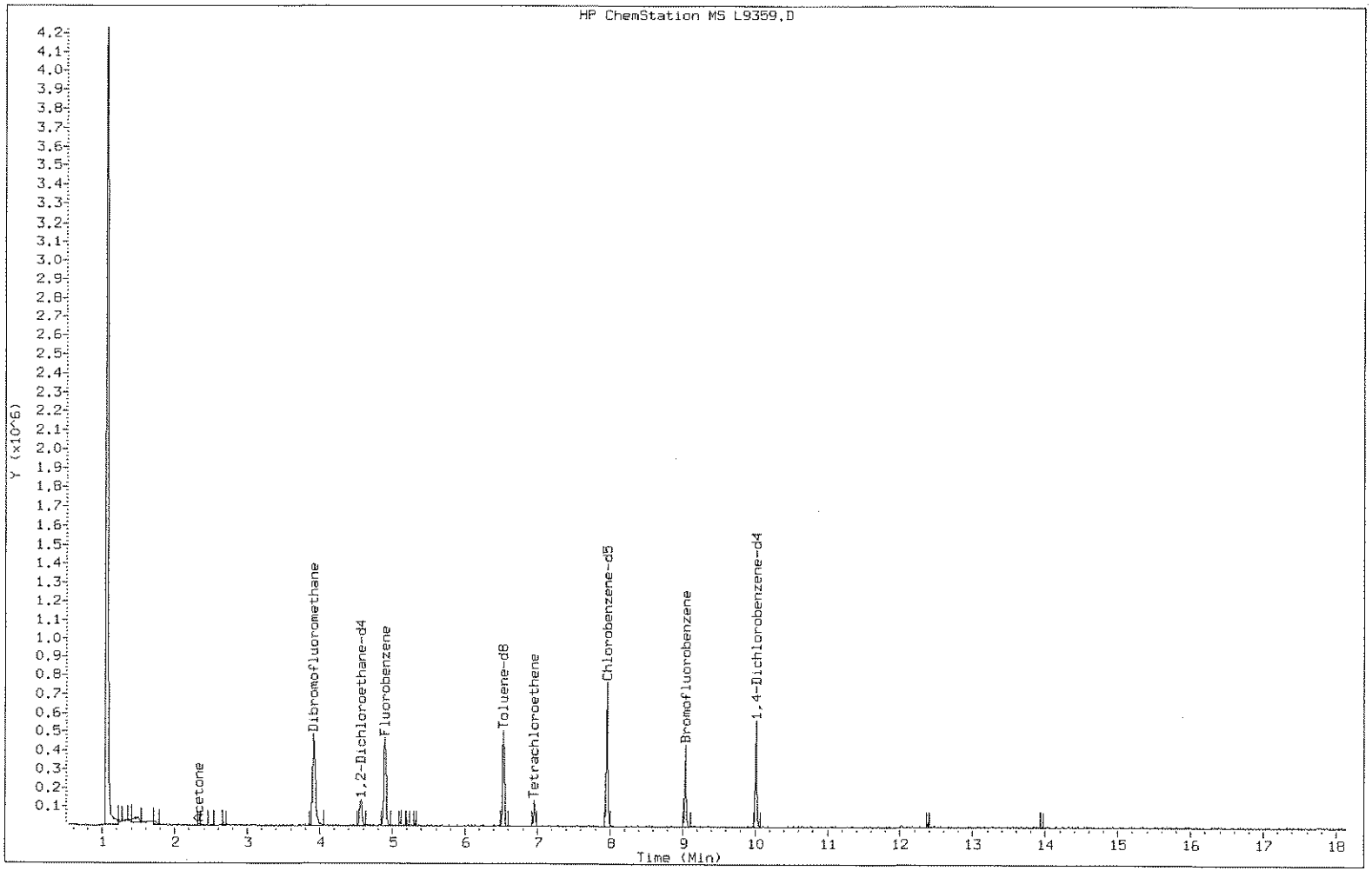
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1	Fluorobenzene	96	4.897	4.908	(1.000)	435324	25.0000	
21	Acetone	43	2.329	2.330	(0.476)	9303	3.36972	3
$ 41	Dibromofluoromethane	111	3.913	3.924	(0.799)	105841	18.8460	19
42	Tetrahydrofuran	42	3.913	3.924	(0.799)	344926	146.236	150
$ 55	1,2-Dichloroethane-d4	65	4.563	4.563	(0.932)	125079	18.2675	18
* 75	Chlorobenzene-d5	117	7.957	7.968	(1.000)	422393	25.0000	
$ 77	Toluene-d8	98	6.531	6.541	(0.821)	304927	20.3024	20
80	Tetrachloroethene	164	6.954	6.964	(0.874)	30920	8.71248	9
* 95	1,4-Dichlorobenzene-d4	152	10.014	10.014	(1.000)	122483	25.0000	
$ 125	Bromofluorobenzene	95	9.040	9.040	(0.903)	126892	27.6151	28

Data File: L9359.D

Date: 01-AUG-2007 18:04

Client ID: MW-113 Instrument: msl.i

Sample Info: 220-2277-A-21 Operator: D. HUMBERT



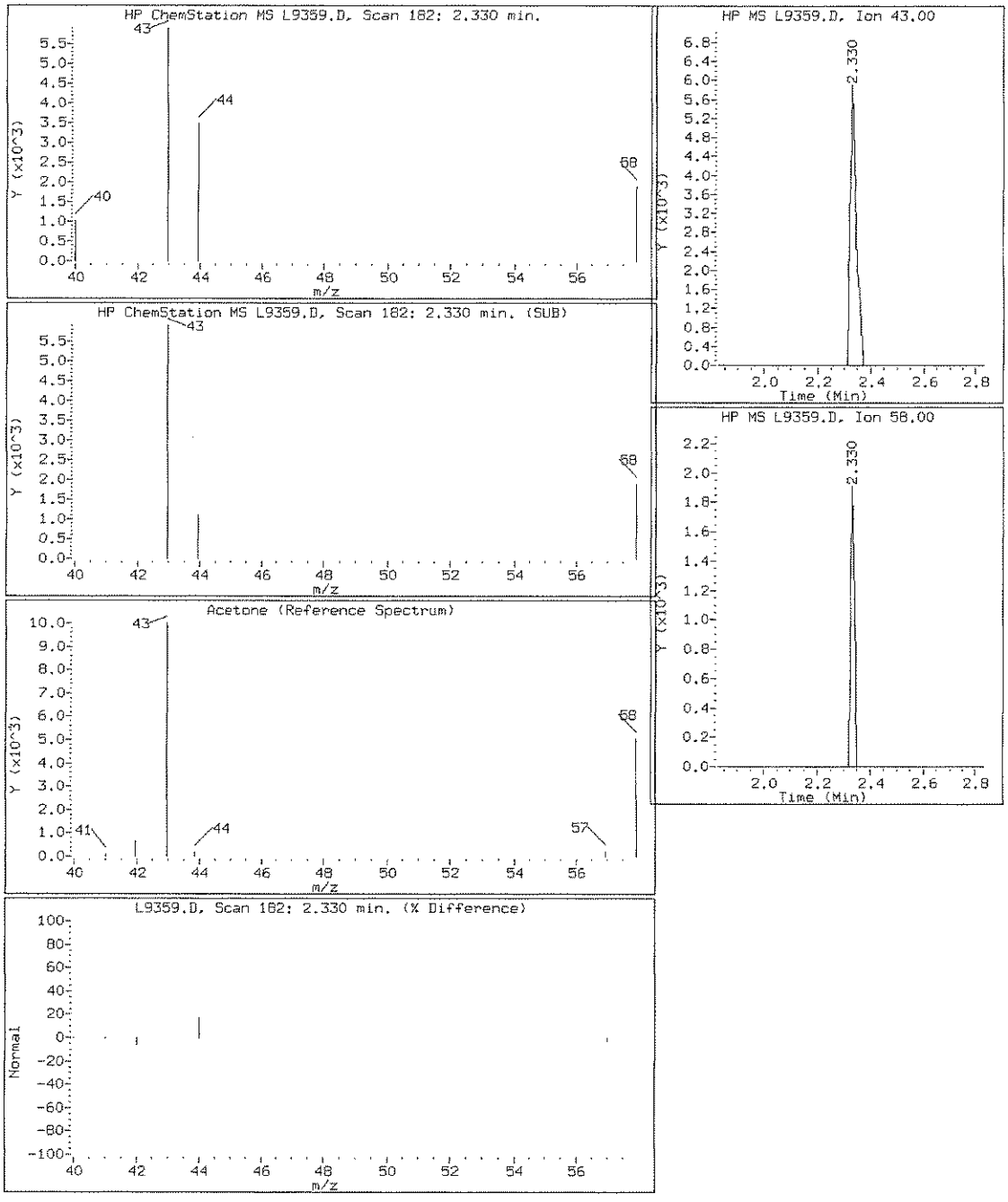
Data File: L9359.D

Date: 01-AUG-2007 18:04

Client ID: MW-113 Instrument: msl.i

Sample Info: 220-2277-A-21 Operator: D. HUMBERT

21 Acetone



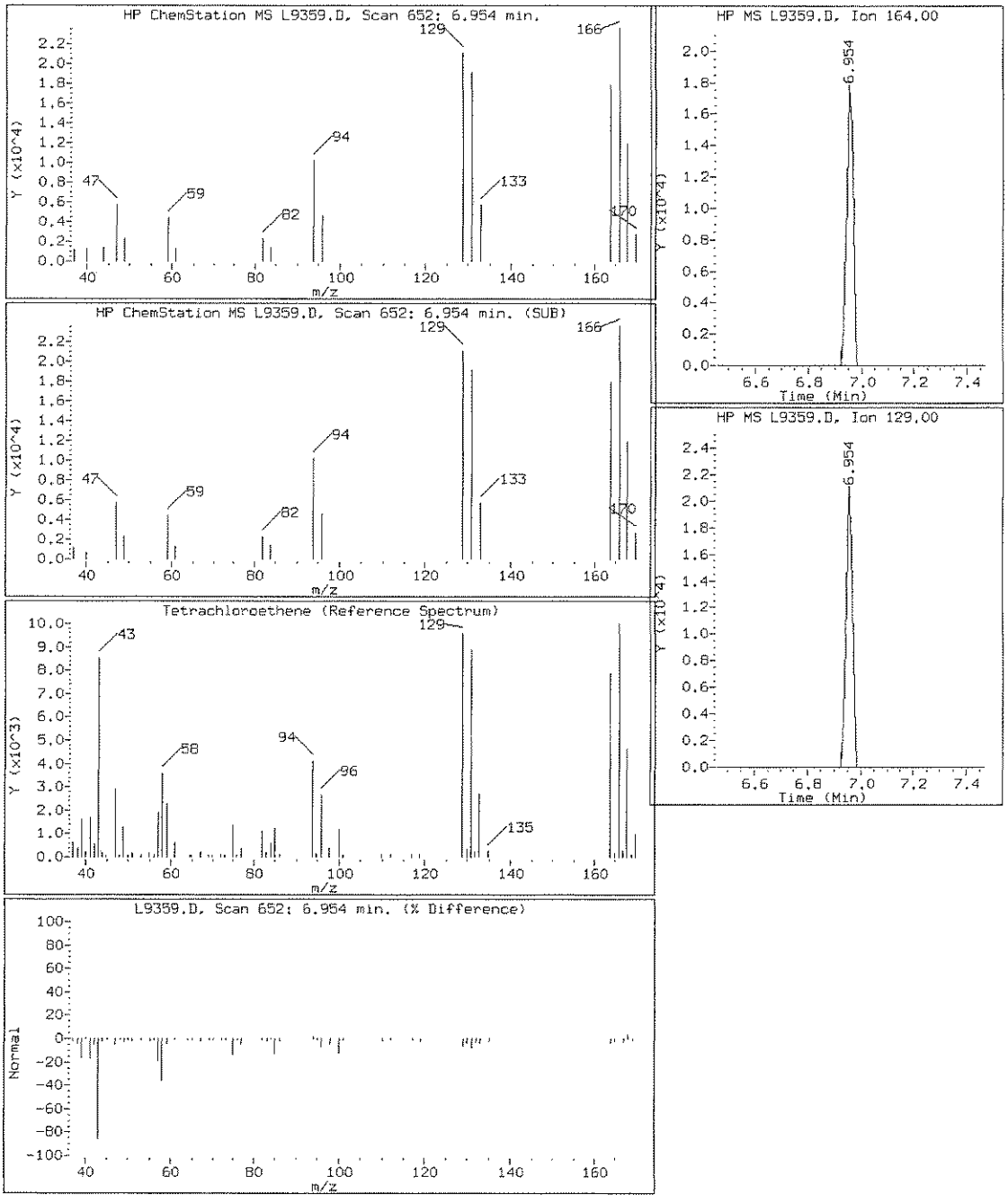
Data File: L9359.D

Date: 01-AUG-2007 18:04

Client ID: MW-113 Instrument: msl.i

Sample Info: 220-2277-A-21 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: MW-112 | Lab Sample ID: 220-2277-22

Matrix: Water | Lab File ID: L9381.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/02/2007 14:36

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8356 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	1.3	J	5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079370.b\L9381.D
Lab Smp Id: 220-2277-B-22 Client Smp ID: MW-112
Inj Date : 02-AUG-2007 14:36 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-B-22
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 63
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

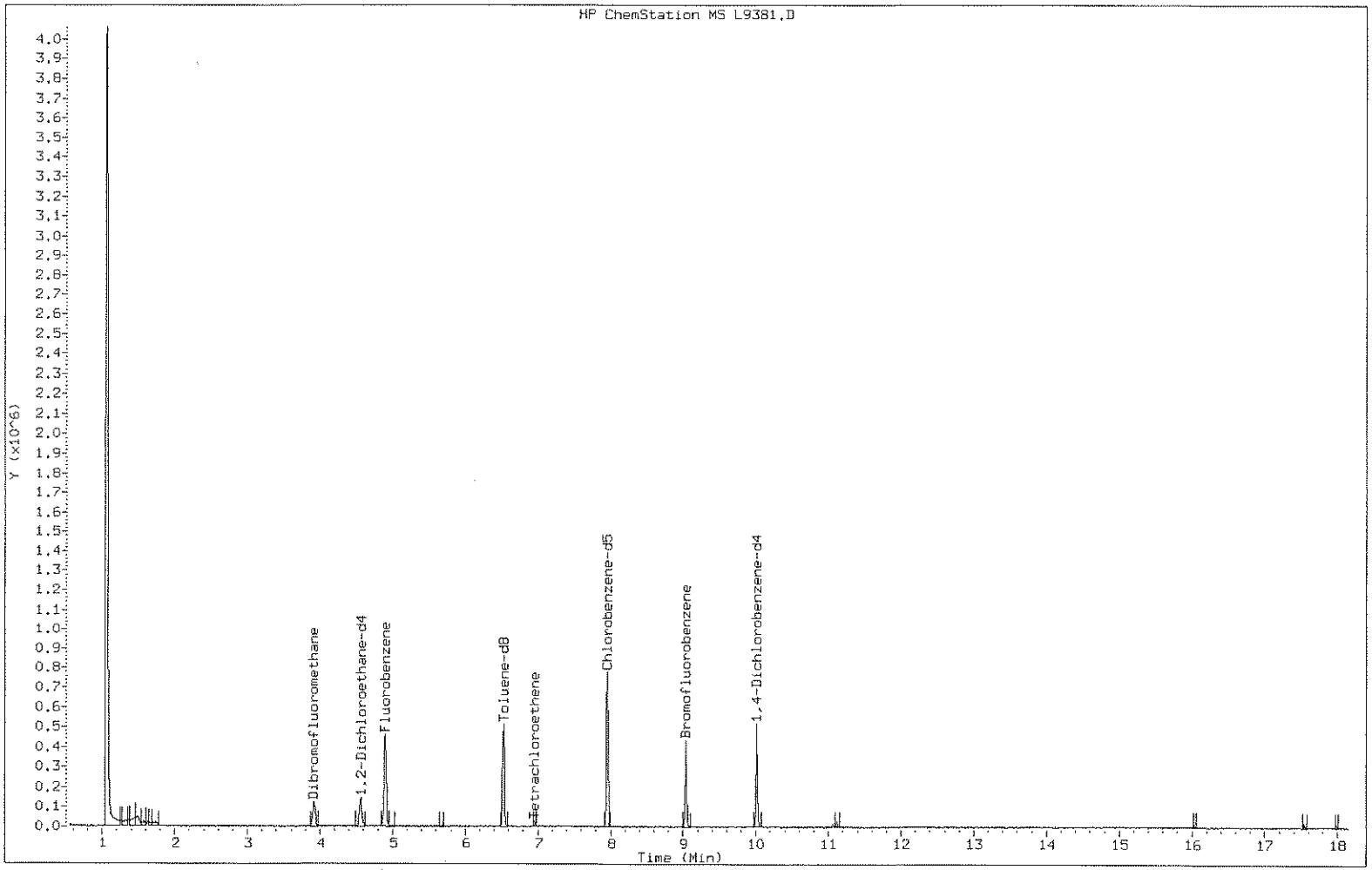
Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1 Fluorobenzene	96	4.896	4.901	(1.000)	452327	25.0000	
$ 41 Dibromofluoromethane	111	3.912	3.927	(0.799)	112430	19.2667	19
$ 55 1,2-Dichloroethane-d4	65	4.562	4.567	(0.932)	133195	18.7216	19
* 75 Chlorobenzene-d5	117	7.956	7.961	(1.000)	435304	25.0000	
$ 77 Toluene-d8	98	6.530	6.535	(0.821)	323853	20.9229	21
80 Tetrachloroethene	164	6.953	6.958	(0.874)	4862	1.32936	1
* 95 1,4-Dichlorobenzene-d4	152	10.013	10.018	(1.000)	121255	25.0000	
$ 125 Bromofluorobenzene	95	9.039	9.043	(0.903)	130839	28.7624	29

Data File: L9381.D

Date: 02-AUG-2007 14:36

Client ID: MW-112 Instrument: msl.i

Sample Info: 220-2277-B-22 Operator: D. HUMBERT



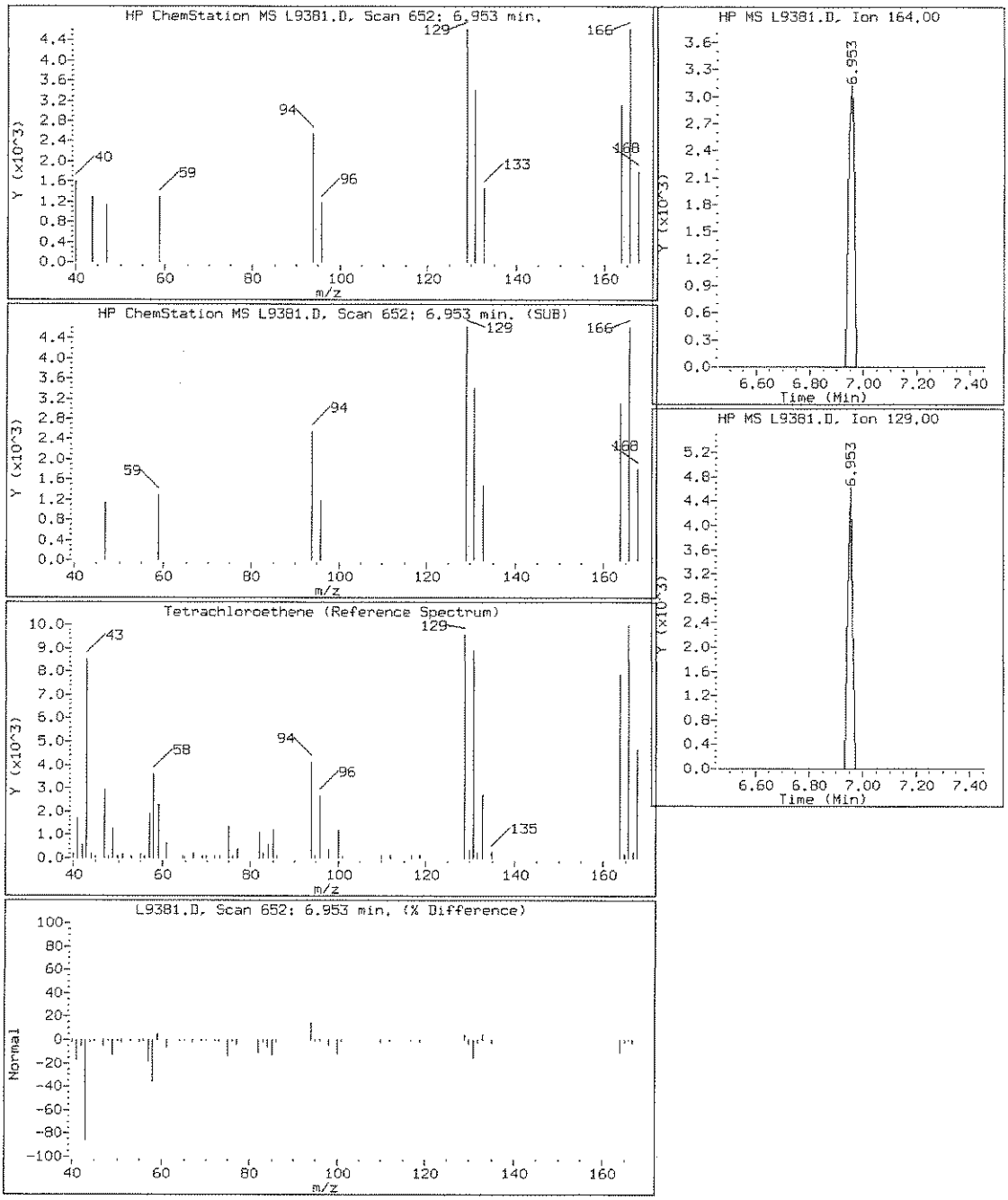
Data File: L9381.D

Date: 02-AUG-2007 14:36

Client ID: MW-112 Instrument: msl.i

Sample Info: 220-2277-B-22 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: ME-8 | Lab Sample ID: 220-2277-23

Matrix: Water | Lab File ID: L9388.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/02/2007 17:37

Level: (low/med) Low | Dilution Factor: 2

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8356 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	20	U	20	2.8
71-43-2	Benzene	10	U	10	0.80
75-27-4	Bromodichloromethane	10	U	10	0.80
75-25-2	Bromoform	10	U	10	1.6
74-83-9	Bromomethane	10	U	10	2.4
78-93-3	2-Butanone (MEK)	20	U	20	2.4
75-15-0	Carbon disulfide	10	U	10	1.8
56-23-5	Carbon tetrachloride	10	U	10	2.0
108-90-7	Chlorobenzene	10	U	10	0.80
75-00-3	Chloroethane	10	U	10	1.6
67-66-3	Chloroform	10	U	10	1.4
74-87-3	Chloromethane	10	U	10	1.0
124-48-1	Dibromochloromethane	10	U	10	1.0
75-34-3	1,1-Dichloroethane	10	U	10	1.2
107-06-2	1,2-Dichloroethane	10	U	10	1.2
75-35-4	1,1-Dichloroethene	10	U	10	1.4
78-87-5	1,2-Dichloropropane	10	U	10	1.8
10061-01-5	cis-1,3-Dichloropropene	10	U	10	1.0
10061-02-6	trans-1,3-Dichloropropene	10	U	10	1.6
100-41-4	Ethylbenzene	10	U	10	2.0
591-78-6	2-Hexanone	20	U	20	1.6
75-09-2	Methylene Chloride	10	U	10	0.80
108-10-1	4-Methyl-2-pentanone (MIBK)	20	U	20	1.4
100-42-5	Styrene	10	U	10	1.0
79-34-5	1,1,2,2-Tetrachloroethane	10	U	10	0.80
127-18-4	Tetrachloroethene	250		10	1.0
108-88-3	Toluene	10	U	10	0.60
71-55-6	1,1,1-Trichloroethane	10	U	10	0.80
79-00-5	1,1,2-Trichloroethane	10	U	10	1.2
79-01-6	Trichloroethene	10	U	10	1.4
75-01-4	Vinyl chloride	10	U	10	1.6
1330-20-7	Xylenes, Total	10	U	10	2.0
156-59-2	cis-1,2-Dichloroethene	10	U	10	1.2
156-60-5	trans-1,2-Dichloroethene	10	U	10	1.0

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\Files\chem\VOA\msl.i\L079370.b\L9388.D
Lab Smp Id:
Inj Date : 02-AUG-2007 17:37 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-B-23
Misc Info : : ;;; ; 8260 ; 2 ; LLW
Comment :
Method : \\target1_ct\Files\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 70
Dil Factor: 2.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONVOAMGR1

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	2.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

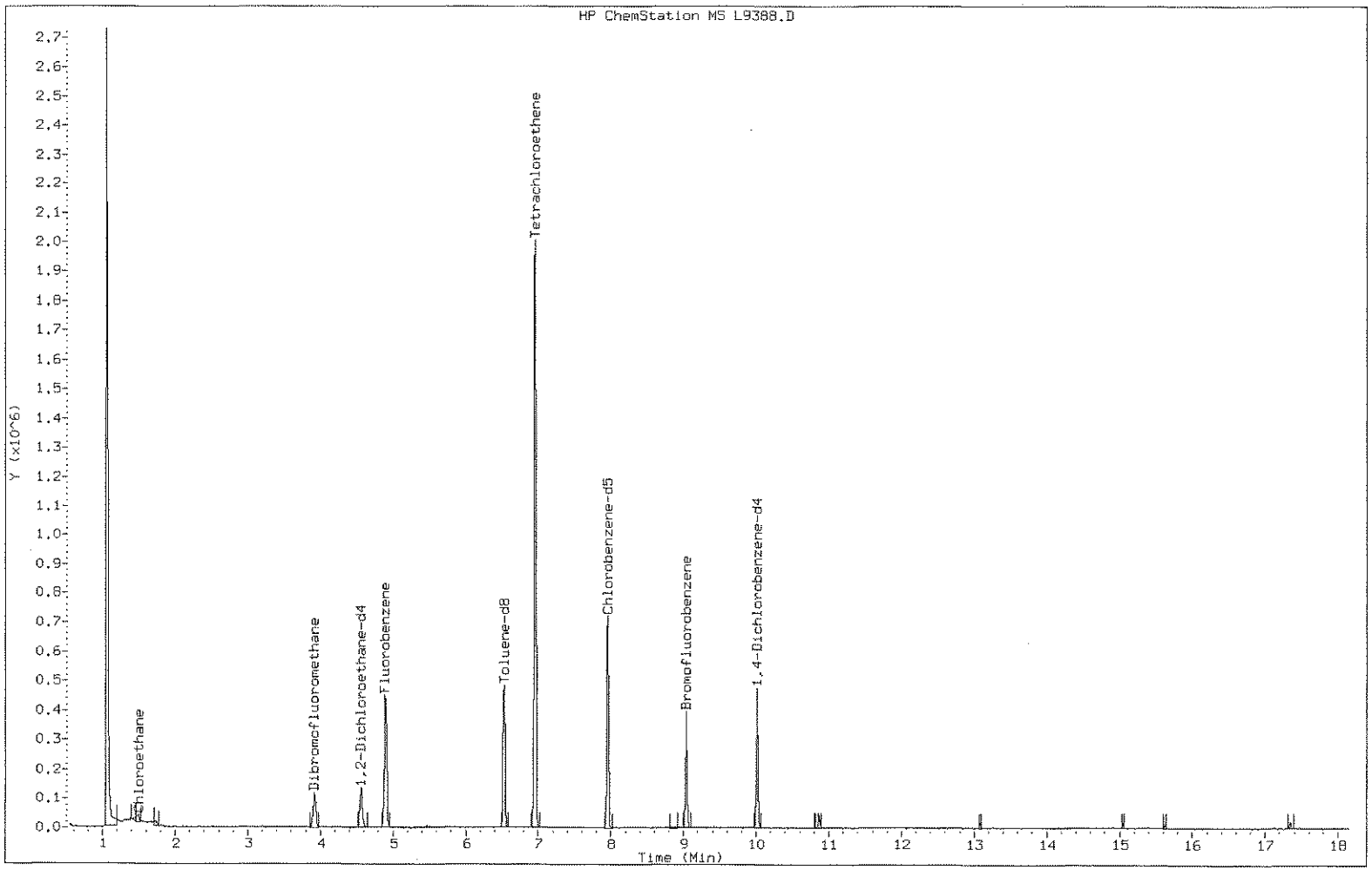
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.889	4.901	(1.000)	440199	25.0000	
	6 Chloroethane	64	1.485	1.546	(0.304)	1270	0.36766	0.7
$	41 Dibromofluoromethane	111	3.915	3.927	(0.801)	104503	18.4017	18
$	55 1,2-Dichloroethane-d4	65	4.554	4.567	(0.932)	124697	18.0100	18
*	75 Chlorobenzene-d5	117	7.959	7.961	(1.000)	436237	25.0000	
$	77 Toluene-d8	98	6.532	6.535	(0.821)	310389	20.0102	20
	80 Tetrachloroethene	164	6.955	6.958	(0.874)	452715	123.516	250
*	95 1,4-Dichlorobenzene-d4	152	10.015	10.018	(1.000)	123622	25.0000	
$	125 Bromofluorobenzene	95	9.041	9.043	(0.903)	125775	27.1198	27

Data File: L9388.D

Date: 02-AUG-2007 17:37

Client ID: ME-8 Instrument: msl.i

Sample Info: 220-2277-B-23 Operator: D. HUMBERT



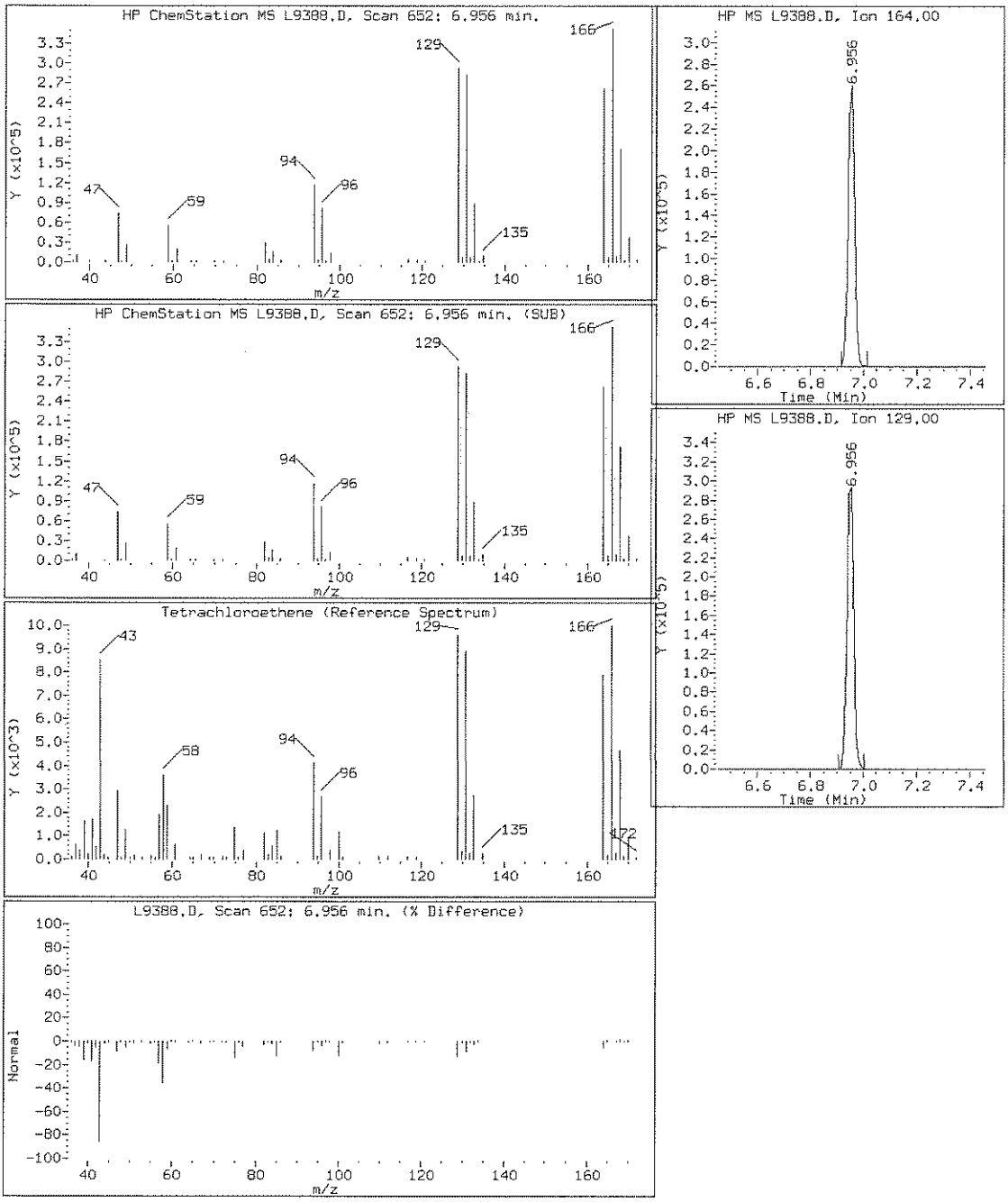
Data File: L9388.D

Date: 02-AUG-2007 17:37

Client ID: ME-8 Instrument: msl.i

Sample Info: 220-2277-B-23 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: ME-13 Lab Sample ID: 220-2277-24

Matrix: Water Lab File ID: L9362.D

Analysis Method: 8260B Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) Date Analyzed: 08/01/2007 19:19

Level: (low/med) Low Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) Soil Aliquot:

Soil Extract Vol.: % Moisture:

Analy. Batch No.: 8321 Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	54		5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079341.b\L9362.D
Lab Smp Id: 220-2277-A-24 Client Smp ID: ME-13
Inj Date : 01-AUG-2007 19:19 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-A-24
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079341.b\L8260BNW.m
Meth Date : 02-Aug-2007 09:12 target Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 49
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

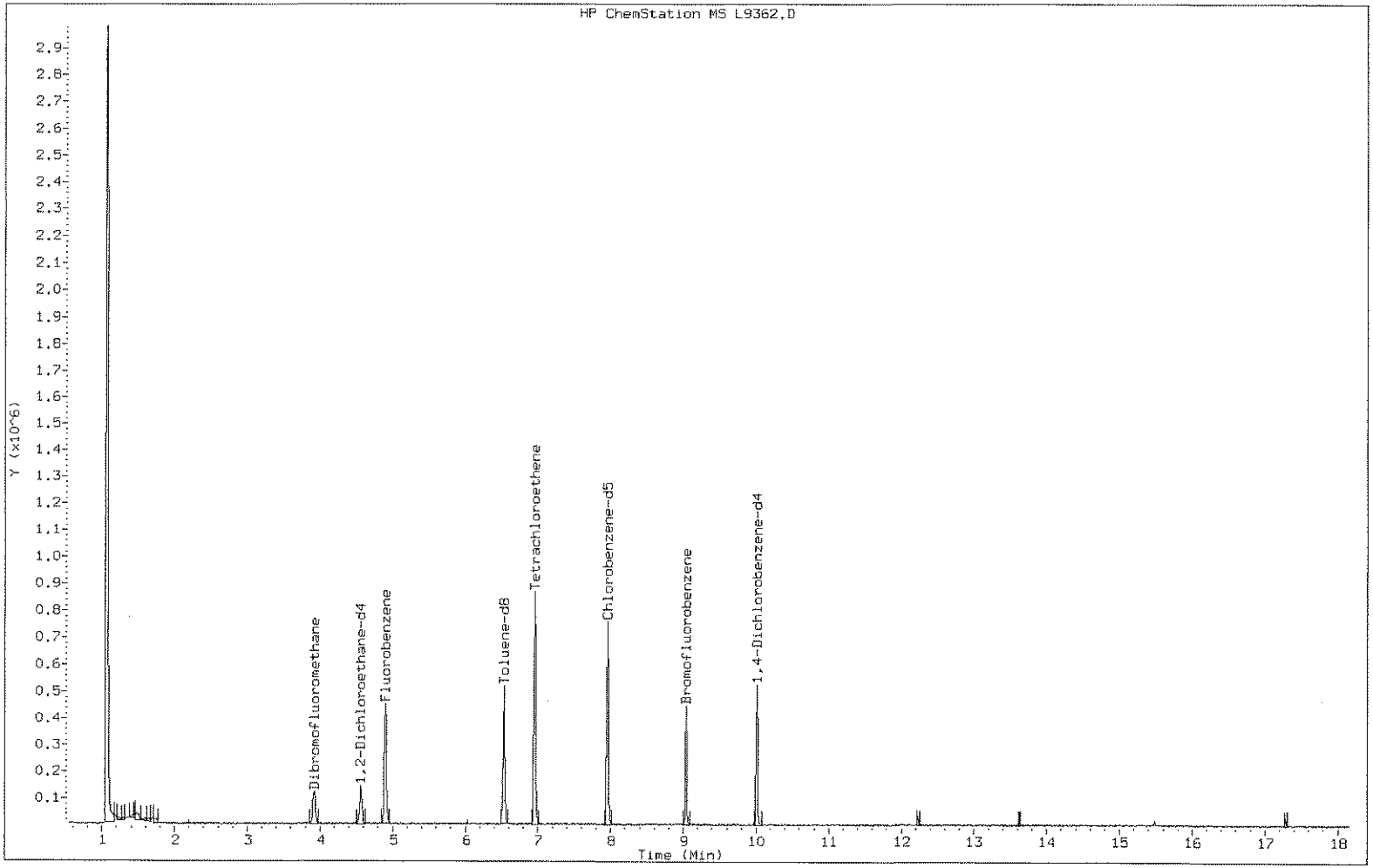
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.896	4.908	(1.000)	424369	25.0000	
$	41 Dibromofluoromethane	111	3.922	3.924	(0.801)	109325	19.9689	20
$	55 1,2-Dichloroethane-d4	65	4.562	4.563	(0.932)	133484	19.9982	20
*	75 Chlorobenzene-d5	117	7.956	7.968	(1.000)	413641	25.0000	
$	77 Toluene-d8	98	6.530	6.541	(0.821)	312744	21.2634	21
	80 Tetrachloroethene	164	6.953	6.964	(0.874)	186035	53.5291	54
*	95 1,4-Dichlorobenzene-d4	152	10.013	10.014	(1.000)	114741	25.0000	
$	125 Bromofluorobenzene	95	9.039	9.040	(0.903)	129012	29.9709	30

Data File: L9362.D

Date: 01-AUG-2007 19:19

Client ID: ME-13 Instrument: msl.i

Sample Info: 220-2277-A-24 Operator: D. HUMBERT



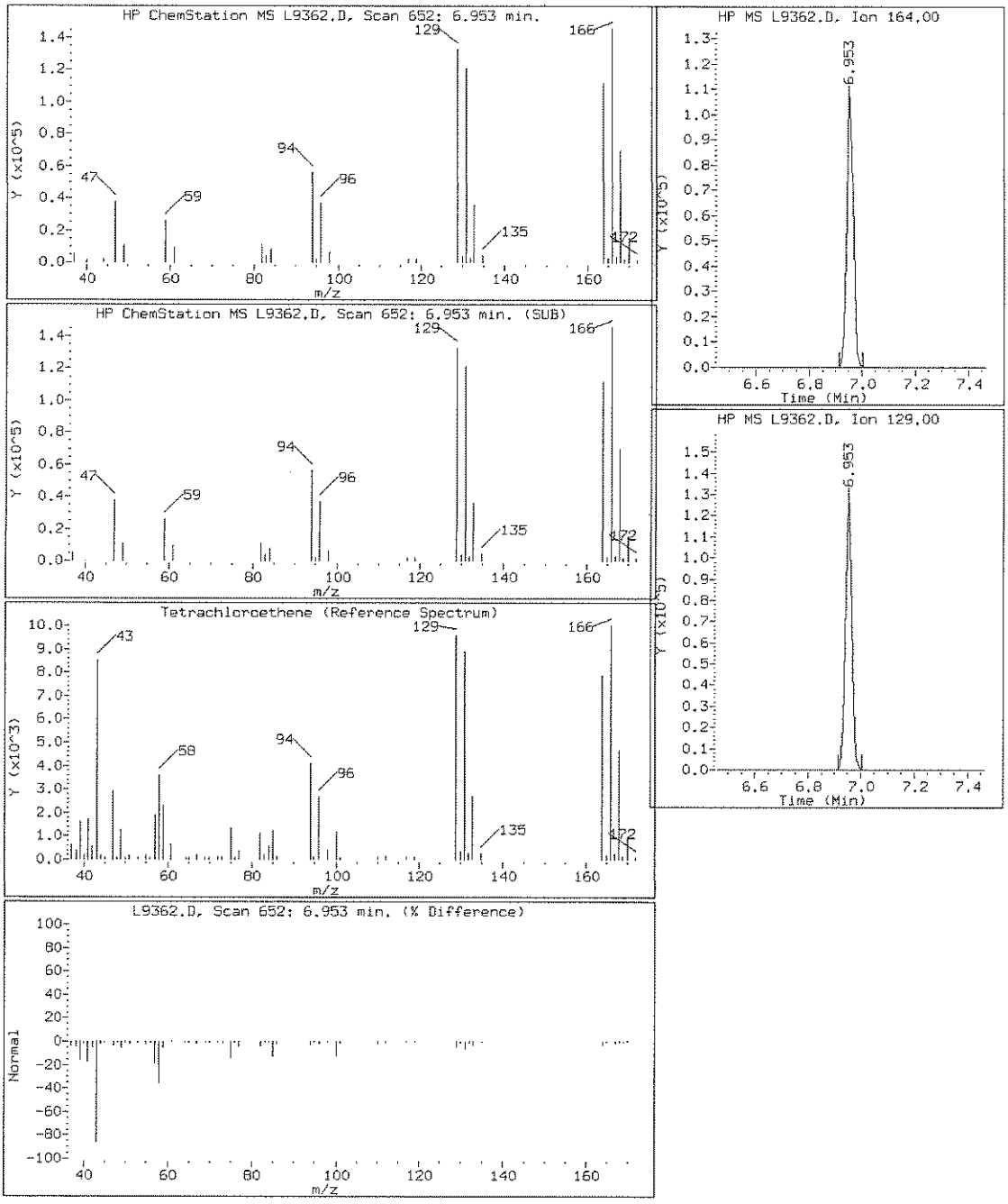
Data File: L9362.D

Date: 01-AUG-2007 19:19

Client ID: ME-13 Instrument: msl.i

Sample Info: 220-2277-A-24 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: FB0726 Lab Sample ID: 220-2277-25

Matrix: Water Lab File ID: L9363.D

Analysis Method: 8260B Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) Date Analyzed: 08/01/2007 19:44

Level: (low/med) Low Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) Soil Aliquot:

Soil Extract Vol.: % Moisture:

Analy. Batch No.: 8321 Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	3.0	J B	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	0.76	J B	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	5.0	U	5.0	0.50
108-88-3	Toluene	0.37	J	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079341.b\L9363.D
Lab Smp Id: 220-2277-A-25 Client Smp ID: FB0726
Inj Date : 01-AUG-2007 19:44 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-A-25
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079341.b\L8260BNW.m
Meth Date : 02-Aug-2007 09:12 target Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 50
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

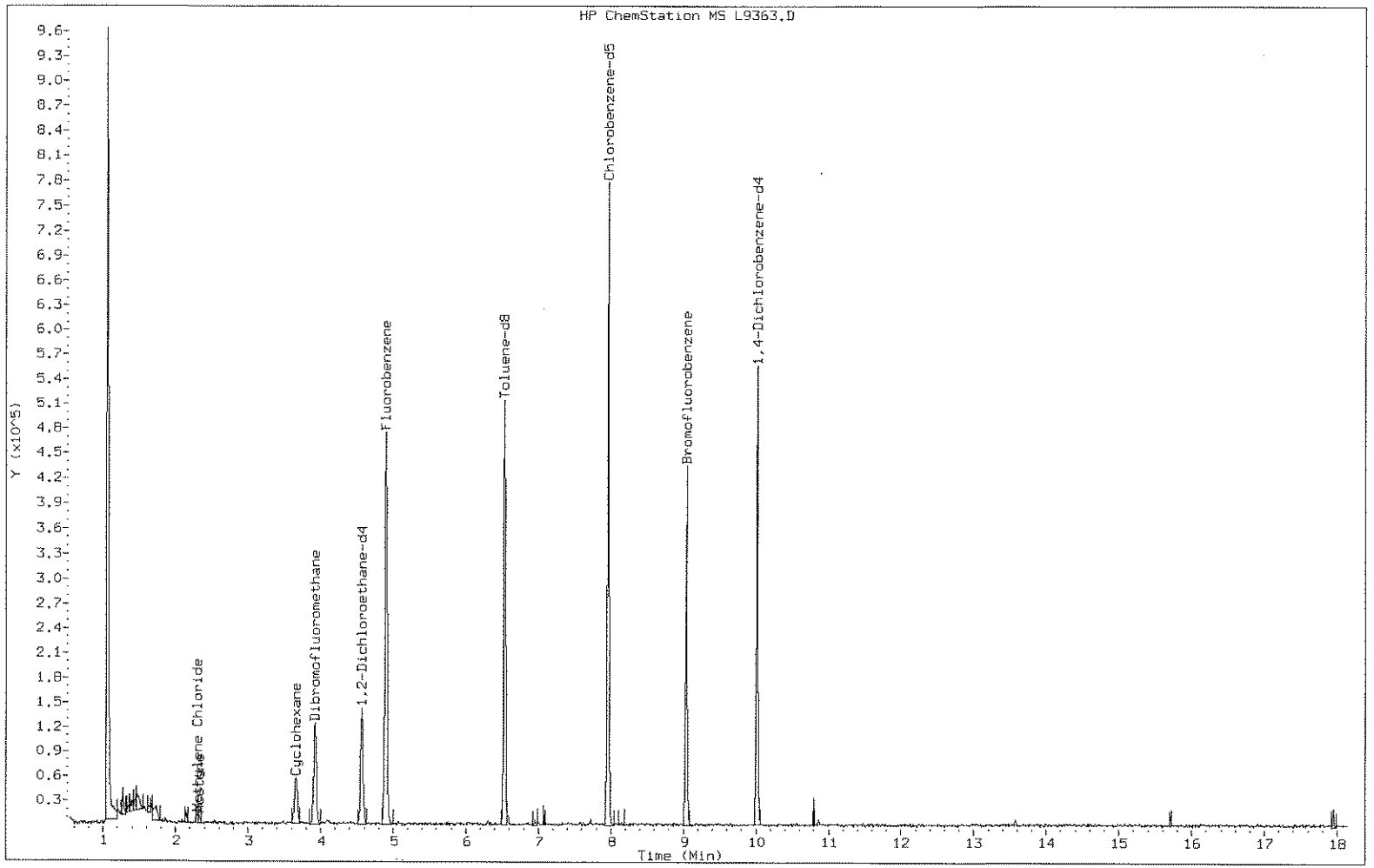
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.898	4.908	(1.000)	445250	25.0000	
	20 Methylene Chloride	84	2.300	2.300	(0.470)	3959	0.75893	0.8
	21 Acetone	43	2.330	2.330	(0.476)	8474	3.00102	3
	37 Cyclohexane	84	3.658	3.658	(0.747)	24900	3.82862	4
$	41 Dibromofluoromethane	111	3.924	3.924	(0.801)	106823	18.5968	18
$	55 1,2-Dichloroethane-d4	65	4.563	4.563	(0.932)	131817	18.8224	19
*	75 Chlorobenzene-d5	117	7.958	7.968	(1.000)	421068	25.0000	
	76 Toluene	91	6.580	6.580	(0.827)	6610	0.36867	0.4
$	77 Toluene-d8	98	6.531	6.541	(0.821)	313102	20.9123	21
*	95 1,4-Dichlorobenzene-d4	152	10.014	10.014	(1.000)	122252	25.0000	
$	125 Bromofluorobenzene	95	9.040	9.040	(0.903)	129750	28.2904	28

Data File: L9363.D

Date: 01-AUG-2007 19:44

Client ID: FB0726 Instrument: msl.i

Sample Info: 220-2277-A-25 Operator: D. HUMBERT



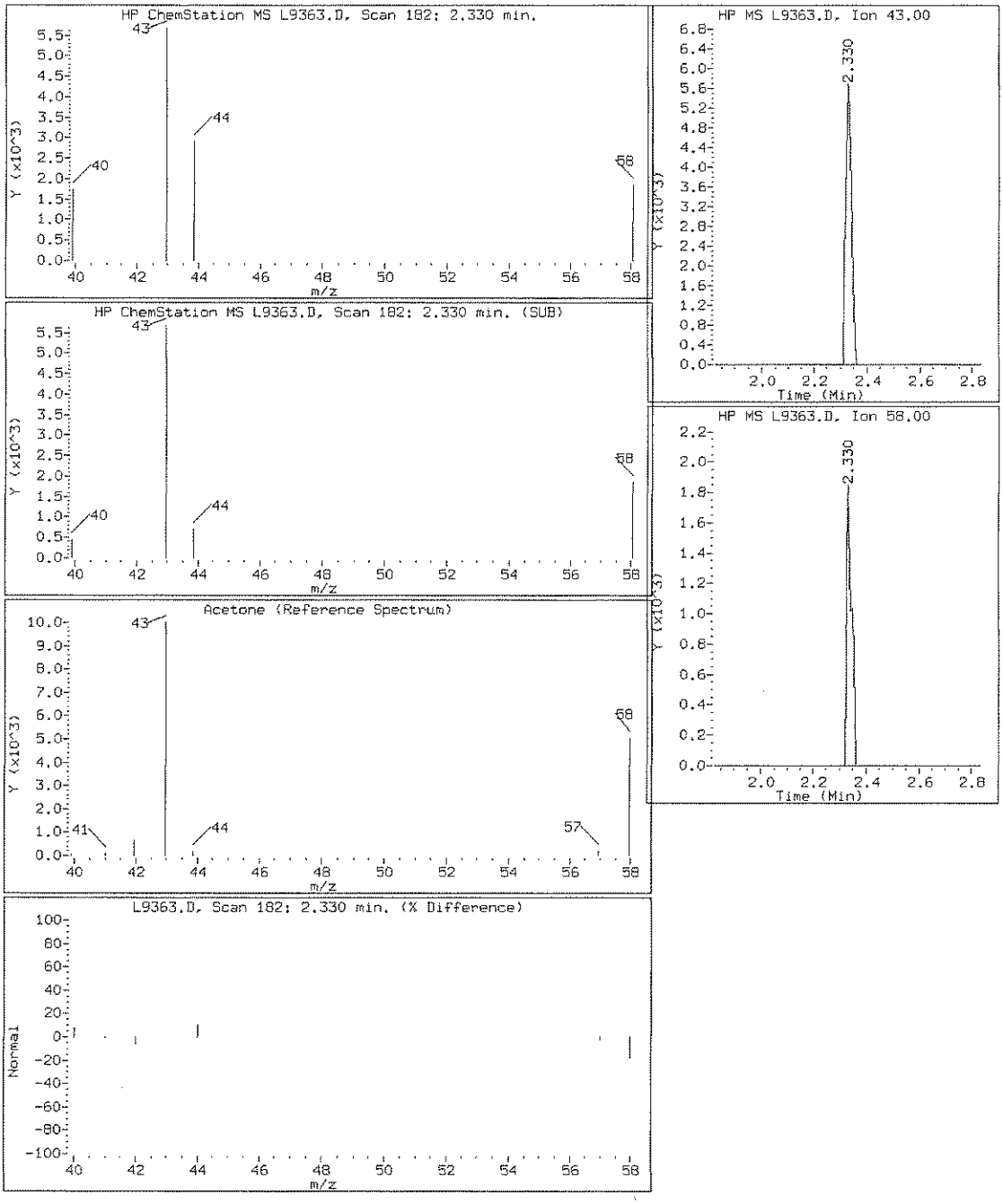
Data File: L9363.D

Date: 01-AUG-2007 19:44

Client ID: FB0726 Instrument: msl.i

Sample Info: 220-2277-A-25 Operator: D. HUMBERT

21 Acetone



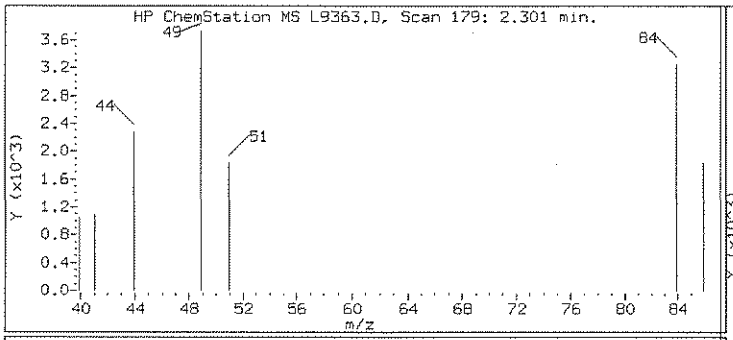
Data File: L9363.D

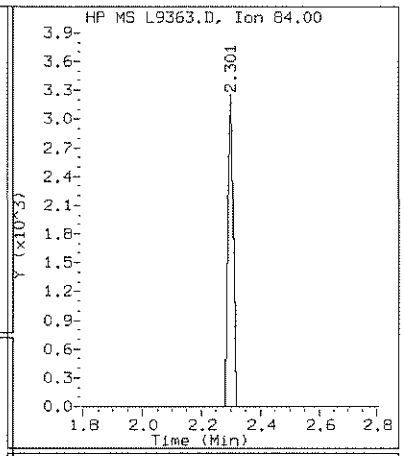
Date: 01-AUG-2007 19:44

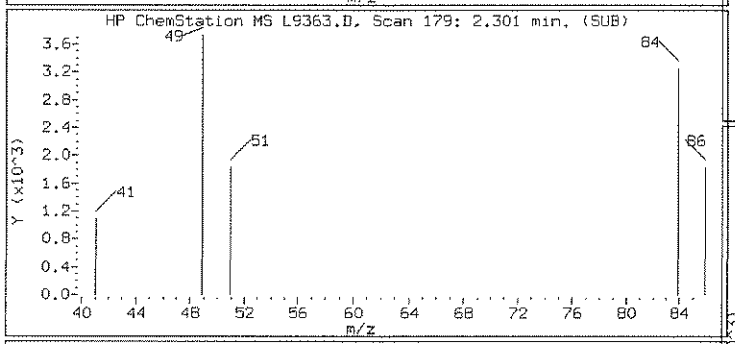
Client ID: FB0726 Instrument: msl.i

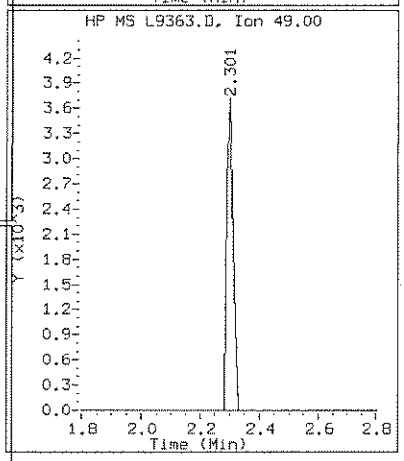
Sample Info: 220-2277-A-25 Operator: D. HUMBERT

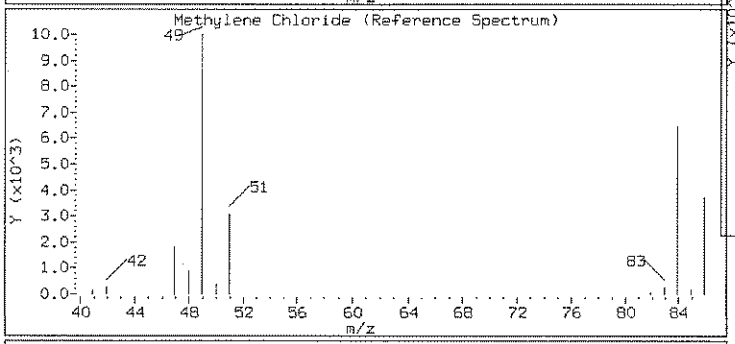
20 Methylene Chloride

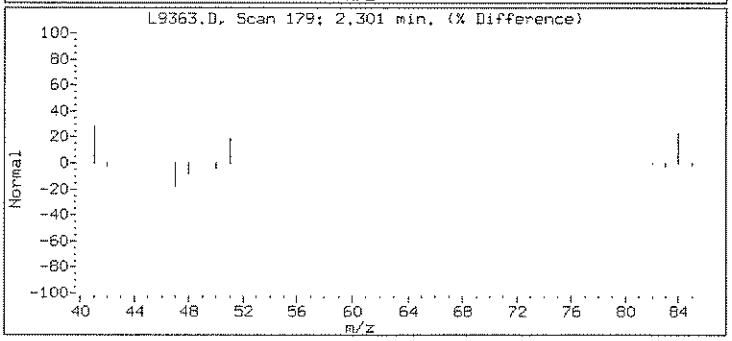












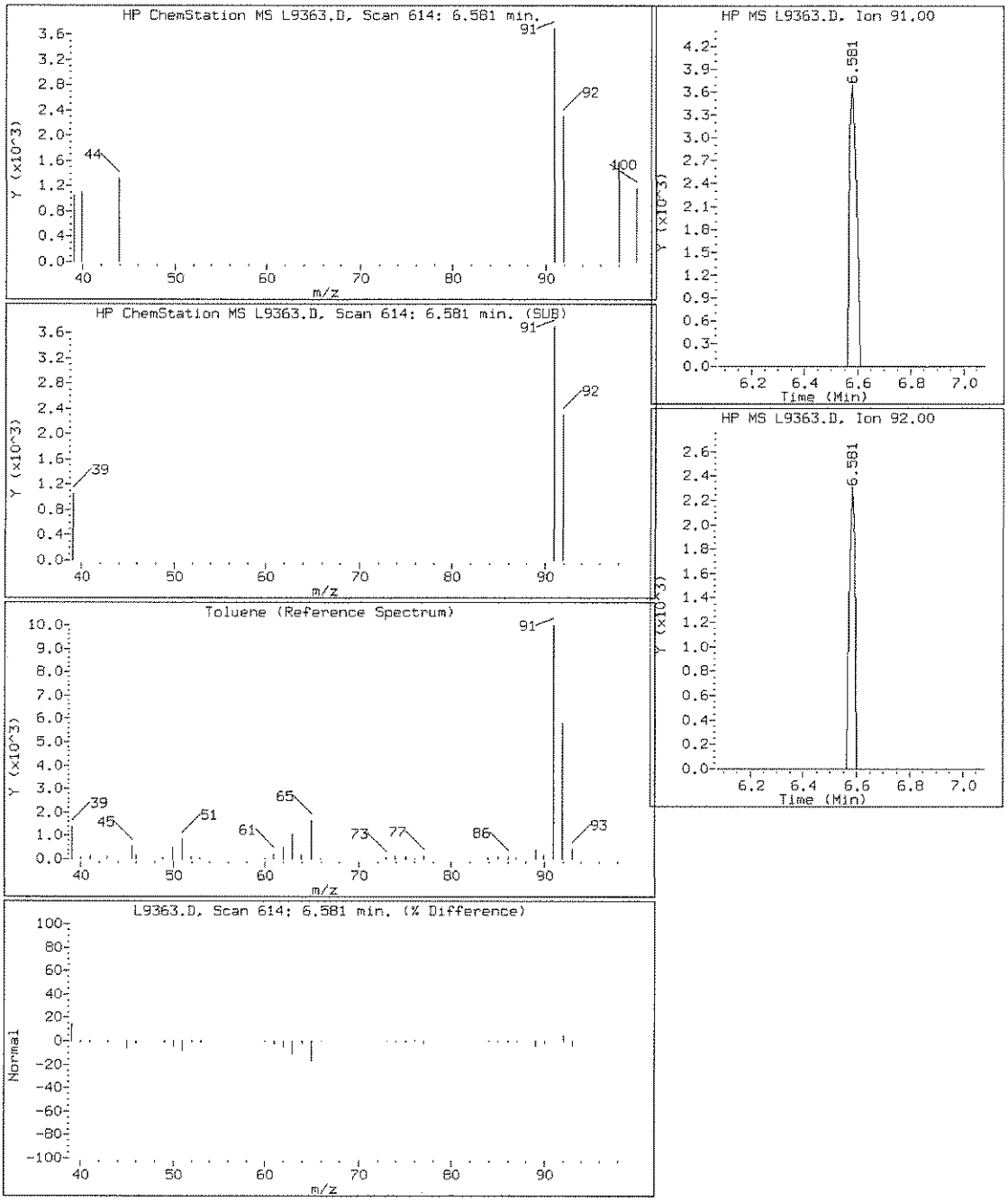
Data File: L9363.D

Date: 01-AUG-2007 19:44

Client ID: FB0726 Instrument: msl.i

Sample Info: 220-2277-A-25 Operator: D. HUMBERT

76 Toluene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-9D | Lab Sample ID: 220-2277-26

Matrix: Water | Lab File ID: L9364.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/01/2007 20:08

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8321 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	22		5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\Files\chem\VOA\msl.i\L079341.b\L9364.D
Lab Smp Id:
Inj Date : 01-AUG-2007 20:08 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-A-26
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\target1_ct\Files\chem\VOA\msl.i\L079341.b\L8260BNW.m
Meth Date : 01-Aug-2007 10:17 ctvoa Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 51
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONVOAMGR1

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
*	1 Fluorobenzene	96	4.899	4.908	(1.000)	445152	25.0000	
	38 Chloroform	83	3.718	3.707	(0.759)	4845	0.52088	0.5
$	41 Dibromofluoromethane	111	3.925	3.924	(0.801)	110875	19.3065	19
$	55 1,2-Dichloroethane-d4	65	4.564	4.563	(0.932)	126192	18.0231	18
*	75 Chlorobenzene-d5	117	7.959	7.968	(1.000)	426544	25.0000	
$	77 Toluene-d8	98	6.532	6.541	(0.821)	315562	20.8060	21
	80 Tetrachloroethene	164	6.955	6.964	(0.874)	80314	22.4102	22
*	95 1,4-Dichlorobenzene-d4	152	10.015	10.014	(1.000)	119100	25.0000	
$	125 Bromofluorobenzene	95	9.041	9.040	(0.903)	133326	29.8394	30

Data File: L9364.D

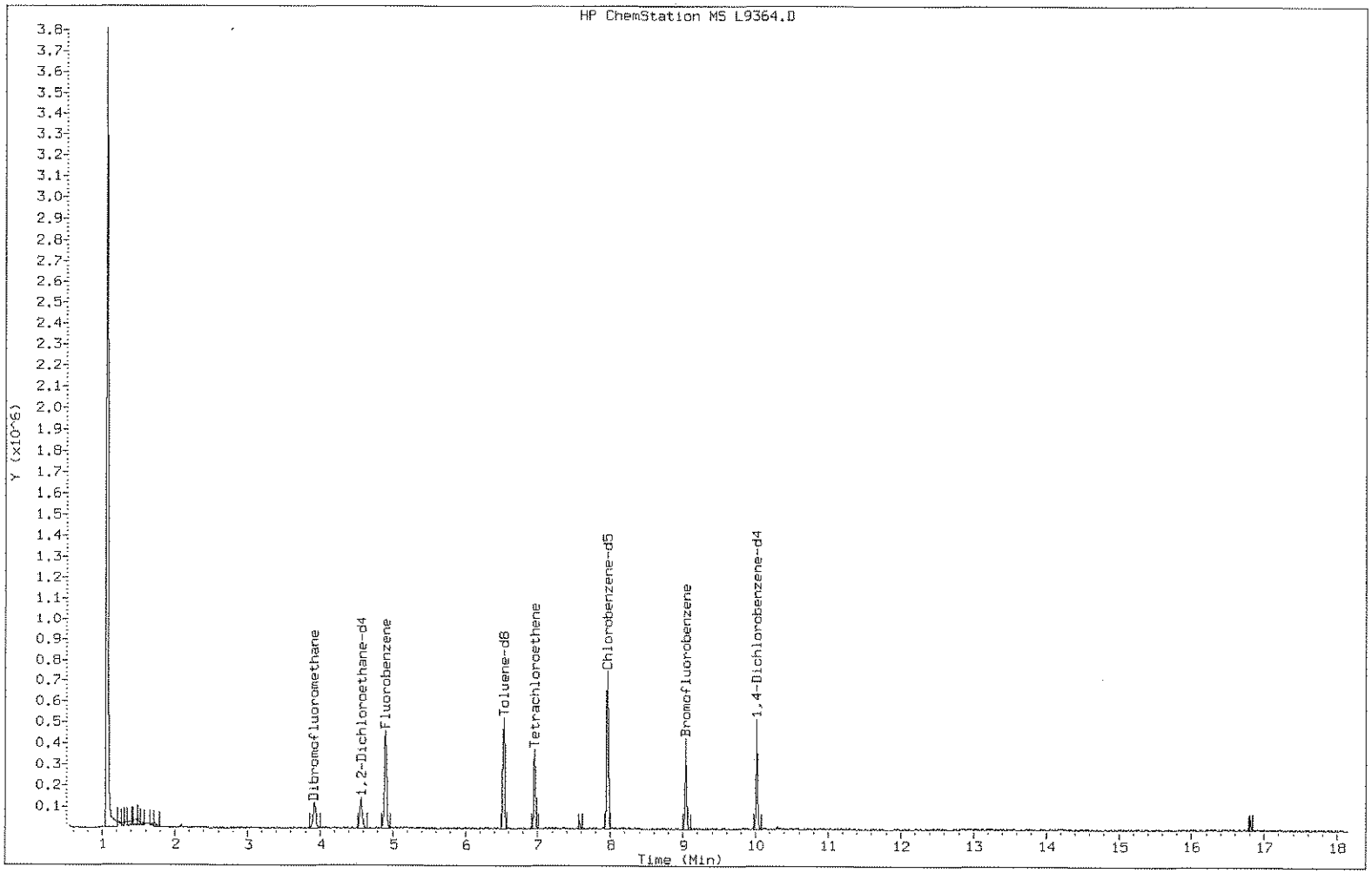
Date: 01-AUG-2007 20:08

Client ID: A-9D

Instrument: msl.i

Sample Info: 220-2277-A-26

Operator: D. HUMBERT



Data File: L9364.D

Date: 01-AUG-2007 20:08

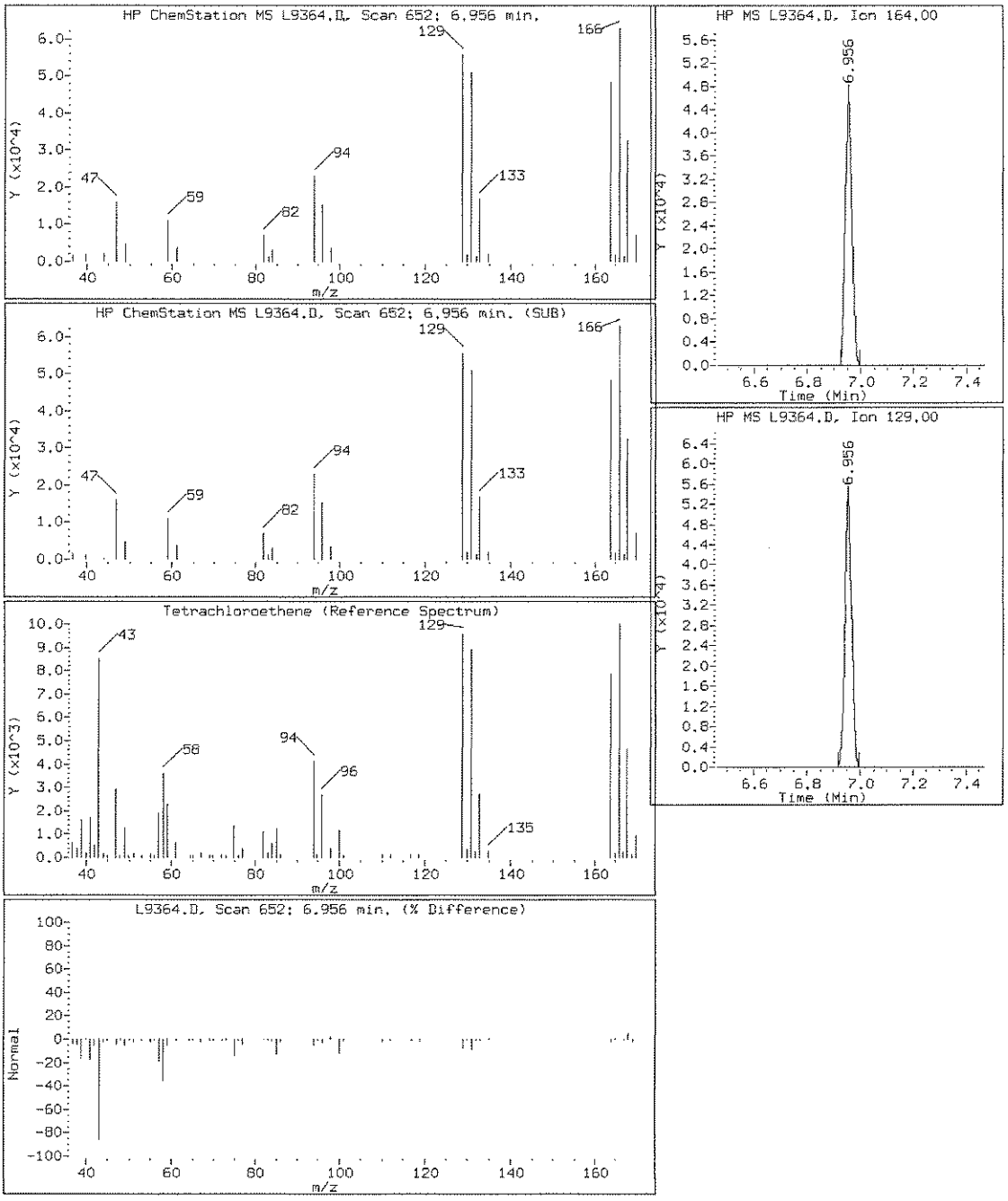
Client ID: A-9D

Instrument: msl.i

Sample Info: 220-2277-A-26

Operator: D. HUMBERT

80 Tetrachloroethene



08/09/2007

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: A-5 | Lab Sample ID: 220-2277-27

Matrix: Water | Lab File ID: L9365.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/01/2007 20:33

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8321 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	39		5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\Files\chem\VOA\msl.i\L079341.b\L9365.D
Lab Smp Id:
Inj Date : 01-AUG-2007 20:33 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-A-27
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\target1_ct\Files\chem\VOA\msl.i\L079341.b\L8260BNW.m
Meth Date : 01-Aug-2007 10:17 ctvoa Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 52
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONVOAMGR1

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

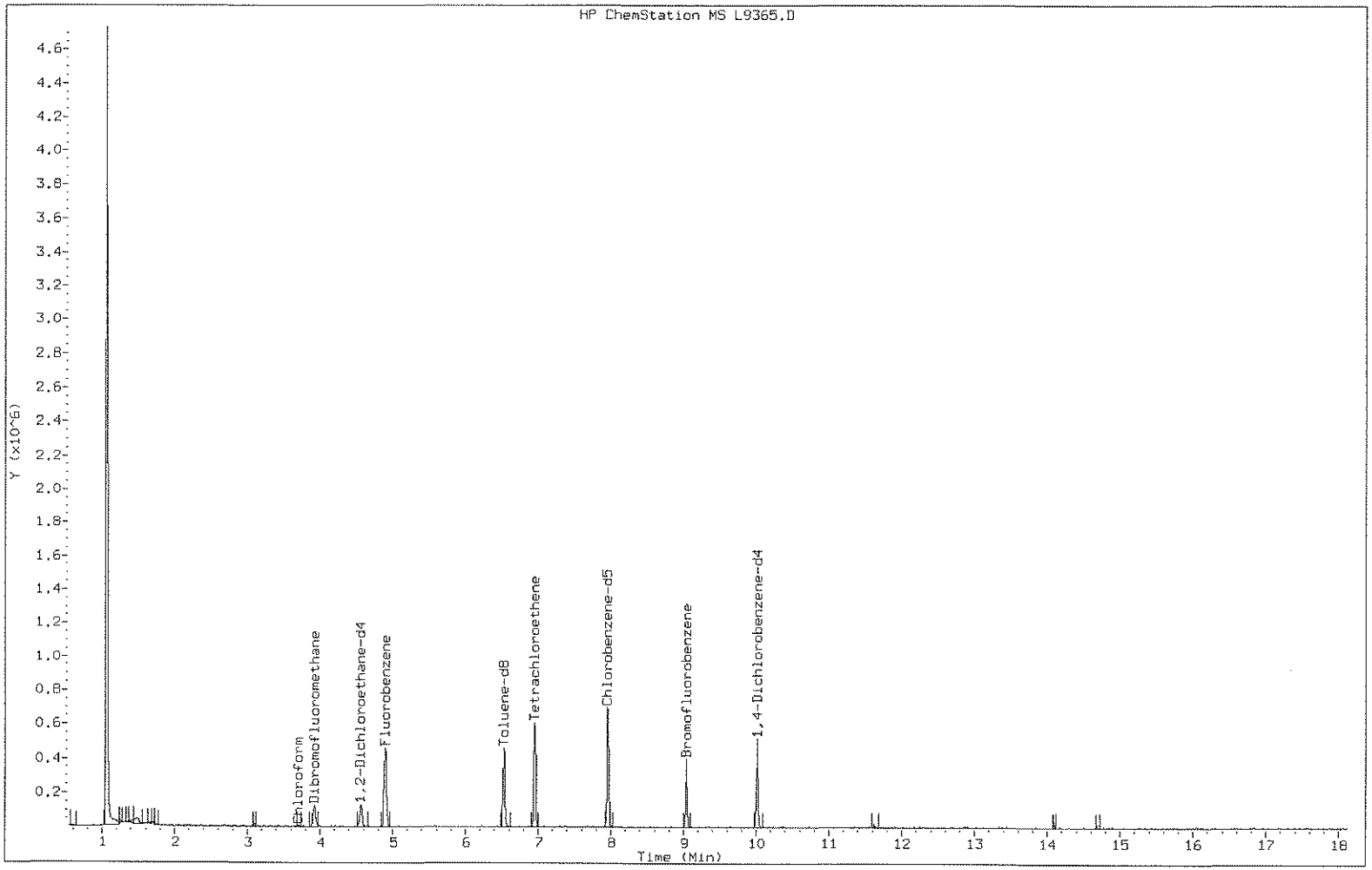
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1	Fluorobenzene	96	4.892	4.908	(1.000)	445548	25.0000	
38	Chloroform	83	3.701	3.707	(0.757)	4376	0.47004	0.5
$ 41	Dibromofluoromethane	111	3.918	3.924	(0.801)	107110	18.6343	19
$ 55	1,2-Dichloroethane-d4	65	4.567	4.563	(0.934)	123916	17.6823	18
* 75	Chlorobenzene-d5	117	7.962	7.968	(1.000)	426795	25.0000	
$ 77	Toluene-d8	98	6.525	6.541	(0.820)	302121	19.9081	20
80	Tetrachloroethene	164	6.958	6.964	(0.874)	141635	39.4976	39
* 95	1,4-Dichlorobenzene-d4	152	10.018	10.014	(1.000)	122697	25.0000	
$ 125	Bromofluorobenzene	95	9.034	9.040	(0.902)	122857	26.6903	27

Data File: L9365.D

Date: 01-AUG-2007 20:33

Client ID: A-5 Instrument: msl.i

Sample Info: 220-2277-A-27 Operator: D. HUMBERT



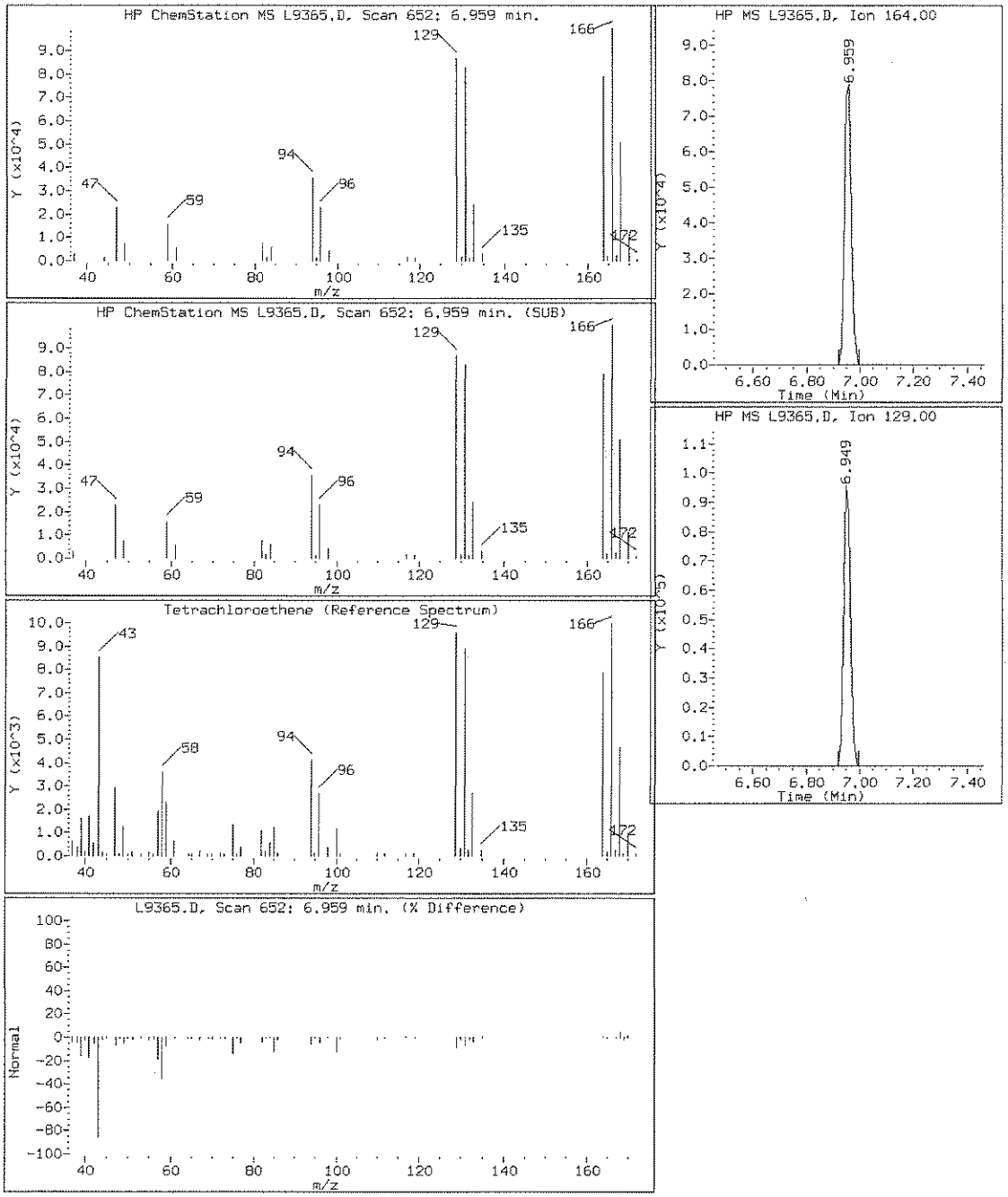
Data File: L9365.D

Date: 01-AUG-2007 20:33

Client ID: A-5 Instrument: msl.i

Sample Info: 220-2277-A-27 Operator: D. HUMBERT

80 Tetrachloroethene



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: ME-15 | Lab Sample ID: 220-2277-28

Matrix: Water | Lab File ID: L9384.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/02/2007 15:58

Level: (low/med) Low | Dilution Factor: 5

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8356 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	50	U	50	7.0
71-43-2	Benzene	25	U	25	2.0
75-27-4	Bromodichloromethane	25	U	25	2.0
75-25-2	Bromoform	25	U	25	4.0
74-83-9	Bromomethane	25	U	25	6.0
78-93-3	2-Butanone (MEK)	50	U	50	6.0
75-15-0	Carbon disulfide	25	U	25	4.5
56-23-5	Carbon tetrachloride	25	U	25	5.0
108-90-7	Chlorobenzene	25	U	25	2.0
75-00-3	Chloroethane	25	U	25	4.0
67-66-3	Chloroform	25	U	25	3.5
74-87-3	Chloromethane	25	U	25	2.5
124-48-1	Dibromochloromethane	25	U	25	2.5
75-34-3	1,1-Dichloroethane	25	U	25	3.0
107-06-2	1,2-Dichloroethane	25	U	25	3.0
75-35-4	1,1-Dichloroethene	25	U	25	3.5
78-87-5	1,2-Dichloropropane	25	U	25	4.5
10061-01-5	cis-1,3-Dichloropropene	25	U	25	2.5
10061-02-6	trans-1,3-Dichloropropene	25	U	25	4.0
100-41-4	Ethylbenzene	25	U	25	5.0
591-78-6	2-Hexanone	50	U	50	4.0
75-09-2	Methylene Chloride	25	U	25	2.0
108-10-1	4-Methyl-2-pentanone (MIBK)	50	U	50	3.5
100-42-5	Styrene	25	U	25	2.5
79-34-5	1,1,2,2-Tetrachloroethane	25	U	25	2.0
127-18-4	Tetrachloroethene	580		25	2.5
108-88-3	Toluene	25	U	25	1.5
71-55-6	1,1,1-Trichloroethane	25	U	25	2.0
79-00-5	1,1,2-Trichloroethane	25	U	25	3.0
79-01-6	Trichloroethene	25	U	25	3.5
75-01-4	Vinyl chloride	25	U	25	4.0
1330-20-7	Xylenes, Total	25	U	25	5.0
156-59-2	cis-1,2-Dichloroethene	25	U	25	3.0
156-60-5	trans-1,2-Dichloroethene	25	U	25	2.5

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079370.b\L9384.D
Lab Smp Id: 220-2277-B-28 Client Smp ID: ME-15
Inj Date : 02-AUG-2007 15:58 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-B-28
Misc Info : : ;;; ; 8260 ; 5 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 66
Dil Factor: 5.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	5.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

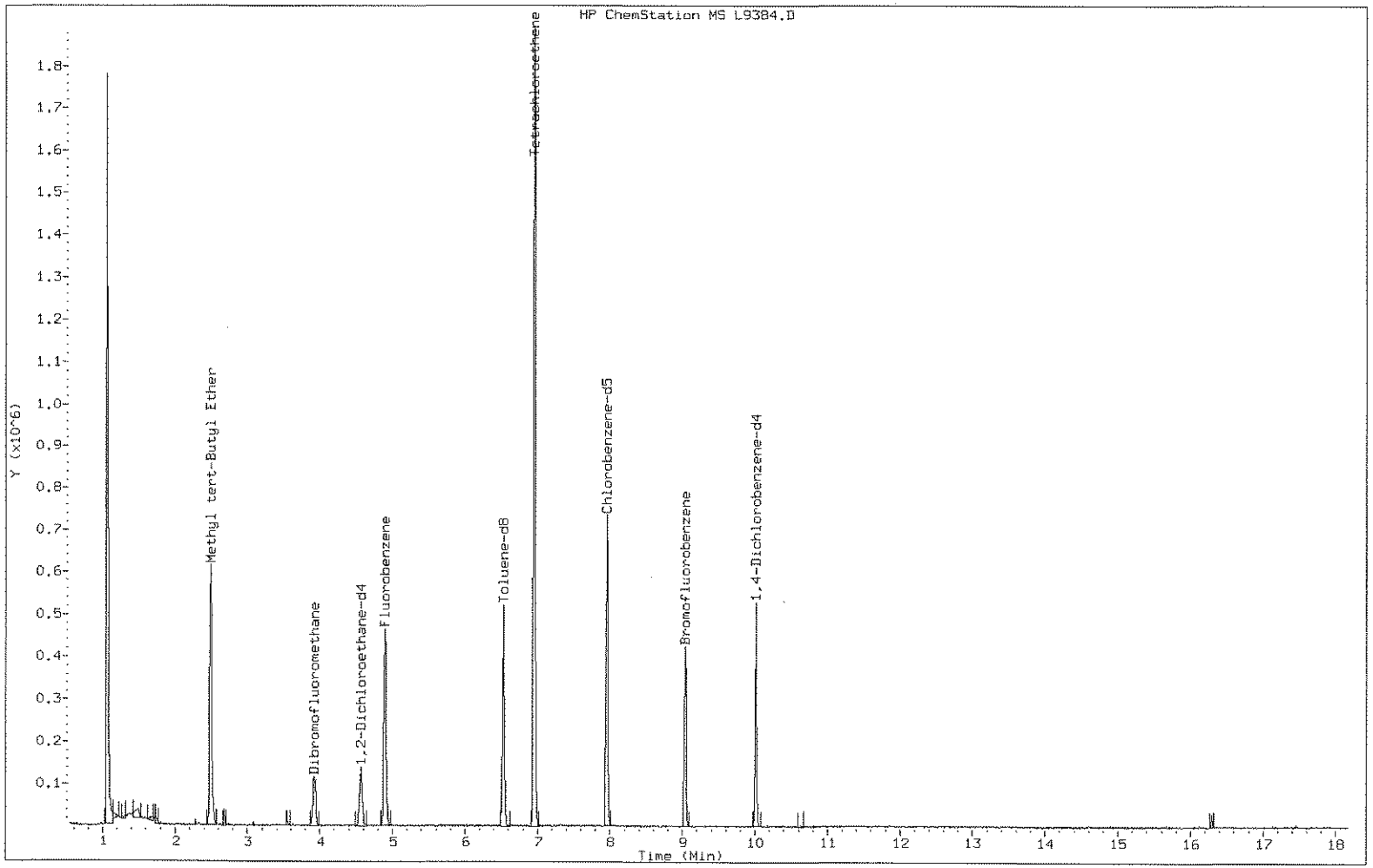
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1	Fluorobenzene	96	4.892	4.901	(1.000)	445725	25.0000	
24	Methyl tert-Butyl Ether	73	2.491	2.491	(0.509)	649362	35.0905	180
$ 41	Dibromofluoromethane	111	3.908	3.927	(0.799)	110340	19.1886	19
$ 55	1,2-Dichloroethane-d4	65	4.557	4.567	(0.932)	131994	18.8275	19
* 75	Chlorobenzene-d5	117	7.952	7.961	(1.000)	425683	25.0000	
$ 77	Toluene-d8	98	6.525	6.535	(0.821)	319754	21.1250	21
80	Tetrachloroethene	164	6.948	6.958	(0.874)	411887	115.162	580
* 95	1,4-Dichlorobenzene-d4	152	10.008	10.018	(1.000)	118200	25.0000	
$ 125	Bromofluorobenzene	95	9.034	9.043	(0.903)	131838	29.7311	30

Data File: L9384.D

Date: 02-AUG-2007 15:58

Client ID: ME-15 Instrument: ms1.i

Sample Info: 220-2277-B-28 Operator: D. HUMBERT



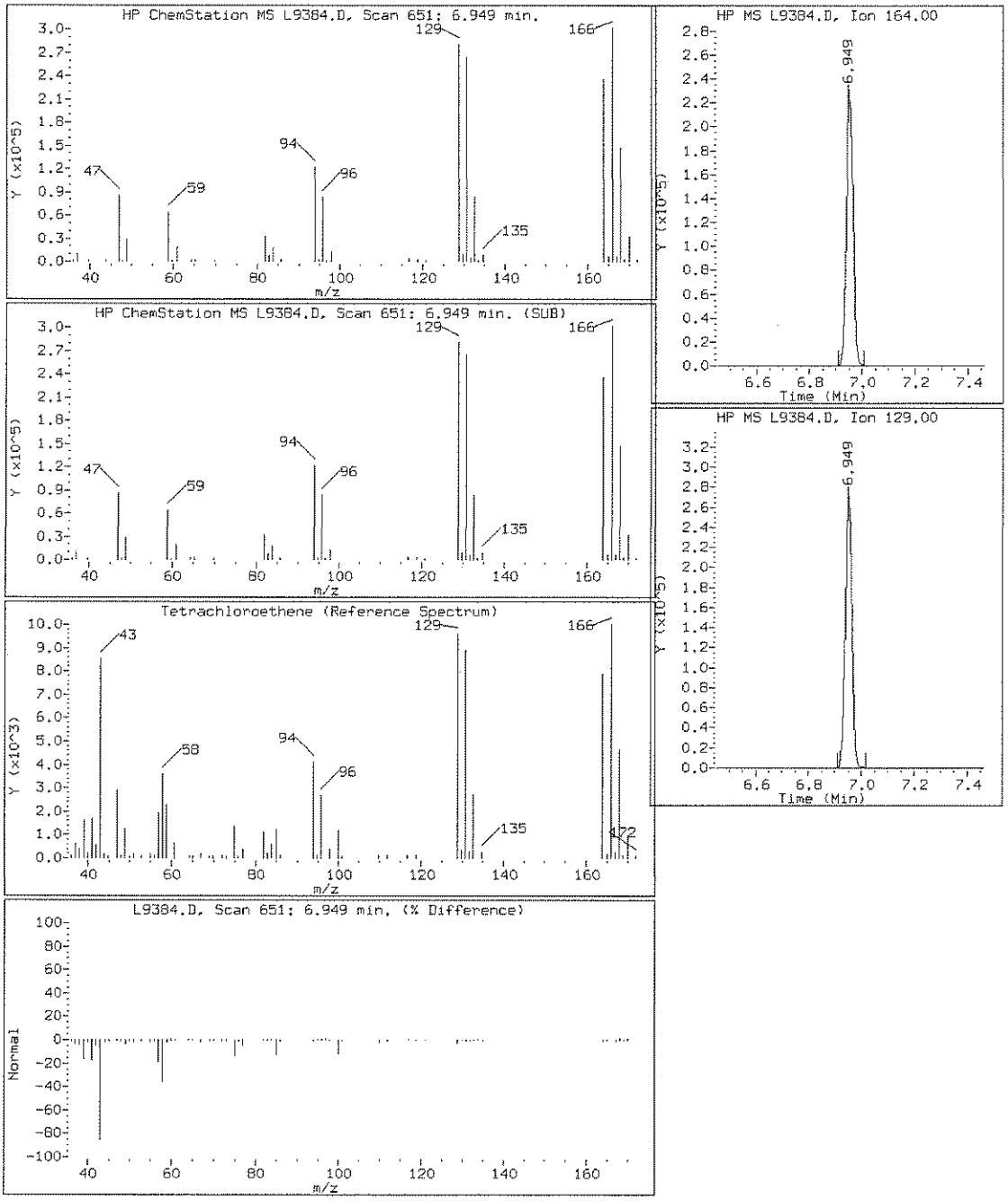
Data File: L9384.D

Date: 02-AUG-2007 15:58

Client ID: ME-15 Instrument: msl.i

Sample Info: 220-2277-B-28 Operator: D. HUMBERT

80 Tetrachloroethene



Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Files:	Lab Sample ID	Lab File ID	Batch	Sample Number
	IC 220-8212/1	L9152.D	8212	1
	IC 220-8212/2	L9153.D	8212	2
	IC 220-8212/3	L9154.D	8212	3
	IC 220-8212/4	L9155.D	8212	4
	IC 220-8212/5	L9156.D	8212	5

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	RRF/RESPONSE					Curve Type	Coefficients		
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5		b	m1	m2
Hexachloroethane	DCB						Ave			
1,2-Dichlorobenzene	DCB	1.3128	1.2962	1.3593	1.3602	1.3715	Ave		1.3400	
Carbon disulfide	FB	0.9713	0.9704	1.0469	1.0917	1.1198	Ave		1.0400	
n-Heptane	FB	0.3206	0.2863	0.3116	0.3111	0.3155	Ave		0.3090	
Iodomethane	FB	0.1394	0.2105	0.3338	0.3435	0.3546	Ave		0.2764	
2-Chlorotoluene	DCB	2.0902	2.3749	2.1233	2.3723	2.4250	Ave		2.2771	
1-Chlorohexane	CBZ	0.3190	0.4127	0.3715	0.3440	0.3555	Ave		0.3605	
trans-1,4-Dichloro-2-butene	DCB	0.2826	0.3410	0.3636	0.3812	0.3623	Ave		0.3461	
1,2,3-Trichloropropane	DCB	0.3296	0.3434	0.3402	0.3388	0.3334	Ave		0.3371	
Carbon tetrachloride	FB	0.3616	0.3777	0.4799	0.4273	0.5041	Ave		0.4301	
Nitrobenzene	DCB	0.0130	0.0268	0.0427	0.0650	0.0890	Ave		0.0473	

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	RRF/RESPONSE					Curve Type	Coefficients		
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5		b	m1	m2
2-Hexanone	CBZ	0.3031	0.3313	0.3573	0.3611	0.3634	Ave		0.3432	
Cyclohexane	FB	0.3284	0.3357	0.3749	0.3868	0.4000	Ave		0.3652	
Xylenes, Total	CBZ	0.4490	0.4495	0.4759	0.4977	0.5000	Ave		0.4744	
3-Chloro-1-propene	FB	0.5167	0.5272	0.5632	0.5712	0.5715	Ave		0.5499	
Ethyl acetate	FB	0.0244	0.0283	0.0288	0.0299	0.0285	Ave		0.0280	
cis-1,3-Dichloropropene	FB	0.5240	0.5286	0.5709	0.5943	0.6141	Ave		0.5664	
n-Propyl acetate	FB	0.0223	0.0318	0.0326	0.0350	0.0364	Ave		0.0316	
Chlorobenzene	CBZ	0.7857	0.8421	0.8745	0.9027	0.9189	Ave		0.8648	
Vinyl chloride	FB	0.3298	0.3675	0.3764	0.3839	0.3931	Ave		0.3701	
Methyl acrylate	FB	0.3516	0.3731	0.4172	0.4295	0.4303	Ave		0.4003	
Propionitrile	FB	0.0576	0.0594	0.0659	0.0666	0.0677	Ave		0.0634	
sec-Butylbenzene	DCB	2.6180	2.5159	2.5996	2.6197	2.6978	Ave		2.6102	
Dibromomethane	FB	0.1815	0.1809	0.1958	0.1978	0.2059	Ave		0.1924	
1,2,4,5-Tetramethylbenzene	FB	0.6166	0.7380	0.8397	0.9621	0.9861	Ave		0.8285	
p-Diethylbenzene	FB	0.4379	0.4907	0.5737	0.6054	0.6150	Ave		0.5445	

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	RRF/RESPONSE					Curve Type	Coefficients		
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5		b	m1	m2
Dibromofluoromethane	FB	0.2874	0.3173	0.3281	0.3300	0.3498	Ave		0.3225	
Acetone	FB	0.1802	0.1500	0.1636	0.1490	0.1500	Ave		0.1585	
Isopropyl alcohol	FB	0.0465	0.0539	0.0564	0.0628	0.0672	Ave		0.0574	
m-Xylene & p-Xylene	CBZ	0.4612	0.4529	0.4831	0.5025	0.5061	Ave		0.4812	
Isopropyl ether	FB	1.1434	1.1642	1.2143	1.2440	1.2710	Ave		1.2074	
1,2,4-Trichlorobenzene	DCB	0.4571	0.6266	0.7807	0.8613	0.9143	Ave		0.7280	
tert-Butyl Formate	FB	0.3358	0.3305	0.3591	0.3625	0.3616	Ave		0.3499	
2-Nitropropane	FB	0.1087	0.1036	0.1107	0.1131	0.1199	Ave		0.1112	
2-Methyl-2-propanol	FB	0.0560	0.0530	0.0577	0.0617	0.0610	Ave		0.0579	
Styrene	CBZ	0.6982	0.7352	0.7731	0.8279	0.8262	Ave		0.7721	
Chlorobromomethane	FB	0.1912	0.1964	0.2108	0.2108	0.2144	Ave		0.2047	
Dichlorobromomethane	FB	0.3253	0.3524	0.3805	0.4027	0.4098	Ave		0.3741	
1,3-Dichlorobenzene	DCB	1.3204	1.3211	1.3913	1.3801	1.3953	Ave		1.3616	
Benzene	FB	1.1785	1.1221	1.2065	1.2324	1.2612	Ave		1.2001	
Pentachloroethane	DCB						Ave			

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	RRF/RESPONSE					Curve Type	Coefficients		
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5		b	m1	m2
Chloroethane	FB	0.2186	0.2428	0.2365	0.1659	0.1170	Ave		0.1962	
2-Chloroethyl vinyl ether	FB	0.0758	0.0917	0.0988	0.1060	0.1102	Ave		0.0965	
trans-1,3-Dichloropropene	FB	0.4479	0.4845	0.5322	0.5428	0.5631	Ave		0.5141	
Acrolein	FB	0.0471	0.0537	0.0612	0.0646	0.0635	Ave		0.0580	
1,2,3-Trichlorobenzene	DCB	0.3916	0.5854	0.7045	0.7887	0.8511	Ave		0.6643	
N-Propylbenzene	DCB	3.2908	3.1601	3.2851	3.3678	3.4158	Ave		3.3039	
o-Xylene	CBZ	0.4246	0.4426	0.4616	0.4881	0.4877	Ave		0.4609	
Tetrahydrofuran	FB	0.1141	0.1351	0.1373	0.1458	0.1450	Ave		0.1355	
Isobutanol	FB	0.0143	0.0148	0.0166	0.0172	0.0179	Ave		0.0162	
2-Chloro-1,3-butadiene	FB	0.1867	0.1796	0.1982	0.1979	0.2051	Ave		0.1935	
Ethanol	FB	0.0166	0.0165	0.0177	0.0187	0.0193	Ave		0.0178	
4-Bromofluorobenzene	DCB	1.0327	0.8850	0.9443	0.8932	0.9343	Ave		0.9379	
4-Isopropyltoluene	DCB	2.7138	2.7223	2.7877	2.8105	2.9382	Ave		2.7945	
n-Butyl acetate	CBZ	0.2373	0.2775	0.2985	0.3236	0.3137	Ave		0.2901	
Methacrylonitrile	FB	0.4605	0.4478	0.4761	0.4788	0.3712	Ave		0.4469	

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	RRF/RESPONSE					Curve Type	Coefficients		
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5		b	m1	m2
n-Butylbenzene	DCB	2.9515	3.0396	3.6489	3.4886	3.7072	Ave		3.3672	
1,1-Dichloropropene	FB	0.4433	0.4490	0.4880	0.5011	0.5068	Ave		0.4776	
cis-1,2-Dichloroethene	FB	0.2566	0.2806	0.2909	0.2983	0.3034	Ave		0.2859	
1,1,2,2-Tetrachloroethane	DCB	1.0402	1.0537	1.0784	1.0704	1.0765	Ave		1.0638	
1,2,4-Trimethylbenzene	DCB	2.4036	2.3988	2.5401	2.4952	2.5646	Ave		2.4805	
Toluene	CBZ	1.0065	1.0242	1.0628	1.1070	1.1221	Ave		1.0645	
Naphthalene	DCB	1.0484	1.7625	2.3560	2.7918	3.0753	Ave		2.2068	
1,3,5-Trimethylbenzene	DCB	2.4933	2.3788	2.4702	2.4914	2.5132	Ave		2.4694	
1,1-Dichloroacetone	CBZ	0.2305	0.2454	0.2651	0.2809	0.2834	Ave		0.2611	
1,3-Dichloropropane	CBZ	0.5503	0.5579	0.5749	0.5939	0.6098	Ave		0.5774	
Chloroform	FB	0.4726	0.5062	0.5351	0.5448	0.5532	Ave		0.5224	
4-Chlorotoluene	DCB	2.1409	2.0536	2.1819	2.1562	2.2007	Ave		2.1467	
Chlorodibromomethane	CBZ	0.3616	0.3747	0.4061	0.4208	0.4367	Ave		0.4000	
Dichlorodifluoromethane	FB	0.1704	0.1909	0.1958	0.2016	0.2011	Ave		0.1920	
1,1,2-Trichloroethane	FB	0.2328	0.2408	0.2499	0.2619	0.2656	Ave		0.2502	

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	RRF/RESPONSE					Curve Type	Coefficients		
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5		b	m1	m2
Tert-amyl methyl ether	FB	0.9514	1.0206	1.0647	1.0980	1.1275	Ave		1.0524	
Toluene-d8 (Surr)	CBZ	0.8677	0.8587	0.8637	0.9102	0.9444	Ave		0.8889	
tert-Butylbenzene	DCB	2.3987	2.2649	2.4139	2.3926	2.4718	Ave		2.3884	
1,1-Dichloro-1-fluoroethane	FB	0.4104	0.4288	0.4513	0.4586	0.4617	Ave		0.4422	
Chloromethane	FB	0.3030	0.3316	0.3381	0.3504	0.3582	Ave		0.3363	
Methylene Chloride	FB	0.2720	0.2789	0.3021	0.3030	0.3085	Ave		0.2929	
Isopropyl acetate	FB	0.1445	0.1001	0.1366	0.1450	0.1447	Ave		0.1342	
4-Ethyltoluene	DCB	3.0082	2.8899	3.0174	3.0358	3.1168	Ave		3.0136	
Methyl methacrylate	FB	0.1515	0.1692	0.1759	0.1845	0.1868	Ave		0.1736	
1,1-Dichloroethene	FB	0.2009	0.2016	0.2143	0.2243	0.2242	Ave		0.2131	
Isopropylbenzene	DCB	3.1784	2.9837	3.0385	3.0780	3.1364	Ave		3.0830	
1,2-Dichloroethane	FB	0.4451	0.4596	0.4815	0.4876	0.4995	Ave		0.4746	
1,2-Dichloroethane-d4 (Surr)	FB	0.3705	0.3793	0.4106	0.3906	0.4151	Ave		0.3932	
Acrylonitrile	FB	0.1628	0.1935	0.1796	0.1910	0.1901	Ave		0.1834	
methyl isobutyl ketone	CBZ	0.4029	0.4534	0.4774	0.4914	0.4761	Ave		0.4602	

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	RRF/RESPONSE					Curve Type	Coefficients		
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5		b	m1	m2
Tetrachloroethene	CBZ	0.1898	0.2036	0.2153	0.2219	0.2197	Ave		0.2100	
1,2-Dichloroethene, Total	FB	0.2581	0.2648	0.2791	0.2884	0.2934	Ave		0.2768	
1,1,1-Trichloroethane	FB	0.3517	0.3633	0.3912	0.4014	0.4121	Ave		0.3840	
2,2-Dichloropropane	FB	0.4963	0.4390	0.4743	0.4679	0.4778	Ave		0.4711	
Benzyl chloride	DCB	0.3454	0.3867	0.4228	0.4521	0.4645	Ave		0.4143	
Ethylene Dibromide	CBZ	0.3022	0.3207	0.3452	0.3567	0.3582	Ave		0.3366	
1,1,2-Trichloro-1,2,2-trifluoroethane	FB	0.2254	0.2339	0.2517	0.2609	0.2655	Ave		0.2475	
Methyl acetate	FB	1.7608	1.7939	1.8847	1.9359	1.9667	Ave		1.8684	
Bromoform	CBZ	0.1752	0.2074	0.2321	0.2425	0.2462	Ave		0.2207	
1,2-Dibromo-3-Chloropropane	DCB	0.1465	0.2121	0.2176	0.2379	0.2480	Ave		0.2124	
Trichlorofluoromethane	FB	0.3238	0.3575	0.3646	0.3682	0.3778	Ave		0.3584	
Trichloroethene	FB	0.3339	0.3329	0.3580	0.3677	0.3728	Ave		0.3531	
Methylcyclohexane	FB	0.3365	0.3349	0.3484	0.3590	0.3705	Ave		0.3498	
Bromobenzene	DCB	0.7967	0.7504	0.7863	0.7772	0.7818	Ave		0.7785	
Methyl Ethyl Ketone	FB	0.2268	0.2186	0.2392	0.2317	0.2345	Ave		0.2301	

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	RRF/RESPONSE					Curve Type	Coefficients		
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5		b	m1	m2
1,2-Dichloropropane	FB	0.3657	0.3847	0.4007	0.4076	0.4213	Ave		0.3960	
1,4-Dioxane	FB	0.0023	0.0025	0.0032	0.0035	0.0035	Ave		0.0030	
1-Chlorobutane	FB	0.6560	0.6445	0.6951	0.7123	0.7194	Ave		0.6854	
n-Butanol	FB	0.0113	0.0117	0.0130	0.0143	0.0148	Ave		0.0130	
Tert-butyl ethyl ether	FB	1.1234	1.1509	1.2528	1.2593	1.2795	Ave		1.2132	
Methyl tert-butyl ether	FB	0.9559	1.0005	1.0565	1.0812	1.0956	Ave		1.0379	
1,4-Dichlorobenzene	DCB	1.3087	1.3501	1.4137	1.4234	1.4528	Ave		1.3897	
1,1,1,2-Tetrachloroethane	CBZ	0.2847	0.3031	0.3144	0.3245	0.3371	Ave		0.3128	
Ethylbenzene	CBZ	0.3376	0.3701	0.3903	0.4124	0.4144	Ave		0.3850	
trans-1,2-Dichloroethene	FB	0.2596	0.2491	0.2674	0.2785	0.2834	Ave		0.2676	
Hexachlorobutadiene	DCB	0.2459	0.2193	0.2439	0.2638	0.2674	Ave		0.2481	
Dichlorofluoromethane	FB	0.6022	0.6197	0.6665	0.6837	0.6979	Ave		0.6540	
1,1-Dichloroethane	FB	0.6336	0.6630	0.6919	0.7005	0.7257	Ave		0.6829	
Ethyl ether	FB	0.2144	0.2254	0.2360	0.2266	0.2250	Ave		0.2255	
1-Bromopropane	FB	0.4959	0.4897	0.5212	0.5292	0.5391	Ave		0.5150	

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	RRF/RESPONSE					Curve Type	Coefficients		
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5		b	m1	m2
Acetonitrile	FB	0.0610	0.0626	0.0676	0.0700	0.0695	Ave		0.0661	
Vinyl acetate	FB	0.8327	0.8699	0.9326	0.9537	0.9678	Ave		0.9114	
Ethyl methacrylate	CBZ	0.5001	0.5606	0.5671	0.6069	0.5986	Ave		0.5666	
Bromomethane	FB	0.1354	0.1203	0.1221	0.1393	0.1292	Ave		0.1293	
Chloroacetonitrile	FB	0.0141	0.0145	0.0164	0.0178	0.0182	Ave		0.0162	

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	Amount ug/L					Curve Evaluation						Min R^2 or COD
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5	Curve Type	Ave RRF	Min Ave RRF	%RSD	Max %RSD	R^2 or COD	
Hexachloroethane	DCB	5.00	20.00	50.00	100.00	200.00	Ave						
1,2-Dichlorobenzene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	1.3400		2.5			
Carbon disulfide	FB	5.00	20.00	50.00	100.00	200.00	Ave	1.0400		6.6			
n-Heptane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.3090		4.3			
Iodomethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.2764		34.8			
2-Chlorotoluene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	2.2771		6.9			
1-Chlorohexane	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	0.3605		9.7			
trans-1,4-Dichloro-2-butene	DCB	10.00	40.00	100.00	200.00	400.00	Ave	0.3461		11.1			
1,2,3-Trichloropropane	DCB	5.00	20.00	50.00	100.00	200.00	Ave	0.3371		1.6			
Carbon tetrachloride	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.4301		14.4			
Nitrobenzene	DCB	50.00	200.00	500.00	1000.00	2000.00	Ave	0.0473		64.0			
2-Hexanone	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	0.3432		7.5			
Cyclohexane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.3652		8.7			
Xylenes, Total	CBZ	15.00	60.00	150.00	300.00	600.00	Ave	0.3953		5.2			

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	Amount ug/L					Curve Evaluation						Min R^2 or COD
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5	Curve Type	Ave RRF	Min Ave RRF	%RSD	Max %RSD	R^2 or COD	
3-Chloro-1-propene	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.5499		4.7			
Ethyl acetate	FB	10.00	40.00	100.00	200.00	400.00	Ave	0.0280		7.5			
cis-1,3-Dichloropropene	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.5664		7.0			
n-Propyl acetate	FB	10.00	40.00	100.00	200.00	400.00	Ave	0.0316		17.4			
Chlorobenzene	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	0.8648	0.3000	6.1			
Vinyl chloride	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.3701		6.6	30.0		
Methyl acrylate	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.4003		9.0			
Propionitrile	FB	50.00	200.00	500.00	1000.00	2000.00	Ave	0.0634		7.3			
sec-Butylbenzene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	2.6102		2.5			
Dibromomethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.1924		5.7			
1,2,4,5-Tetramethylbenzene	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.8285		18.7			
p-Diethylbenzene	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.5445		14.2			
Dibromofluoromethane	FB	5.00	20.00	25.00	100.00	200.00	Ave	0.3225		7.1			
Acetone	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.1585		8.5			
Isopropyl alcohol	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.0574		14.0			

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	Amount ug/L					Curve Evaluation						Min R^2 or COD
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5	Curve Type	Ave RRF	Min Ave RRF	%RSD	Max %RSD	R^2 or COD	
m-Xylene & p-Xylene	CBZ	10.00	40.00	100.00	200.00	400.00	Ave	0.4812		5.0			
Isopropyl ether	FB	5.00	20.00	50.00	100.00	200.00	Ave	1.2074		4.4			
1,2,4-Trichlorobenzene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	0.7280		25.6			
tert-Butyl Formate	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.3499		4.4			
2-Nitropropane	FB	10.00	40.00	100.00	200.00	400.00	Ave	0.1112		5.4			
2-Methyl-2-propanol	FB	25.00	100.00	250.00	500.00	1000.00	Ave	0.0579		6.2			
Styrene	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	0.7721		7.3			
Chlorobromomethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.2047		5.0			
Dichlorobromomethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.3741		9.4			
1,3-Dichlorobenzene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	1.3616		2.8			
Benzene	FB	5.00	20.00	50.00	100.00	200.00	Ave	1.2001		4.4			
Pentachloroethane	DCB	5.00	20.00	50.00	100.00	200.00	Ave						
Chloroethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.1962		27.3			
2-Chloroethyl vinyl ether	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.0965		14.0			
trans-1,3-Dichloropropene	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.5141		9.1			

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	Amount ug/L					Curve Evaluation						Min R^2 or COD
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5	Curve Type	Ave RRF	Min Ave RRF	%RSD	Max %RSD	R^2 or COD	
Acrolein	FB	25.00	100.00	250.00	500.00	1000.00	Ave	0.0580		12.8			
1,2,3-Trichlorobenzene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	0.6643		27.4			
N-Propylbenzene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	3.3039		2.9			
o-Xylene	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	0.4609		6.0			
Tetrahydrofuran	FB	10.00	40.00	100.00	200.00	400.00	Ave	0.1355		9.5			
Isobutanol	FB	50.00	200.00	500.00	1000.00	2000.00	Ave	0.0162		9.5			
2-Chloro-1,3-butadiene	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.1935		5.3			
Ethanol	FB	50.00	200.00	500.00	1000.00	2000.00	Ave	0.0178		6.9			
4-Bromofluorobenzene	DCB	5.00	20.00	25.00	100.00	200.00	Ave	0.9379		6.3			
4-Isopropyltoluene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	2.7945		3.2			
n-Butyl acetate	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	0.2901		11.8			
Methacrylonitrile	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.4469		9.9			
n-Butylbenzene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	3.3672		10.4			
1,1-Dichloropropene	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.4776		6.2			
cis-1,2-Dichloroethene	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.2859		6.5			

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	Amount ug/L					Curve Evaluation						Min R^2 or COD
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5	Curve Type	Ave RRF	Min Ave RRF	%RSD	Max %RSD	R^2 or COD	
1,1,2,2-Tetrachloroethane	DCB	5.00	20.00	50.00	100.00	200.00	Ave	1.0638	0.3000	1.5			
1,2,4-Trimethylbenzene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	2.4805		3.1			
Toluene	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	1.0645		4.7	30.0		
Naphthalene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	2.2068		36.9			
1,3,5-Trimethylbenzene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	2.4694		2.1			
1,1-Dichloroacetone	CBZ	25.00	100.00	250.00	500.00	1000.00	Ave	0.2611		8.8			
1,3-Dichloropropane	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	0.5774		4.3			
Chloroform	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.5224		6.3	30.0		
4-Chlorotoluene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	2.1467		2.7			
Chlorodibromomethane	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	0.4000		7.8			
Dichlorodifluoromethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.1920		6.7			
1,1,2-Trichloroethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.2502		5.5			
Tert-amyl methyl ether	FB	5.00	20.00	50.00	100.00	200.00	Ave	1.0524		6.6			
Toluene-d8 (Surr)	CBZ	5.00	20.00	25.00	100.00	200.00	Ave	0.8889		4.2			
tert-Butylbenzene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	2.3884		3.2			

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	Amount ug/L					Curve Evaluation						Min R^2 or COD
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5	Curve Type	Ave RRF	Min Ave RRF	%RSD	Max %RSD	R^2 or COD	
1,1-Dichloro-1-fluoroethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.4422		5.0			
Chloromethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.3363	0.1000	6.3			
Methylene Chloride	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.2929		5.6			
Isopropyl acetate	FB	10.00	40.00	100.00	200.00	400.00	Ave	0.1342		14.4			
4-Ethyltoluene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	3.0136		2.7			
Methyl methacrylate	FB	10.00	40.00	100.00	200.00	400.00	Ave	0.1736		8.2			
1,1-Dichloroethene	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.2131		5.4	30.0		
Isopropylbenzene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	3.0830		2.5			
1,2-Dichloroethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.4746		4.6			
1,2-Dichloroethane-d4 (Surr)	FB	5.00	20.00	25.00	100.00	200.00	Ave	0.3932		4.9			
Acrylonitrile	FB	10.00	40.00	100.00	200.00	400.00	Ave	0.1834		6.9			
methyl isobutyl ketone	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	0.4602		7.6			
Tetrachloroethene	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	0.2100		6.3			
1,2-Dichloroethene, Total	FB	10.00	40.00	100.00	200.00	400.00	Ave	0.2306		5.4			
1,1,1-Trichloroethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.3840		6.7			

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	Amount ug/L					Curve Evaluation						Min R^2 or COD
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5	Curve Type	Ave RRF	Min Ave RRF	%RSD	Max %RSD	R^2 or COD	
2,2-Dichloropropane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.4711		4.4			
Benzyl chloride	DCB	5.00	20.00	50.00	100.00	200.00	Ave	0.4143		11.8			
Ethylene Dibromide	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	0.3366		7.3			
1,1,2-Trichloro-1,2,2-trifluoroethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.2475		7.0			
Methyl acetate	FB	5.00	20.00	50.00	100.00	200.00	Ave	1.8684		4.8			
Bromoform	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	0.2207	0.1000	13.4			
1,2-Dibromo-3-Chloropropane	DCB	5.00	20.00	50.00	100.00	200.00	Ave	0.2124		18.7			
Trichlorofluoromethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.3584		5.8			
Trichloroethene	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.3531		5.3			
Methylcyclohexane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.3498		4.3			
Bromobenzene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	0.7785		2.2			
Methyl Ethyl Ketone	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.2301		3.4			
1,2-Dichloropropane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.3960		5.4	30.0		
1,4-Dioxane	FB	50.00	200.00	500.00	1000.00	2000.00	Ave	0.0030		18.1			
1-Chlorobutane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.6854		4.9			

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	Amount ug/L					Curve Evaluation						Min R^2 or COD
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5	Curve Type	Ave RRF	Min Ave RRF	%RSD	Max %RSD	R^2 or COD	
n-Butanol	FB	50.00	200.00	500.00	1000.00	2000.00	Ave	0.0130		11.7			
Tert-butyl ethyl ether	FB	5.00	20.00	50.00	100.00	200.00	Ave	1.2132		5.8			
Methyl tert-butyl ether	FB	5.00	20.00	50.00	100.00	200.00	Ave	1.0379		5.6			
1,4-Dichlorobenzene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	1.3897		4.2			
1,1,1,2-Tetrachloroethane	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	0.3128		6.4			
Ethylbenzene	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	0.3850		8.3	30.0		
trans-1,2-Dichloroethene	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.2676		5.2			
Hexachlorobutadiene	DCB	5.00	20.00	50.00	100.00	200.00	Ave	0.2481		7.7			
Dichlorofluoromethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.6540		6.3			
1,1-Dichloroethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.6829	0.1000	5.2			
Ethyl ether	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.2255		3.4			
1-Bromopropane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.5150		4.2			
Acetonitrile	FB	50.00	200.00	500.00	1000.00	2000.00	Ave	0.0661		6.2			
Vinyl acetate	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.9114		6.3			
Ethyl methacrylate	CBZ	5.00	20.00	50.00	100.00	200.00	Ave	0.5666		7.4			

Volatile Initial Calibration Data

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1 Cal ID: 241

SDG No.: 220-2277

Instrument ID: MSL Column: RTX-VMS Heated Purge: (Y/N) N

Calibration Dates: 07/26/2007 13:32 07/26/2007 15:11

Analyte:	ISTD Ref	Amount ug/L					Curve Evaluation						Min R^2 or COD
		IC 220-8212/1	IC 220-8212/2	IC 220-8212/3	IC 220-8212/4	IC 220-8212/5	Curve Type	Ave RRF	Min Ave RRF	%RSD	Max %RSD	R^2 or COD	
Bromomethane	FB	5.00	20.00	50.00	100.00	200.00	Ave	0.1293		6.4			
Chloroacetonitrile	FB	50.00	200.00	500.00	1000.00	2000.00	Ave	0.0162		11.6			

Curve Types:
Ave = Average
Lin = Linear
Quad = Quadratic

STL-CT

Volatile Report SW-846 Method 8260B

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079145.b\L9152.D
Lab Smp Id: IC;5 Client Smp ID: IC;5
Inj Date : 26-JUL-2007 13:32 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : IC;5
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079145.b\L8260BNW.m
Meth Date : 26-Jul-2007 15:45 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 13:32 Cal File: L9152.D
Als bottle: 13 Calibration Sample, Level: 1
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	ON-COL (ug/L)	
* 1	Fluorobenzene	96	4.899	4.882	(1.000)	475766	25.0000		
2	Dichlorodifluoromethane	85	1.151	1.143	(0.235)	16212	5.00000		
3	Chloromethane	50	1.269	1.261	(0.259)	28831	5.00000		
4	Vinyl Chloride	62	1.298	1.301	(0.265)	31378	5.00000		
5	Bromomethane	94	1.485	1.468	(0.303)	12882	5.00000		
6	Chloroethane	64	1.554	1.517	(0.317)	20802	5.00000		
7	Trichlorofluoromethane	101	1.633	1.596	(0.333)	30814	5.00000		
8	Dichlorofluoromethane	67	1.652	1.625	(0.337)	57305	5.00000		(T)
9	Ethyl Ether	45	1.800	1.783	(0.367)	20403	5.00000		
10	Ethanol	45	1.859	1.842	(0.380)	15824	50.0000		
11	Freon 141	81	1.859	1.842	(0.380)	39048	5.00000		
12	Freon 123a	67	1.652	1.625	(0.337)	57305	5.00000		
13	Trichlorotrifluoroethane	101	1.948	1.930	(0.398)	21450	5.00000		
14	1,1-Dichloroethene	96	1.938	1.921	(0.396)	19116	5.00000		
15	Carbon Disulfide	76	1.977	1.960	(0.404)	92427	5.00000		
16	Iodomethane	142	2.036	2.019	(0.416)	13266	5.00000		
17	Acrolein	56	2.135	2.117	(0.436)	22421	25.0000		
18	2-Propanol	45	2.213	2.235	(0.452)	4423	5.00000		(M)
19	3-Chloro-1-Propene	41	2.223	2.216	(0.454)	49164	5.00000		
20	Methylene Chloride	84	2.302	2.285	(0.470)	25882	5.00000		
21	Acetone	43	2.321	2.314	(0.474)	17143	5.00000		
22	trans-1,2-Dichloroethene	96	2.420	2.412	(0.494)	24700	5.00000		
23	Methyl Acetate	43	2.410	2.393	(0.492)	167548	5.00000		

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	AMOUNTS ON-COL (ug/L)	
	24 Methyl tert-Butyl Ether	73	2.499	2.481	(0.510)	90955	5.00000		
	25 tert-Butyl alcohol	59	2.538	2.540	(0.518)	26664	25.0000		
	27 Isopropyl ether	45	2.784	2.767	(0.568)	108795	5.00000		
	28 tert-Butyl ethyl ether	59	3.118	3.091	(0.637)	106892	5.00000		
	29 2-Chloro-1,3-Butadiene	88	2.882	2.875	(0.588)	17761	5.00000		
	30 Acrylonitrile	53	2.922	2.914	(0.596)	30987	10.0000		
	31 1,1-Dichloroethane	63	2.902	2.885	(0.592)	60288	5.00000		
	32 Vinyl Acetate	43	3.109	3.091	(0.635)	79237	5.00000		
	33 cis-1,2-Dichloroethene	96	3.414	3.396	(0.697)	24420	5.00000		
	34 2,2-Dichloropropane	77	3.522	3.505	(0.719)	47220	5.00000		
	35 Bromochloromethane	128	3.630	3.613	(0.741)	18190	5.00000		
	36 1-Bromopropane	43	3.620	3.603	(0.739)	47182	5.00000		
	37 Cyclohexane	84	3.650	3.632	(0.745)	31245	5.00000		
	38 Chloroform	83	3.699	3.692	(0.755)	44969	5.00000		
	39 Ethyl Acetate	43	3.866	3.829	(0.789)	4639	10.0000		(M)
	40 Methyl Acrylate	55	3.856	3.839	(0.787)	33455	5.00000		
$	41 Dibromofluoromethane	111	3.915	3.908	(0.799)	27351	5.00000		
	42 Tetrahydrofuran	42	3.935	3.888	(0.803)	21716	10.0000		
	43 Carbon Tetrachloride	117	3.896	3.878	(0.795)	34409	5.00000		
	44 1,1,1-Trichloroethane	97	3.965	3.947	(0.809)	33468	5.00000		
	45 2-Butanone	43	4.073	4.046	(0.831)	21580	5.00000		
	46 1,1-Dichloropropene	75	4.112	4.105	(0.839)	42178	5.00000		
	47 tert-Amyl methyl ether	73	4.584	4.548	(0.936)	90528	5.00000		
	48 tert-Butyl formate	57	3.109	3.091	(0.635)	31950	5.00000		
	49 1-Chlorobutane	56	4.171	4.154	(0.851)	62421	5.00000		
	50 Heptane	43	4.398	4.380	(0.898)	30506	5.00000		
	51 Propionitrile	54	4.407	4.390	(0.900)	54782	50.0000		
	52 Benzene	78	4.407	4.400	(0.900)	112138	5.00000		
	53 2-Methyl-2-Propenenitrile	41	4.437	4.420	(0.906)	43819	5.00000		
$	55 1,2-Dichloroethane-d4	65	4.565	4.548	(0.932)	35255	5.00000		
	56 1,2-Dichloroethane	62	4.643	4.626	(0.948)	42349	5.00000		
	59 Methyl Cyclohexane	83	5.096	5.079	(1.040)	32018	5.00000		
	60 Trichloroethene	130	5.096	5.089	(1.040)	31768	5.00000		
	61 Isopropyl Acetate	43	4.693	4.695	(0.958)	27503	10.0000		
	62 N-Butanol	56	5.490	5.482	(1.120)	10799	50.0000		
	63 Dibromomethane	93	5.529	5.522	(1.129)	17268	5.00000		
	64 1,2-Dichloropropane	63	5.627	5.620	(1.149)	34797	5.00000		(T)
	65 Bromodichloromethane	83	5.706	5.709	(1.165)	30949	5.00000		
	66 Methyl Methacrylate	69	5.893	5.886	(1.203)	28827	10.0000		
	67 1,4-Dioxane	58	5.942	5.935	(1.213)	2233	50.0000		(M)
	68 N-Propyl Acetate	43	6.296	6.289	(1.285)	4252	10.0000		(M)
	69 2-Chloroethylvinylether	63	6.296	6.289	(1.285)	7215	5.00000		(M)
	70 cis-1,3-Dichloropropene	75	6.346	6.338	(1.295)	49856	5.00000		
	71 Chloroacetonitrile	48	6.690	6.692	(1.365)	13401	50.0000		
	72 2-Nitropropane	41	6.769	6.761	(1.382)	20683	10.0000		(T)
	73 trans-1,3-Dichloropropene	75	6.975	6.968	(1.424)	42622	5.00000		
	74 1,1,2-Trichloroethane	97	7.123	7.115	(1.454)	22152	5.00000		
*	75 Chlorobenzene-d5	117	7.959	7.952	(1.000)	449491	25.0000		
	76 Toluene	91	6.582	6.574	(0.827)	90482	5.00000		
$	77 Toluene-d8	98	6.533	6.525	(0.821)	78005	5.00000		
	78 1,1-Dichloro-2-propanone	43	6.798	6.791	(0.854)	103609	25.0000		
	79 4-Methyl-2-Pentanone	43	6.936	6.929	(0.871)	36222	5.00000		
	80 Tetrachloroethene	164	6.956	6.948	(0.874)	17064	5.00000		
	81 Ethyl Methacrylate	69	7.143	7.135	(0.897)	44954	5.00000		

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	ON-COL (ug/L)	
82	Dibromochloromethane	129	7.290	7.283	(0.916)	32507	5.00000		
83	1,3-Dichloropropane	76	7.359	7.352	(0.925)	49474	5.00000		
84	1,2-Dibromoethane	107	7.487	7.480	(0.941)	27164	5.00000		
85	n-Butyl Acetate	56	7.644	7.637	(0.960)	21332	5.00000		
86	2-Hexanone	43	7.713	7.696	(0.969)	27248	5.00000		
87	1-Chlorohexane	91	7.969	7.962	(1.001)	28680	5.00000		
88	Chlorobenzene	112	7.969	7.962	(1.001)	70637	5.00000		
89	1,1,1,2-Tetrachloroethane	131	8.038	8.031	(1.010)	25594	5.00000		(M)
90	Ethylbenzene	106	8.008	8.001	(1.006)	30354	5.00000		
91	Xylene (total)mp	106	8.136	8.129	(1.022)	82928	10.0000		
92	Xylene (total)o	106	8.510	8.513	(1.069)	38171	5.00000		
93	Styrene	104	8.559	8.552	(1.075)	62763	5.00000		
94	Bromoform	173	8.579	8.581	(1.078)	15753	5.00000		
* 95	1,4-Dichlorobenzene-d4	152	10.016	10.008	(1.000)	150172	25.0000		
96	Isopropylbenzene	105	8.796	8.788	(0.878)	95462	5.00000		
97	Bromobenzene	156	9.120	9.123	(0.911)	23927	5.00000		
98	1,1,2,2-Tetrachloroethane	83	9.219	9.221	(0.920)	31243	5.00000		
99	4-Ethyltoluene	105	9.258	9.251	(0.924)	90350	5.00000		
100	1,2,3-Trichloropropane	110	9.327	9.319	(0.931)	9898	5.00000		
101	trans-1,4-Dichloro-2-Butene	53	9.366	9.359	(0.935)	16977	10.0000		
102	n-Propylbenzene	91	9.160	9.152	(0.915)	98837	5.00000		
103	2-Chlorotoluene	91	9.287	9.280	(0.927)	62778	5.00000		
104	4-Chlorotoluene	91	9.425	9.428	(0.941)	64301	5.00000		
105	1,3,5-Trimethylbenzene	105	9.327	9.329	(0.931)	74886	5.00000		
106	tert-Butylbenzene	119	9.602	9.605	(0.959)	72043	5.00000		
107	1,2,4-Trimethylbenzene	105	9.671	9.664	(0.966)	72192	5.00000		
108	sec-Butylbenzene	105	9.760	9.752	(0.974)	78630	5.00000		
109	4-Isopropyltoluene	119	9.888	9.880	(0.987)	81506	5.00000		
110	1,3-Dichlorobenzene	146	9.947	9.939	(0.993)	39658	5.00000		
111	1,4-Dichlorobenzene	146	10.025	10.018	(1.001)	39307	5.00000		
112	1,2-Dichlorobenzene	146	10.389	10.382	(1.037)	39430	5.00000		
113	Benzyl Chloride	126	10.232	10.234	(1.022)	10375	5.00000		
114	1,4-Diethylbenzene	119	10.203	10.205	(2.082)	41670	5.00000		
115	n-Butylbenzene	91	10.252	10.244	(1.024)	88645	5.00000		
118	1,2,4,5-Tetramethylbenzene	119	10.911	10.904	(2.227)	58670	5.00000		
119	1,2-Dibromo-3-chloropropane	75	11.078	11.081	(1.106)	4400	5.00000		
120	Nitrobenzene	77	11.570	11.563	(1.155)	3903	50.0000		
121	1,2,4-Trichlorobenzene	180	11.688	11.681	(1.167)	13729	5.00000		
122	Hexachlorobutadiene	225	11.669	11.671	(1.165)	7385	5.00000		
123	Naphthalene	128	11.964	11.956	(1.194)	31487	5.00000		
124	1,2,3-Trichlorobenzene	180	12.131	12.133	(1.211)	11761	5.00000		
$ 125	Bromofluorobenzene	95	9.032	9.034	(0.902)	31017	5.00000		
M 126	1,2-Dichloroethene (total)	100				49120	10.0000		
M 127	Xylene (total)	100				121099	15.0000		

QC Flag Legend

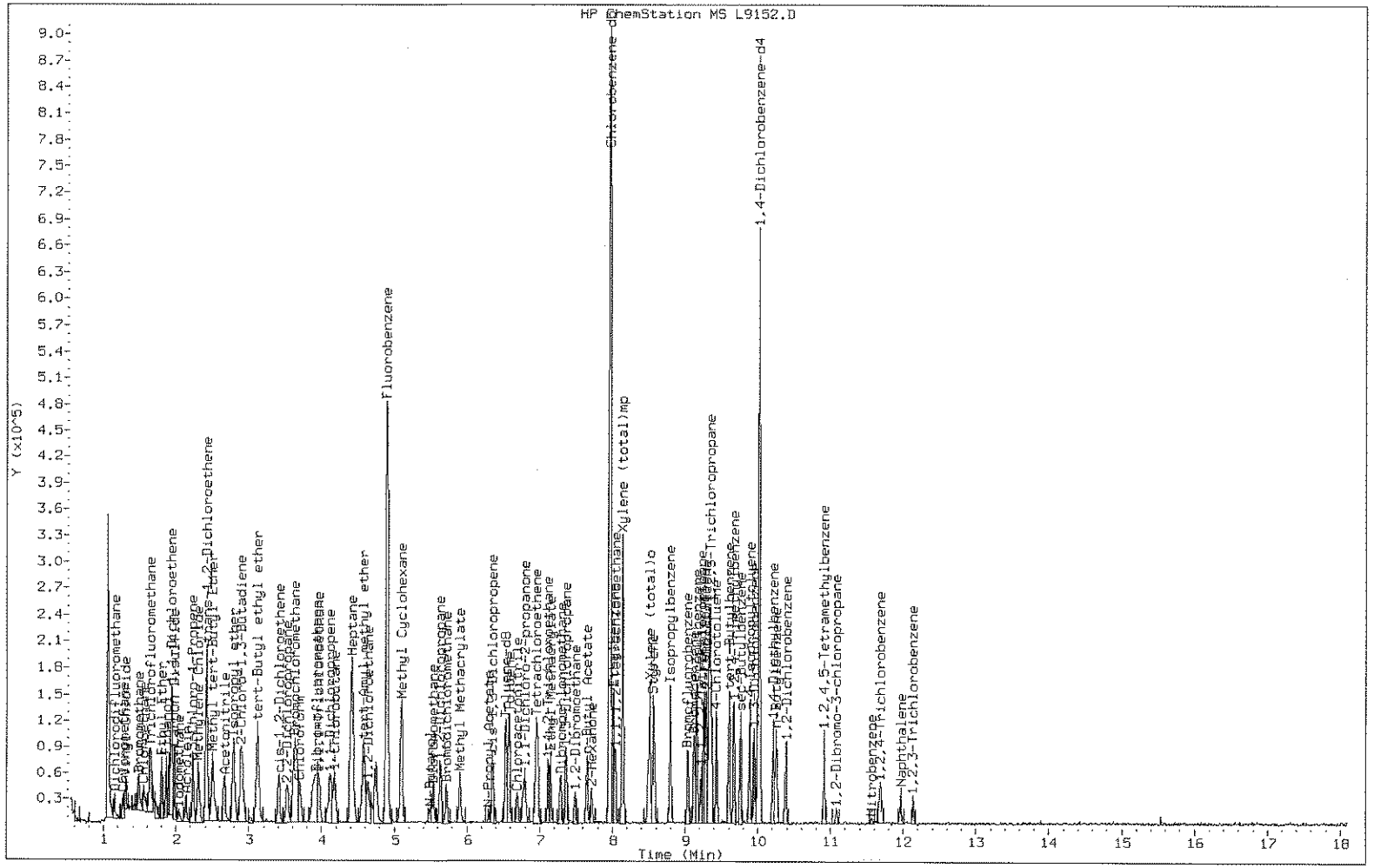
T - Target compound detected outside RT window.
M - Compound response manually integrated.

Data File: L9152.D

Date: 26-JUL-2007 13:32

Client ID: IC;5 Instrument: msl.i

Sample Info: IC;5 Operator: D. HUMBERT



08/09/2007

STL-CT

Volatile Report SW-846 Method 8260B

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079145.b\L9153.D
Lab Smp Id: IC;20 Client Smp ID: IC;20
Inj Date : 26-JUL-2007 13:57 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : IC;20
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079145.b\L8260BNW.m
Meth Date : 26-Jul-2007 15:45 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 13:57 Cal File: L9153.D
Als bottle: 14 Calibration Sample, Level: 2
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	AMOUNTS ON-COL (ug/L)
* 1	Fluorobenzene	96	4.890	4.882	(1.000)	472080	25.0000	
2	Dichlorodifluoromethane	85	1.142	1.143	(0.234)	72110	20.0000	22
3	Chloromethane	50	1.270	1.261	(0.260)	125228	20.0000	22
4	Vinyl Chloride	62	1.299	1.301	(0.266)	138805	20.0000	22
5	Bromomethane	94	1.476	1.468	(0.302)	45420	20.0000	18(M)
6	Chloroethane	64	1.545	1.517	(0.316)	91681	20.0000	22
7	Trichlorofluoromethane	101	1.624	1.596	(0.332)	135020	20.0000	22
8	Dichlorofluoromethane	67	1.644	1.625	(0.336)	234030	20.0000	20
9	Ethyl Ether	45	1.791	1.783	(0.366)	85126	20.0000	21
10	Ethanol	45	1.850	1.842	(0.378)	62344	200.000	200
11	Freon 141	81	1.850	1.842	(0.378)	161955	20.0000	21
12	Freon 123a	67	1.644	1.625	(0.336)	234030	20.0000	20
13	Trichlorotrifluoroethane	101	1.939	1.930	(0.396)	88332	20.0000	21
14	1,1-Dichloroethene	96	1.929	1.921	(0.394)	76133	20.0000	20
15	Carbon Disulfide	76	1.968	1.960	(0.403)	366490	20.0000	20
16	Iodomethane	142	2.027	2.019	(0.415)	79483	20.0000	30
17	Acrolein	56	2.126	2.117	(0.435)	101390	100.000	110
18	2-Propanol	45	2.214	2.235	(0.453)	20365	20.0000	23(H)
19	3-Chloro-1-Propene	41	2.224	2.216	(0.455)	199100	20.0000	20
20	Methylene Chloride	84	2.293	2.285	(0.469)	105312	20.0000	20
21	Acetone	43	2.322	2.314	(0.475)	56640	20.0000	17
22	trans-1,2-Dichloroethene	96	2.411	2.412	(0.493)	94065	20.0000	19
23	Methyl Acetate	43	2.401	2.393	(0.491)	677474	20.0000	20

			QUANT SIG					AMOUNTS	
		Compounds	MASS	RT	EXP RT	REL RT	RESPONSE	CAL-AMT (ug/L)	ON-COL (ug/L)
	24	Methyl tert-Butyl Ether	73	2.490	2.481	(0.509)	377858	20.0000	21
	25	tert-Butyl alcohol	59	2.529	2.540	(0.517)	100091	100.000	94
	26	Acetonitrile	41	2.657	2.658	(0.543)	236520	200.000	200
	27	Isopropyl ether	45	2.775	2.767	(0.567)	439684	20.0000	20
	28	tert-Butyl ethyl ether	59	3.110	3.091	(0.636)	434663	20.0000	20
	29	2-Chloro-1,3-Butadiene	88	2.873	2.875	(0.588)	67831	20.0000	19
	30	Acrylonitrile	53	2.913	2.914	(0.596)	146120	40.0000	48
	31	1,1-Dichloroethane	63	2.893	2.885	(0.592)	250389	20.0000	21
	32	Vinyl Acetate	43	3.100	3.091	(0.634)	328518	20.0000	21
	33	cis-1,2-Dichloroethene	96	3.405	3.396	(0.696)	105969	20.0000	22
	34	2,2-Dichloropropane	77	3.513	3.505	(0.718)	165808	20.0000	18
	35	Bromochloromethane	128	3.611	3.613	(0.738)	74181	20.0000	20
	36	1-Bromopropane	43	3.611	3.603	(0.738)	184942	20.0000	20
	37	Cyclohexane	84	3.641	3.632	(0.745)	126792	20.0000	20
	38	Chloroform	83	3.690	3.692	(0.755)	191184	20.0000	21
	39	Ethyl Acetate	43	3.847	3.829	(0.787)	21359	40.0000	46
	40	Methyl Acrylate	55	3.847	3.839	(0.787)	140894	20.0000	21
$	41	Dibromofluoromethane	111	3.907	3.908	(0.799)	119842	20.0000	22
	42	Tetrahydrofuran	42	3.916	3.888	(0.801)	102061	40.0000	47
	43	Carbon Tetrachloride	117	3.887	3.878	(0.795)	142637	20.0000	21
	44	1,1,1-Trichloroethane	97	3.956	3.947	(0.809)	137208	20.0000	21
	45	2-Butanone	43	4.054	4.046	(0.829)	82547	20.0000	19
	46	1,1-Dichloropropene	75	4.103	4.105	(0.839)	169560	20.0000	20
	47	tert-Amyl methyl ether	73	4.566	4.548	(0.934)	385429	20.0000	21
	48	tert-Butyl formate	57	3.100	3.091	(0.634)	124826	20.0000	20
	49	1-Chlorobutane	56	4.162	4.154	(0.851)	243396	20.0000	20
	50	Heptane	43	4.389	4.380	(0.897)	108120	20.0000	18
	51	Propionitrile	54	4.389	4.390	(0.897)	224269	200.000	210
	52	Benzene	78	4.398	4.400	(0.899)	423764	20.0000	19
	53	2-Methyl-2-Propenenitrile	41	4.418	4.420	(0.903)	169132	20.0000	19
	54	Isobutyl alcohol	42	4.674	4.695	(0.956)	55820	200.000	210
$	55	1,2-Dichloroethane-d4	65	4.546	4.548	(0.930)	143243	20.0000	20
	56	1,2-Dichloroethane	62	4.625	4.626	(0.946)	173568	20.0000	21
	59	Methyl Cyclohexane	83	5.077	5.079	(1.038)	126470	20.0000	20
	60	Trichloroethene	130	5.087	5.089	(1.040)	125730	20.0000	20
	61	Isopropyl Acetate	43	4.674	4.695	(0.956)	75602	40.0000	28
	62	N-Butanol	56	5.471	5.482	(1.119)	44001	200.000	200
	63	Dibromomethane	93	5.520	5.522	(1.129)	68308	20.0000	20
	64	1,2-Dichloropropane	63	5.619	5.620	(1.149)	145305	20.0000	21
	65	Bromodichloromethane	83	5.707	5.709	(1.167)	133083	20.0000	22
	66	Methyl Methacrylate	69	5.884	5.886	(1.203)	127836	40.0000	45
	67	1,4-Dioxane	58	5.933	5.935	(1.213)	9467	200.000	210
	68	N-Propyl Acetate	43	6.288	6.289	(1.286)	24024	40.0000	57
	69	2-Chloroethylvinylether	63	6.288	6.289	(1.286)	34616	20.0000	24
	70	cis-1,3-Dichloropropene	75	6.337	6.338	(1.296)	199651	20.0000	20
	71	Chloroacetonitrile	48	6.681	6.692	(1.366)	54725	200.000	200
	72	2-Nitropropane	41	6.760	6.761	(1.382)	78256	40.0000	38
	73	trans-1,3-Dichloropropene	75	6.966	6.968	(1.424)	182966	20.0000	22
	74	1,1,2-Trichloroethane	97	7.114	7.115	(1.455)	90934	20.0000	21
*	75	Chlorobenzene-d5	117	7.950	7.952	(1.000)	449699	25.0000	
	76	Toluene	91	6.573	6.574	(0.827)	368460	20.0000	20
$	77	Toluene-d8	98	6.524	6.525	(0.821)	308915	20.0000	20
	78	1,1-Dichloro-2-propanone	43	6.789	6.791	(0.854)	441352	100.000	110
	79	4-Methyl-2-Pentanone	43	6.927	6.929	(0.871)	163108	20.0000	22

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	ON-COL (ug/L)
80	Tetrachloroethene	164	6.947	6.948	(0.874)	73241	20.0000	21
81	Ethyl Methacrylate	69	7.134	7.135	(0.897)	201679	20.0000	22
82	Dibromochloromethane	129	7.281	7.283	(0.916)	134793	20.0000	21
83	1,3-Dichloropropane	76	7.360	7.352	(0.926)	200701	20.0000	20
84	1,2-Dibromoethane	107	7.478	7.480	(0.941)	115370	20.0000	21
85	n-Butyl Acetate	56	7.635	7.637	(0.960)	99825	20.0000	23
86	2-Hexanone	43	7.704	7.696	(0.969)	119181	20.0000	22
87	1-Chlorohexane	91	7.960	7.962	(1.001)	148459	20.0000	26
88	Chlorobenzene	112	7.970	7.962	(1.002)	302937	20.0000	21
89	1,1,1,2-Tetrachloroethane	131	8.029	8.031	(1.010)	109035	20.0000	21
90	Ethylbenzene	106	8.000	8.001	(1.006)	133140	20.0000	22
91	Xylene (total)mp	106	8.127	8.129	(1.022)	325887	40.0000	39
92	Xylene (total)o	106	8.511	8.513	(1.071)	159229	20.0000	21
93	Styrene	104	8.560	8.552	(1.077)	264511	20.0000	21
94	Bromoform	173	8.580	8.581	(1.079)	74614	20.0000	24
* 95	1,4-Dichlorobenzene-d4	152	10.007	10.008	(1.000)	162525	25.0000	
96	Isopropylbenzene	105	8.787	8.788	(0.878)	387937	20.0000	19
97	Bromobenzene	156	9.121	9.123	(0.912)	97563	20.0000	19
98	1,1,2,2-Tetrachloroethane	83	9.210	9.221	(0.920)	137004	20.0000	20
99	4-Ethyltoluene	105	9.249	9.251	(0.924)	375738	20.0000	19
100	1,2,3-Trichloropropane	110	9.318	9.319	(0.931)	44647	20.0000	21
101	trans-1,4-Dichloro-2-Butene	53	9.357	9.359	(0.935)	88666	40.0000	48
102	n-Propylbenzene	91	9.151	9.152	(0.914)	410870	20.0000	19
103	2-Chlorotoluene	91	9.279	9.280	(0.927)	308780	20.0000	23
104	4-Chlorotoluene	91	9.426	9.428	(0.942)	267004	20.0000	19
105	1,3,5-Trimethylbenzene	105	9.328	9.329	(0.932)	309296	20.0000	19
106	tert-Butylbenzene	119	9.603	9.605	(0.960)	294479	20.0000	19
107	1,2,4-Trimethylbenzene	105	9.662	9.664	(0.966)	311887	20.0000	20
108	sec-Butylbenzene	105	9.751	9.752	(0.974)	327119	20.0000	19
109	4-Isopropyltoluene	119	9.879	9.880	(0.987)	353954	20.0000	20
110	1,3-Dichlorobenzene	146	9.938	9.939	(0.993)	171768	20.0000	20
111	1,4-Dichlorobenzene	146	10.017	10.018	(1.001)	175534	20.0000	21
112	1,2-Dichlorobenzene	146	10.381	10.382	(1.037)	168529	20.0000	20
113	Benzyl Chloride	126	10.233	10.234	(1.023)	50283	20.0000	22
114	1,4-Diethylbenzene	119	10.203	10.205	(2.086)	185331	20.0000	22
115	n-Butylbenzene	91	10.243	10.244	(1.024)	395205	20.0000	20
118	1,2,4,5-Tetramethylbenzene	119	10.902	10.904	(2.229)	278703	20.0000	24
119	1,2-Dibromo-3-chloropropane	75	11.079	11.081	(1.107)	27578	20.0000	29
120	Nitrobenzene	77	11.561	11.563	(1.155)	34805	200.000	410
121	1,2,4-Trichlorobenzene	180	11.679	11.681	(1.167)	81473	20.0000	27
122	Hexachlorobutadiene	225	11.669	11.671	(1.166)	28509	20.0000	18
123	Naphthalene	128	11.955	11.956	(1.195)	229159	20.0000	34
124	1,2,3-Trichlorobenzene	180	12.132	12.133	(1.212)	76112	20.0000	30
$ 125	Bromofluorobenzene	95	9.033	9.034	(0.903)	115064	20.0000	17
M 126	1,2-Dichloroethene (total)	100				200034	40.0000	
M 127	Xylene (total)	100				485116	60.0000	

QC Flag Legend

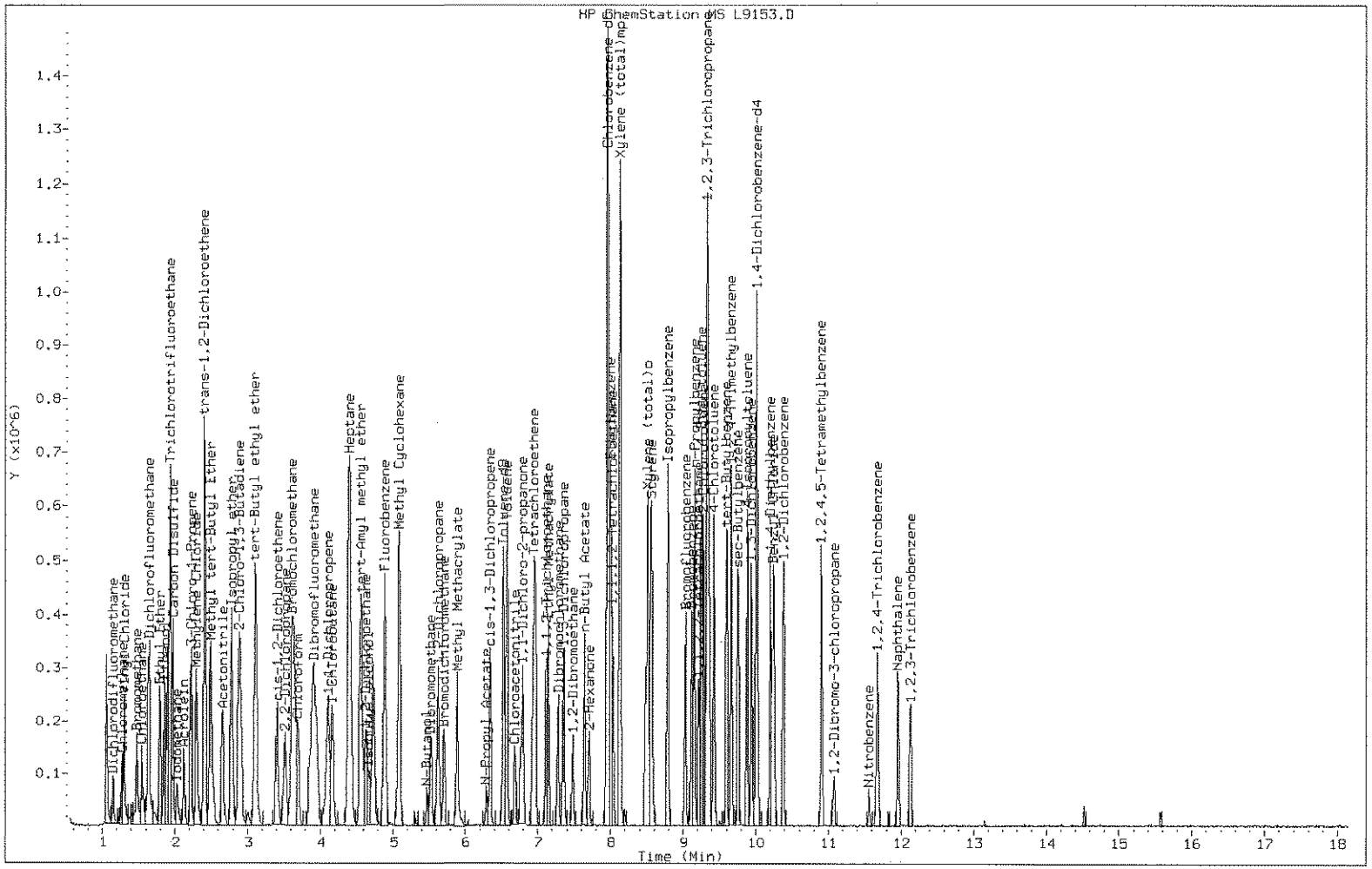
M - Compound response manually integrated.
H - Operator selected an alternate compound hit.

Data File: L9153.D

Date: 26-JUL-2007 13:57

Client ID: IC;20 Instrument: msl.i

Sample Info: IC;20 Operator: D. HUMBERT



STL-CT

Volatile Report SW-846 Method 8260B

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079145.b\L9154.D
Lab Smp Id: IC;50 Client Smp ID: IC;50
Inj Date : 26-JUL-2007 14:22 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : IC;50
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079145.b\L8260BNW.m
Meth Date : 26-Jul-2007 15:45 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 14:22 Cal File: L9154.D
Als bottle: 15 Calibration Sample, Level: 3
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	AMOUNTS ON-COL (ug/L)
* 1	Fluorobenzene	96	4.885	4.882	(1.000)	463133	25.0000	
2	Dichlorodifluoromethane	85	1.146	1.143	(0.235)	181338	50.0000	54
3	Chloromethane	50	1.264	1.261	(0.259)	313165	50.0000	53
4	Vinyl Chloride	62	1.304	1.301	(0.267)	348675	50.0000	54
5	Bromomethane	94	1.471	1.468	(0.301)	113131	50.0000	48(M)
6	Chloroethane	64	1.540	1.517	(0.315)	219075	50.0000	51
7	Trichlorofluoromethane	101	1.619	1.596	(0.331)	337670	50.0000	54
8	Dichlorofluoromethane	67	1.638	1.625	(0.335)	617375	50.0000	54
9	Ethyl Ether	45	1.796	1.783	(0.368)	218611	50.0000	54
10	Ethanol	45	1.855	1.842	(0.380)	164271	500.000	540
11	Freon 141	81	1.855	1.842	(0.380)	417999	50.0000	54
12	Freon 123a	67	1.638	1.625	(0.335)	617375	50.0000	54
13	Trichlorotrifluoroethane	101	1.933	1.930	(0.396)	233148	50.0000	55
14	1,1-Dichloroethene	96	1.924	1.921	(0.394)	198480	50.0000	53
15	Carbon Disulfide	76	1.963	1.960	(0.402)	969738	50.0000	54
16	Iodomethane	142	2.032	2.019	(0.416)	309220	50.0000	95
17	Acrolein	56	2.120	2.117	(0.434)	283448	250.000	300
18	2-Propanol	45	2.209	2.235	(0.452)	52218	50.0000	56
19	3-Chloro-1-Propene	41	2.219	2.216	(0.454)	521673	50.0000	54
20	Methylene Chloride	84	2.288	2.285	(0.468)	279838	50.0000	55
21	Acetone	43	2.317	2.314	(0.474)	151536	50.0000	50
22	trans-1,2-Dichloroethene	96	2.416	2.412	(0.495)	247669	50.0000	52
23	Methyl Acetate	43	2.396	2.393	(0.490)	1745753	50.0000	53

		Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	AMOUNTS ON-COL (ug/L)
	24	Methyl tert-Butyl Ether	73	2.484	2.481	(0.509)	978615	50.0000	54
	25	tert-Butyl alcohol	59	2.524	2.540	(0.517)	267292	250.000	260
	26	Acetonitrile	41	2.652	2.658	(0.543)	626539	500.000	550
	27	Isopropyl ether	45	2.770	2.767	(0.567)	1124714	50.0000	53
	28	tert-Butyl ethyl ether	59	3.104	3.091	(0.635)	1160436	50.0000	55
	29	2-Chloro-1,3-Butadiene	88	2.878	2.875	(0.589)	183557	50.0000	54
	30	Acrylonitrile	53	2.907	2.914	(0.595)	332799	100.000	100
	31	1,1-Dichloroethane	63	2.888	2.885	(0.591)	640865	50.0000	53
	32	Vinyl Acetate	43	3.094	3.091	(0.633)	863863	50.0000	55
	33	cis-1,2-Dichloroethene	96	3.399	3.396	(0.696)	269414	50.0000	54
	34	2,2-Dichloropropane	77	3.518	3.505	(0.720)	439361	50.0000	51
	35	Bromochloromethane	128	3.616	3.613	(0.740)	195300	50.0000	54
	36	1-Bromopropane	43	3.606	3.603	(0.738)	482802	50.0000	53
	37	Cyclohexane	84	3.645	3.632	(0.746)	347286	50.0000	56
	38	Chloroform	83	3.695	3.692	(0.756)	495681	50.0000	55
	39	Ethyl Acetate	43	3.832	3.829	(0.785)	53278	100.000	110
	40	Methyl Acrylate	55	3.842	3.839	(0.787)	386392	50.0000	58
$	41	Dibromofluoromethane	111	3.911	3.908	(0.801)	151938	25.0000	27
	42	Tetrahydrofuran	42	3.901	3.888	(0.799)	254358	100.000	110
	43	Carbon Tetrachloride	117	3.882	3.878	(0.795)	444509	50.0000	65
	44	1,1,1-Trichloroethane	97	3.950	3.947	(0.809)	362380	50.0000	55
	45	2-Butanone	43	4.049	4.046	(0.829)	221560	50.0000	54
	46	1,1-Dichloropropene	75	4.108	4.105	(0.841)	452006	50.0000	55
	47	tert-Amyl methyl ether	73	4.551	4.548	(0.932)	986156	50.0000	54
	48	tert-Butyl formate	57	3.104	3.091	(0.635)	332603	50.0000	54
	49	1-Chlorobutane	56	4.157	4.154	(0.851)	643857	50.0000	53
	50	Heptane	43	4.383	4.380	(0.897)	288619	50.0000	51
	51	Propionitrile	54	4.383	4.390	(0.897)	610664	500.000	560
	52	Benzene	78	4.403	4.400	(0.901)	1117566	50.0000	52
	53	2-Methyl-2-Propenenitrile	41	4.423	4.420	(0.905)	441023	50.0000	52
	54	Isobutyl alcohol	42	4.669	4.695	(0.956)	153541	500.000	570
$	55	1,2-Dichloroethane-d4	65	4.551	4.548	(0.932)	190180	25.0000	27
	56	1,2-Dichloroethane	62	4.629	4.626	(0.948)	445956	50.0000	53
	59	Methyl Cyclohexane	83	5.082	5.079	(1.040)	322681	50.0000	52
	60	Trichloroethene	130	5.082	5.089	(1.040)	331611	50.0000	54
	61	Isopropyl Acetate	43	4.669	4.695	(0.956)	252997	100.000	110
	62	N-Butanol	56	5.466	5.482	(1.119)	120808	500.000	570
	63	Dibromomethane	93	5.525	5.522	(1.131)	181340	50.0000	54
	64	1,2-Dichloropropane	63	5.623	5.620	(1.151)	371118	50.0000	53
	65	Bromodichloromethane	83	5.702	5.709	(1.167)	352418	50.0000	56
	66	Methyl Methacrylate	69	5.879	5.886	(1.203)	325785	100.000	110
	67	1,4-Dioxane	58	5.918	5.935	(1.211)	29249	500.000	650
	68	N-Propyl Acetate	43	6.292	6.289	(1.288)	60440	100.000	120
	69	2-Chloroethylvinylether	63	6.292	6.289	(1.288)	91478	50.0000	59
	70	cis-1,3-Dichloropropene	75	6.341	6.338	(1.298)	528803	50.0000	54
	71	Chloroacetonitrile	48	6.676	6.692	(1.367)	151888	500.000	570
	72	2-Nitropropane	41	6.764	6.761	(1.385)	204987	100.000	100
	73	trans-1,3-Dichloropropene	75	6.961	6.968	(1.425)	492914	50.0000	57
	74	1,1,2-Trichloroethane	97	7.109	7.115	(1.455)	231500	50.0000	53
*	75	Chlorobenzene-d5	117	7.945	7.952	(1.000)	456972	25.0000	
	76	Toluene	91	6.568	6.574	(0.827)	971370	50.0000	52
$	77	Toluene-d8	98	6.518	6.525	(0.820)	394689	25.0000	25
	78	1,1-Dichloro-2-propanone	43	6.784	6.791	(0.854)	1211371	250.000	280
	79	4-Methyl-2-Pentanone	43	6.932	6.929	(0.872)	436285	50.0000	56

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	ON-COL (ug/L)
80	Tetrachloroethene	164	6.941	6.948	(0.874)	196791	50.0000	55
81	Ethyl Methacrylate	69	7.138	7.135	(0.898)	518290	50.0000	53
82	Dibromochloromethane	129	7.276	7.283	(0.916)	371171	50.0000	55
83	1,3-Dichloropropane	76	7.355	7.352	(0.926)	525409	50.0000	52
84	1,2-Dibromoethane	107	7.483	7.480	(0.942)	315483	50.0000	55
85	n-Butyl Acetate	56	7.640	7.637	(0.962)	272791	50.0000	58
86	2-Hexanone	43	7.699	7.696	(0.969)	326541	50.0000	56
87	1-Chlorohexane	91	7.955	7.962	(1.001)	339575	50.0000	51
88	Chlorobenzene	112	7.965	7.962	(1.002)	799284	50.0000	54
89	1,1,1,2-Tetrachloroethane	131	8.024	8.031	(1.010)	287333	50.0000	53
90	Ethylbenzene	106	7.994	8.001	(1.006)	356676	50.0000	55
91	Xylene (total)mp	106	8.132	8.129	(1.024)	882967	100.000	100
92	Xylene (total)o	106	8.506	8.513	(1.071)	421892	50.0000	53
93	Styrene	104	8.555	8.552	(1.077)	706604	50.0000	54
94	Bromoform	173	8.575	8.581	(1.079)	212127	50.0000	61
* 95	1,4-Dichlorobenzene-d4	152	10.001	10.008	(1.000)	167085	25.0000	
96	Isopropylbenzene	105	8.791	8.788	(0.879)	1015372	50.0000	49
97	Bromobenzene	156	9.116	9.123	(0.911)	262763	50.0000	51
98	1,1,2,2-Tetrachloroethane	83	9.214	9.221	(0.921)	360375	50.0000	52
99	4-Ethyltoluene	105	9.254	9.251	(0.925)	1008337	50.0000	51
100	1,2,3-Trichloropropane	110	9.323	9.319	(0.932)	113672	50.0000	50
101	trans-1,4-Dichloro-2-Butene	53	9.362	9.359	(0.936)	243030	100.000	120
102	n-Propylbenzene	91	9.155	9.152	(0.915)	1097768	50.0000	51
103	2-Chlorotoluene	91	9.283	9.280	(0.928)	709538	50.0000	48
104	4-Chlorotoluene	91	9.421	9.428	(0.942)	729131	50.0000	52
105	1,3,5-Trimethylbenzene	105	9.323	9.329	(0.932)	825460	50.0000	51
106	tert-Butylbenzene	119	9.598	9.605	(0.960)	806648	50.0000	52
107	1,2,4-Trimethylbenzene	105	9.667	9.664	(0.967)	848818	50.0000	53
108	sec-Butylbenzene	105	9.755	9.752	(0.975)	868721	50.0000	51
109	4-Isopropyltoluene	119	9.883	9.880	(0.988)	931551	50.0000	51
110	1,3-Dichlorobenzene	146	9.942	9.939	(0.994)	464940	50.0000	53
111	1,4-Dichlorobenzene	146	10.021	10.018	(1.002)	472405	50.0000	53
112	1,2-Dichlorobenzene	146	10.385	10.382	(1.038)	454247	50.0000	52
113	Benzyl Chloride	126	10.228	10.234	(1.023)	141280	50.0000	58
114	1,4-Diethylbenzene	119	10.198	10.205	(2.087)	531372	50.0000	62
115	n-Butylbenzene	91	10.247	10.244	(1.025)	1219361	50.0000	61
118	1,2,4,5-Tetramethylbenzene	119	10.907	10.904	(2.232)	777768	50.0000	62
119	1,2-Dibromo-3-chloropropane	75	11.074	11.081	(1.107)	72717	50.0000	61
120	Nitrobenzene	77	11.566	11.563	(1.156)	142798	500.000	1100
121	1,2,4-Trichlorobenzene	180	11.684	11.681	(1.168)	260899	50.0000	72
122	Hexachlorobutadiene	225	11.664	11.671	(1.166)	81517	50.0000	52
123	Naphthalene	128	11.959	11.956	(1.196)	787295	50.0000	84
124	1,2,3-Trichlorobenzene	180	12.127	12.133	(1.212)	235406	50.0000	72
$ 125	Bromofluorobenzene	95	9.027	9.034	(0.903)	157780	25.0000	25
M 126	1,2-Dichloroethene (total)	100				517083	100.000	
M 127	Xylene (total)	100				1304859	150.000	

QC Flag Legend

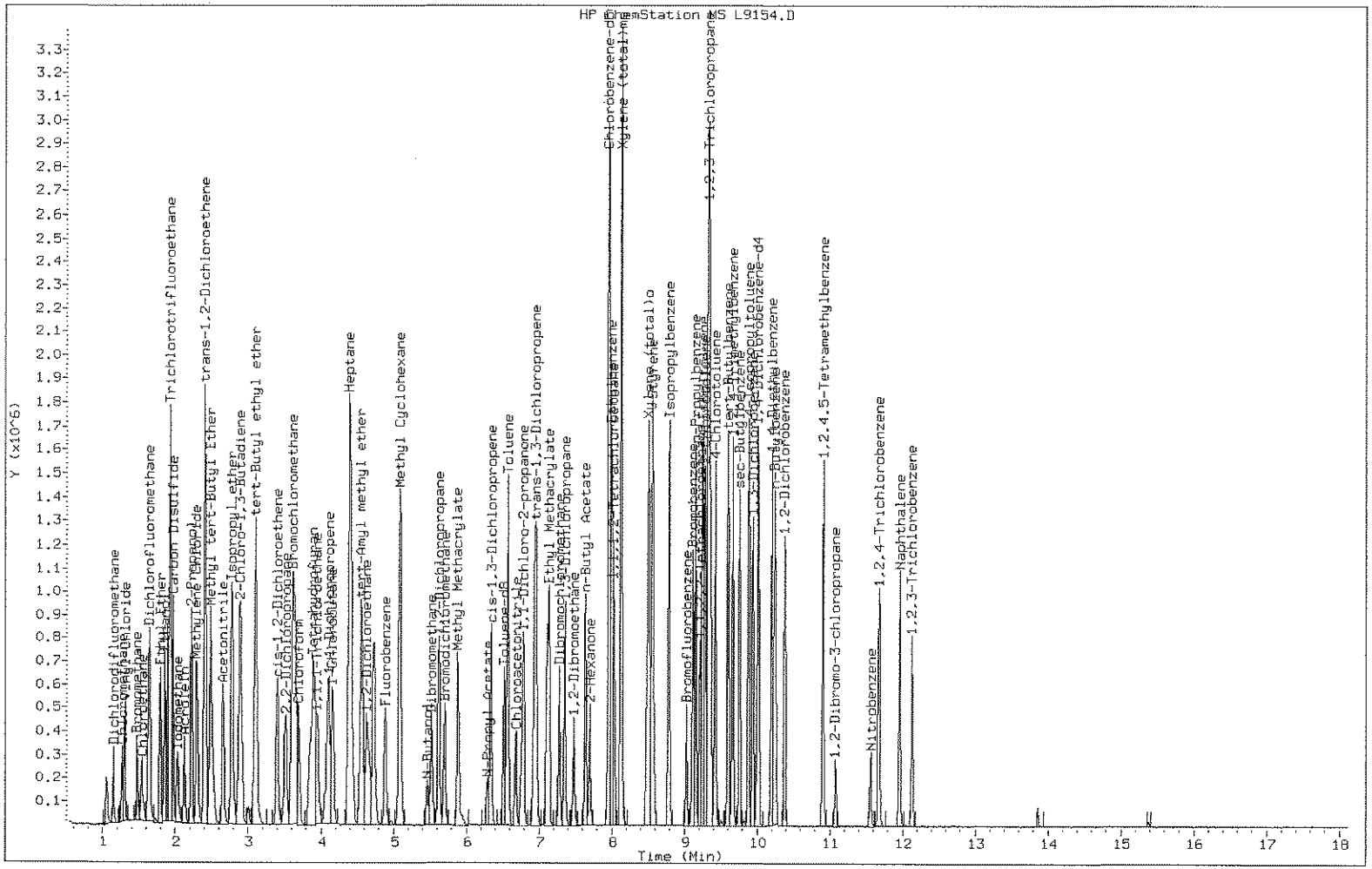
M - Compound response manually integrated.

Data File: L9154.D

Date: 26-JUL-2007 14:22

Client ID: IC;50 Instrument: msl.i

Sample Info: IC;50 Operator: D. HUMBERT



STL-CT

Volatile Report SW-846 Method 8260B

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079145.b\L9155.D
Lab Smp Id: IC;100 Client Smp ID: IC;100
Inj Date : 26-JUL-2007 14:47 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : IC;100
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079145.b\L8260BNW.m
Meth Date : 26-Jul-2007 15:45 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 14:47 Cal File: L9155.D
Als bottle: 16 Calibration Sample, Level: 4
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	AMOUNTS ON-COL (ug/L)
* 1	Fluorobenzene	96	4.882	4.882	(1.000)	442478	25.0000	
2	Dichlorodifluoromethane	85	1.143	1.143	(0.234)	356791	100.000	110
3	Chloromethane	50	1.261	1.261	(0.258)	620154	100.000	110
4	Vinyl Chloride	62	1.301	1.301	(0.266)	679439	100.000	110
5	Bromomethane	94	1.468	1.468	(0.301)	246624	100.000	110
6	Chloroethane	64	1.527	1.517	(0.313)	293707	100.000	71
7	Trichlorofluoromethane	101	1.606	1.596	(0.329)	651712	100.000	100
8	Dichlorofluoromethane	67	1.635	1.625	(0.335)	1210169	100.000	110
9	Ethyl Ether	45	1.792	1.783	(0.367)	401011	100.000	100
10	Ethanol	45	1.852	1.842	(0.379)	330278	1000.00	1100
11	Freon 141	81	1.852	1.842	(0.379)	811736	100.000	110
12	Freon 123a	67	1.635	1.625	(0.335)	1210169	100.000	110
13	Trichlorotrifluoroethane	101	1.930	1.930	(0.395)	461687	100.000	110
14	1,1-Dichloroethene	96	1.920	1.921	(0.393)	397036	100.000	110
15	Carbon Disulfide	76	1.960	1.960	(0.401)	1932253	100.000	110
16	Iodomethane	142	2.029	2.019	(0.416)	607881	100.000	150
17	Acrolein	56	2.127	2.117	(0.436)	571889	500.000	600
18	2-Propanol	45	2.216	2.235	(0.454)	111177	100.000	120
19	3-Chloro-1-Propene	41	2.216	2.216	(0.454)	1010971	100.000	110
20	Methylene Chloride	84	2.294	2.285	(0.470)	536290	100.000	110
21	Acetone	43	2.314	2.314	(0.474)	263638	100.000	90
22	trans-1,2-Dichloroethene	96	2.412	2.412	(0.494)	492941	100.000	110
23	Methyl Acetate	43	2.393	2.393	(0.490)	3426336	100.000	110

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	ON-COL (ug/L)
	24 Methyl tert-Butyl Ether	73	2.481	2.481	(0.508)	1913661	100.000	110
	25 tert-Butyl alcohol	59	2.530	2.540	(0.518)	546100	500.000	560
	26 Acetonitrile	41	2.648	2.658	(0.543)	1238371	1000.00	1100
	27 Isopropyl ether	45	2.767	2.767	(0.567)	2201842	100.000	100
	28 tert-Butyl ethyl ether	59	3.101	3.091	(0.635)	2228795	100.000	110
	29 2-Chloro-1,3-Butadiene	88	2.875	2.875	(0.589)	350323	100.000	100
	30 Acrylonitrile	53	2.914	2.914	(0.597)	676109	200.000	210
	31 1,1-Dichloroethane	63	2.885	2.885	(0.591)	1239866	100.000	100
	32 Vinyl Acetate	43	3.091	3.091	(0.633)	1688017	100.000	110
	33 cis-1,2-Dichloroethene	96	3.396	3.396	(0.696)	527903	100.000	110
	34 2,2-Dichloropropane	77	3.514	3.505	(0.720)	828211	100.000	100
	35 Bromochloromethane	128	3.613	3.613	(0.740)	373036	100.000	100
	36 1-Bromopropane	43	3.603	3.603	(0.738)	936559	100.000	100
	37 Cyclohexane	84	3.642	3.632	(0.746)	684576	100.000	110
	38 Chloroform	83	3.691	3.692	(0.756)	964204	100.000	110
	39 Ethyl Acetate	43	3.829	3.829	(0.784)	105777	200.000	220
	40 Methyl Acrylate	55	3.839	3.839	(0.786)	760136	100.000	110
$	41 Dibromofluoromethane	111	3.908	3.908	(0.800)	584066	100.000	110
	42 Tetrahydrofuran	42	3.898	3.888	(0.798)	515961	200.000	230
	43 Carbon Tetrachloride	117	3.878	3.878	(0.794)	756214	100.000	100
	44 1,1,1-Trichloroethane	97	3.947	3.947	(0.809)	710514	100.000	110
	45 2-Butanone	43	4.046	4.046	(0.829)	410047	100.000	100
	46 1,1-Dichloropropene	75	4.105	4.105	(0.841)	886915	100.000	110
	47 tert-Amyl methyl ether	73	4.557	4.548	(0.933)	1943373	100.000	110
	48 tert-Butyl formate	57	3.101	3.091	(0.635)	641545	100.000	110
	49 1-Chlorobutane	56	4.164	4.154	(0.853)	1260636	100.000	110
	50 Heptane	43	4.380	4.380	(0.897)	550540	100.000	100
	51 Propionitrile	54	4.380	4.390	(0.897)	1178501	1000.00	1100
	52 Benzene	78	4.400	4.400	(0.901)	2181153	100.000	100
	53 2-Methyl-2-Propenenitrile	41	4.420	4.420	(0.905)	847360	100.000	100
	54 Isobutyl alcohol	42	4.665	4.695	(0.956)	303812	1000.00	1100
$	55 1,2-Dichloroethane-d4	65	4.547	4.548	(0.931)	691324	100.000	100
	56 1,2-Dichloroethane	62	4.626	4.626	(0.948)	862943	100.000	100
	59 Methyl Cyclohexane	83	5.079	5.079	(1.040)	635389	100.000	100
	60 Trichloroethene	130	5.089	5.089	(1.042)	650766	100.000	110
	61 Isopropyl Acetate	43	4.675	4.695	(0.958)	513356	200.000	230
	62 N-Butanol	56	5.472	5.482	(1.121)	253242	1000.00	1200
	63 Dibromomethane	93	5.521	5.522	(1.131)	350026	100.000	110
	64 1,2-Dichloropropane	63	5.620	5.620	(1.151)	721454	100.000	110
	65 Bromodichloromethane	83	5.708	5.709	(1.169)	712704	100.000	110
	66 Methyl Methacrylate	69	5.886	5.886	(1.206)	653271	200.000	220
	67 1,4-Dioxane	58	5.925	5.935	(1.214)	61135	1000.00	1300
	68 N-Propyl Acetate	43	6.289	6.289	(1.288)	123827	200.000	240
	69 2-Chloroethylvinylether	63	6.289	6.289	(1.288)	187548	100.000	120
	70 cis-1,3-Dichloropropene	75	6.338	6.338	(1.298)	1051789	100.000	110
	71 Chloroacetonitrile	48	6.682	6.692	(1.369)	315510	1000.00	1200
	72 2-Nitropropane	41	6.761	6.761	(1.385)	400395	200.000	210
	73 trans-1,3-Dichloropropene	75	6.968	6.968	(1.427)	960720	100.000	110
	74 1,1,2-Trichloroethane	97	7.115	7.115	(1.457)	463542	100.000	110
*	75 Chlorobenzene-d5	117	7.952	7.952	(1.000)	426698	25.0000	
	76 Toluene	91	6.574	6.574	(0.827)	1889431	100.000	110
$	77 Toluene-d8	98	6.525	6.525	(0.821)	1553484	100.000	100
	78 1,1-Dichloro-2-propanone	43	6.791	6.791	(0.854)	2397294	500.000	570
	79 4-Methyl-2-Pentanone	43	6.928	6.929	(0.871)	838654	100.000	110

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	ON-COL (ug/L)
	80 Tetrachloroethene	164	6.948	6.948	(0.874)	378675	100.000	110
	81 Ethyl Methacrylate	69	7.135	7.135	(0.897)	1035803	100.000	110
	82 Dibromochloromethane	129	7.283	7.283	(0.916)	718297	100.000	110
	83 1,3-Dichloropropane	76	7.361	7.352	(0.926)	1013743	100.000	100
	84 1,2-Dibromoethane	107	7.479	7.480	(0.941)	608776	100.000	110
	85 n-Butyl Acetate	56	7.637	7.637	(0.960)	552358	100.000	120
	86 2-Hexanone	43	7.696	7.696	(0.968)	616249	100.000	110
	87 1-Chlorohexane	91	7.962	7.962	(1.001)	587051	100.000	94
	88 Chlorobenzene	112	7.971	7.962	(1.002)	1540664	100.000	110
	89 1,1,1,2-Tetrachloroethane	131	8.030	8.031	(1.010)	553854	100.000	110
	90 Ethylbenzene	106	8.001	8.001	(1.006)	703843	100.000	110
	91 Xylene (total)mp	106	8.129	8.129	(1.022)	1715186	200.000	220
	92 Xylene (total)o	106	8.513	8.513	(1.071)	833064	100.000	110
	93 Styrene	104	8.562	8.552	(1.077)	1412985	100.000	110
	94 Bromoform	173	8.581	8.581	(1.079)	413918	100.000	120
*	95 1,4-Dichlorobenzene-d4	152	10.008	10.008	(1.000)	165859	25.0000	
	96 Isopropylbenzene	105	8.788	8.788	(0.878)	2042067	100.000	100
	97 Bromobenzene	156	9.123	9.123	(0.912)	515637	100.000	100
	98 1,1,2,2-Tetrachloroethane	83	9.211	9.221	(0.920)	710126	100.000	100
	99 4-Ethyltoluene	105	9.250	9.251	(0.924)	2014031	100.000	100
	100 1,2,3-Trichloropropane	110	9.319	9.319	(0.931)	224742	100.000	100
	101 trans-1,4-Dichloro-2-Butene	53	9.359	9.359	(0.935)	505844	200.000	230
	102 n-Propylbenzene	91	9.152	9.152	(0.914)	2234294	100.000	100
	103 2-Chlorotoluene	91	9.280	9.280	(0.927)	1573858	100.000	110
	104 4-Chlorotoluene	91	9.428	9.428	(0.942)	1430506	100.000	100
	105 1,3,5-Trimethylbenzene	105	9.329	9.329	(0.932)	1652875	100.000	100
	106 tert-Butylbenzene	119	9.605	9.605	(0.960)	1587348	100.000	100
	107 1,2,4-Trimethylbenzene	105	9.664	9.664	(0.966)	1655398	100.000	100
	108 sec-Butylbenzene	105	9.762	9.752	(0.975)	1738017	100.000	100
	109 4-Isopropyltoluene	119	9.890	9.880	(0.988)	1864570	100.000	100
	110 1,3-Dichlorobenzene	146	9.939	9.939	(0.993)	915574	100.000	100
	111 1,4-Dichlorobenzene	146	10.018	10.018	(1.001)	944301	100.000	100
	112 1,2-Dichlorobenzene	146	10.382	10.382	(1.037)	902425	100.000	100
	113 Benzyl Chloride	126	10.234	10.234	(1.023)	299966	100.000	120
	114 1,4-Diethylbenzene	119	10.205	10.205	(2.090)	1071506	100.000	120
	115 n-Butylbenzene	91	10.244	10.244	(1.024)	2314449	100.000	110
	118 1,2,4,5-Tetramethylbenzene	119	10.903	10.904	(2.233)	1702829	100.000	130
	119 1,2-Dibromo-3-chloropropane	75	11.081	11.081	(1.107)	157805	100.000	120
	120 Nitrobenzene	77	11.563	11.563	(1.155)	431146	1000.00	2400(A)
	121 1,2,4-Trichlorobenzene	180	11.681	11.681	(1.167)	571406	100.000	140
	122 Hexachlorobutadiene	225	11.671	11.671	(1.166)	174987	100.000	110
	123 Naphthalene	128	11.966	11.956	(1.196)	1852158	100.000	160
	124 1,2,3-Trichlorobenzene	180	12.133	12.133	(1.212)	523254	100.000	140
$	125 Bromofluorobenzene	95	9.034	9.034	(0.903)	592570	100.000	94
M	126 1,2-Dichloroethene (total)	100				1020844	200.000	
M	127 Xylene (total)	100				2548250	300.000	

QC Flag Legend

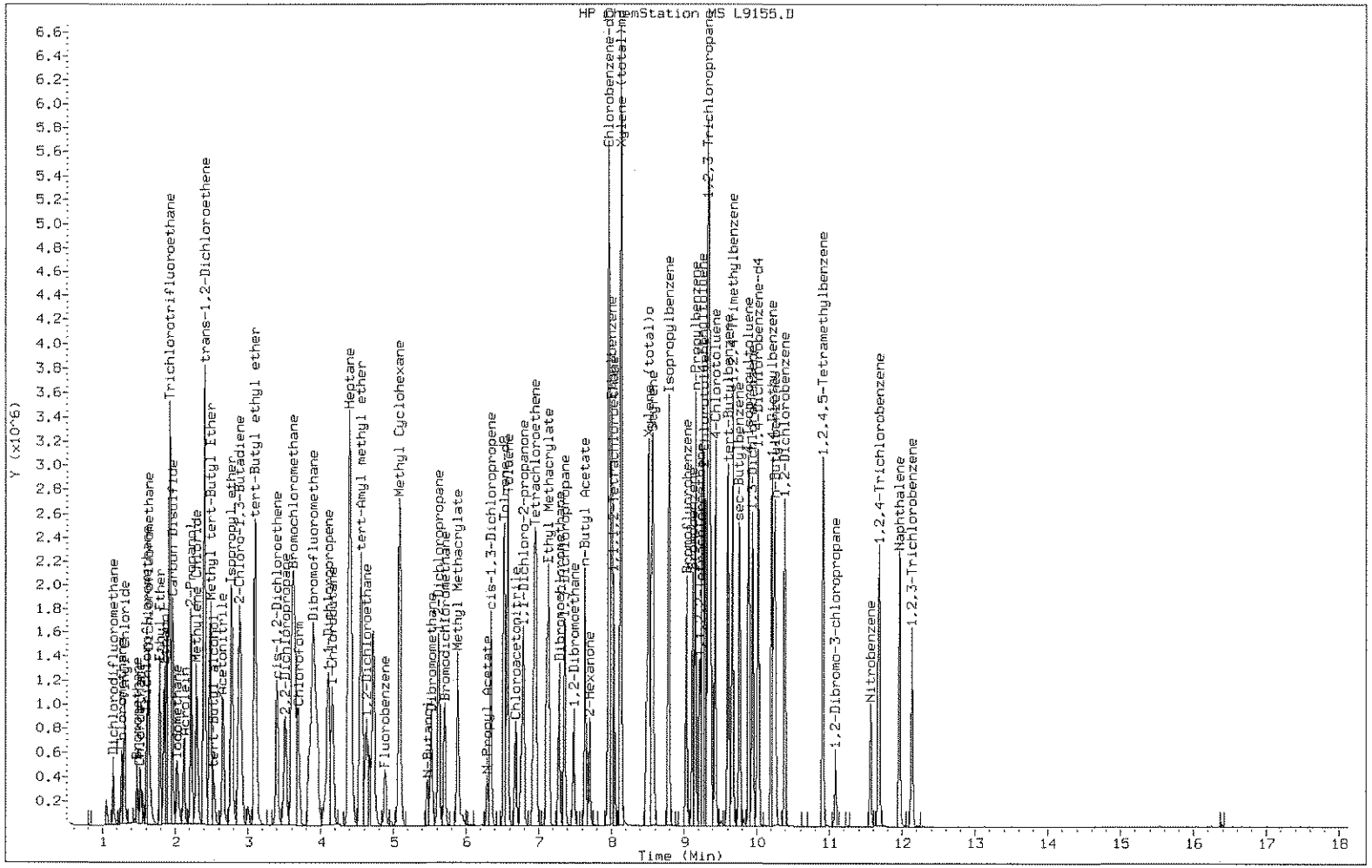
A - Target compound detected but, quantitated amount exceeded maximum amount.

Data File: L9155.D

Date: 26-JUL-2007 14:47

Client ID: IC;100 Instrument: msl.i

Sample Info: IC;100 Operator: D. HUMBERT



STL-CT

Volatile Report SW-846 Method 8260B

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079145.b\L9156.D
Lab Smp Id: IC;200 Client Smp ID: IC;200
Inj Date : 26-JUL-2007 15:11 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : IC;200
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079145.b\L8260BNW.m
Meth Date : 26-Jul-2007 15:45 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 17 Calibration Sample, Level: 5
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	ON-COL (ug/L)
* 1	Fluorobenzene	96	4.882	4.882	(1.000)	443394	25.0000	
2	Dichlorodifluoromethane	85	1.143	1.143	(0.234)	713398	200.000	210(A)
3	Chloromethane	50	1.261	1.261	(0.258)	1270571	200.000	220(A)
4	Vinyl Chloride	62	1.301	1.301	(0.266)	1394375	200.000	220(A)
5	Bromomethane	94	1.468	1.468	(0.301)	458216	200.000	200
6	Chloroethane	64	1.517	1.517	(0.311)	415164	200.000	110
7	Trichlorofluoromethane	101	1.596	1.596	(0.327)	1340237	200.000	210(A)
8	Dichlorofluoromethane	67	1.625	1.625	(0.333)	2475717	200.000	220(A)
9	Ethyl Ether	45	1.783	1.783	(0.365)	798164	200.000	200
10	Ethanol	45	1.842	1.842	(0.377)	685100	2000.00	2200(A)
11	Freon 141	81	1.842	1.842	(0.377)	1637853	200.000	210(A)
12	Freon 123a	67	1.625	1.625	(0.333)	2475717	200.000	220(A)
13	Trichlorotrifluoroethane	101	1.930	1.930	(0.395)	941828	200.000	220(A)
14	1,1-Dichloroethene	96	1.921	1.921	(0.393)	795431	200.000	210(A)
15	Carbon Disulfide	76	1.960	1.960	(0.402)	3972137	200.000	220(A)
16	Iodomethane	142	2.019	2.019	(0.414)	1257993	200.000	280(A)
17	Acrolein	56	2.117	2.117	(0.434)	1126607	1000.00	1100(A)
18	2-Propanol	45	2.235	2.235	(0.458)	238541	200.000	240(A)
19	3-Chloro-1-Propene	41	2.216	2.216	(0.454)	2027040	200.000	210(A)
20	Methylene Chloride	84	2.285	2.285	(0.468)	1094414	200.000	210(A)
21	Acetone	43	2.314	2.314	(0.474)	532218	200.000	190
22	trans-1,2-Dichloroethene	96	2.412	2.412	(0.494)	1005327	200.000	220(A)
23	Methyl Acetate	43	2.393	2.393	(0.490)	6976078	200.000	210(A)

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	ON-COL (ug/L)
	24 Methyl tert-Butyl Ether	73	2.481	2.481	(0.508)	3886103	200.000	210(A)
	25 tert-Butyl alcohol	59	2.540	2.540	(0.520)	1082658	1000.00	1100(AM)
	26 Acetonitrile	41	2.658	2.658	(0.545)	2464678	2000.00	2100(A)
	27 Isopropyl ether	45	2.767	2.767	(0.567)	4508414	200.000	210(A)
	28 tert-Butyl ethyl ether	59	3.091	3.091	(0.633)	4538425	200.000	210(A)
	29 2-Chloro-1,3-Butadiene	88	2.875	2.875	(0.589)	727607	200.000	220(A)
	30 Acrylonitrile	53	2.914	2.914	(0.597)	1348639	400.000	420(A)
	31 1,1-Dichloroethane	63	2.885	2.885	(0.591)	2574069	200.000	220(A)
	32 Vinyl Acetate	43	3.091	3.091	(0.633)	3433072	200.000	220(A)
	33 cis-1,2-Dichloroethene	96	3.396	3.396	(0.696)	1076027	200.000	220(A)
	34 2,2-Dichloropropane	77	3.505	3.505	(0.718)	1694681	200.000	200(A)
	35 Bromochloromethane	128	3.613	3.613	(0.740)	760506	200.000	210(A)
	36 1-Bromopropane	43	3.603	3.603	(0.738)	1912344	200.000	210(A)
	37 Cyclohexane	84	3.632	3.632	(0.744)	1418972	200.000	220(A)
	38 Chloroform	83	3.692	3.692	(0.756)	1962248	200.000	210(A)
	39 Ethyl Acetate	43	3.829	3.829	(0.784)	202164	400.000	410(A)
	40 Methyl Acrylate	55	3.839	3.839	(0.786)	1526298	200.000	220(A)
$	41 Dibromofluoromethane	111	3.908	3.908	(0.800)	1240771	200.000	220(A)
	42 Tetrahydrofuran	42	3.888	3.888	(0.796)	1028591	400.000	440(A)
	43 Carbon Tetrachloride	117	3.878	3.878	(0.794)	1788034	200.000	240(A)
	44 1,1,1-Trichloroethane	97	3.947	3.947	(0.809)	1461749	200.000	220(A)
	45 2-Butanone	43	4.046	4.046	(0.829)	831761	200.000	200(A)
	46 1,1-Dichloropropene	75	4.105	4.105	(0.841)	1797720	200.000	220(A)
	47 tert-Amyl methyl ether	73	4.548	4.548	(0.931)	3999220	200.000	220(A)
	48 tert-Butyl formate	57	3.091	3.091	(0.633)	1282744	200.000	210(A)
	49 1-Chlorobutane	56	4.154	4.154	(0.851)	2551737	200.000	210(A)
	50 Heptane	43	4.380	4.380	(0.897)	1119150	200.000	200(A)
	51 Propionitrile	54	4.390	4.390	(0.899)	2400934	2000.00	2200(A)
	52 Benzene	78	4.400	4.400	(0.901)	4473567	200.000	210(A)
	53 2-Methyl-2-Propenenitrile	41	4.420	4.420	(0.905)	1316560	200.000	160
	54 Isobutyl alcohol	42	4.695	4.695	(0.962)	636170	2000.00	2300(A)
$	55 1,2-Dichloroethane-d4	65	4.548	4.548	(0.931)	1472282	200.000	210(A)
	56 1,2-Dichloroethane	62	4.626	4.626	(0.948)	1771871	200.000	210(A)
	59 Methyl Cyclohexane	83	5.079	5.079	(1.040)	1314289	200.000	210(A)
	60 Trichloroethene	130	5.089	5.089	(1.042)	1322325	200.000	210(A)
	61 Isopropyl Acetate	43	4.695	4.695	(0.962)	1026258	400.000	440(A)
	62 N-Butanol	56	5.482	5.482	(1.123)	523419	2000.00	2300(A)
	63 Dibromomethane	93	5.522	5.522	(1.131)	730368	200.000	220(A)
	64 1,2-Dichloropropane	63	5.620	5.620	(1.151)	1494377	200.000	220(A)
	65 Bromodichloromethane	83	5.709	5.709	(1.169)	1453591	200.000	220(A)
	66 Methyl Methacrylate	69	5.886	5.886	(1.206)	1325545	400.000	440(A)
	67 1,4-Dioxane	58	5.935	5.935	(1.216)	124966	2000.00	2400
	68 N-Propyl Acetate	43	6.289	6.289	(1.288)	257957	400.000	480(A)
	69 2-Chloroethylvinylether	63	6.289	6.289	(1.288)	390909	200.000	240(A)
	70 cis-1,3-Dichloropropene	75	6.338	6.338	(1.298)	2178368	200.000	220(A)
	71 Chloroacetonitrile	48	6.692	6.692	(1.371)	645921	2000.00	2300
	72 2-Nitropropane	41	6.761	6.761	(1.385)	850899	400.000	440(A)
	73 trans-1,3-Dichloropropene	75	6.968	6.968	(1.427)	1997362	200.000	220(A)
	74 1,1,2-Trichloroethane	97	7.115	7.115	(1.457)	942127	200.000	220(A)
*	75 Chlorobenzene-d5	117	7.952	7.952	(1.000)	433942	25.0000	
	76 Toluene	91	6.574	6.574	(0.827)	3895334	200.000	210(A)
$	77 Toluene-d8	98	6.525	6.525	(0.821)	3278667	200.000	220(A)
	78 1,1-Dichloro-2-propanone	43	6.791	6.791	(0.854)	4919221	1000.00	1100(A)
	79 4-Methyl-2-Pentanone	43	6.929	6.929	(0.871)	1652708	200.000	210(A)

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	ON-COL (ug/L)
80	Tetrachloroethene	164	6.948	6.948	(0.874)	762569	200.000	210(A)
81	Ethyl Methacrylate	69	7.135	7.135	(0.897)	2078100	200.000	210(A)
82	Dibromochloromethane	129	7.283	7.283	(0.916)	1515965	200.000	220(A)
83	1,3-Dichloropropane	76	7.352	7.352	(0.925)	2116884	200.000	210(A)
84	1,2-Dibromoethane	107	7.480	7.480	(0.941)	1243668	200.000	220(A)
85	n-Butyl Acetate	56	7.637	7.637	(0.960)	1089091	200.000	220(A)
86	2-Hexanone	43	7.696	7.696	(0.968)	1261486	200.000	210(A)
87	1-Chlorohexane	91	7.962	7.962	(1.001)	1234202	200.000	200
88	Chlorobenzene	112	7.962	7.962	(1.001)	3190028	200.000	220(A)
89	1,1,1,2-Tetrachloroethane	131	8.031	8.031	(1.010)	1170387	200.000	220(A)
90	Ethylbenzene	106	8.001	8.001	(1.006)	1438586	200.000	220(A)
91	Xylene (total)mp	106	8.129	8.129	(1.022)	3514018	400.000	430(A)
92	Xylene (total)o	106	8.513	8.513	(1.071)	1692965	200.000	210(A)
93	Styrene	104	8.552	8.552	(1.075)	2868190	200.000	220(A)
94	Bromoform	173	8.581	8.581	(1.079)	854613	200.000	230(A)
* 95	1,4-Dichlorobenzene-d4	152	10.008	10.008	(1.000)	163940	25.0000	
96	Isopropylbenzene	105	8.788	8.788	(0.878)	4113409	200.000	200(A)
97	Bromobenzene	156	9.123	9.123	(0.912)	1025285	200.000	200(A)
98	1,1,2,2-Tetrachloroethane	83	9.221	9.221	(0.921)	1411843	200.000	200(A)
99	4-Ethyltoluene	105	9.251	9.251	(0.924)	4087786	200.000	210(A)
100	1,2,3-Trichloropropane	110	9.319	9.319	(0.931)	437299	200.000	200
101	trans-1,4-Dichloro-2-Butene	53	9.359	9.359	(0.935)	950270	400.000	420(A)
102	n-Propylbenzene	91	9.152	9.152	(0.914)	4479848	200.000	210(A)
103	2-Chlorotoluene	91	9.280	9.280	(0.927)	3180389	200.000	220(A)
104	4-Chlorotoluene	91	9.428	9.428	(0.942)	2886209	200.000	210(A)
105	1,3,5-Trimethylbenzene	105	9.329	9.329	(0.932)	3296100	200.000	200(A)
106	tert-Butylbenzene	119	9.605	9.605	(0.960)	3241838	200.000	210(A)
107	1,2,4-Trimethylbenzene	105	9.664	9.664	(0.966)	3363577	200.000	210(A)
108	sec-Butylbenzene	105	9.752	9.752	(0.974)	3538238	200.000	210(A)
109	4-Isopropyltoluene	119	9.880	9.880	(0.987)	3853442	200.000	210(A)
110	1,3-Dichlorobenzene	146	9.939	9.939	(0.993)	1829937	200.000	210(A)
111	1,4-Dichlorobenzene	146	10.018	10.018	(1.001)	1905362	200.000	210(A)
112	1,2-Dichlorobenzene	146	10.382	10.382	(1.037)	1798756	200.000	200(A)
113	Benzyl Chloride	126	10.234	10.234	(1.023)	609218	200.000	230(A)
114	1,4-Diethylbenzene	119	10.205	10.205	(2.090)	2181355	200.000	230(A)
115	n-Butylbenzene	91	10.244	10.244	(1.024)	4862125	200.000	220(A)
118	1,2,4,5-Tetramethylbenzene	119	10.904	10.904	(2.233)	3497909	200.000	250(A)
119	1,2-Dibromo-3-chloropropane	75	11.081	11.081	(1.107)	325207	200.000	240(A)
120	Nitrobenzene	77	11.563	11.563	(1.155)	1166858	2000.00	4800(A)
121	1,2,4-Trichlorobenzene	180	11.681	11.681	(1.167)	1199072	200.000	270(A)
122	Hexachlorobutadiene	225	11.671	11.671	(1.166)	350760	200.000	220(A)
123	Naphthalene	128	11.956	11.956	(1.195)	4033323	200.000	310(A)
124	1,2,3-Trichlorobenzene	180	12.133	12.133	(1.212)	1116273	200.000	280(A)
$ 125	Bromofluorobenzene	95	9.034	9.034	(0.903)	1225319	200.000	200
M 126	1,2-Dichloroethene (total)	100				2081354	400.000	
M 127	Xylene (total)	100				5206983	600.000	

QC Flag Legend

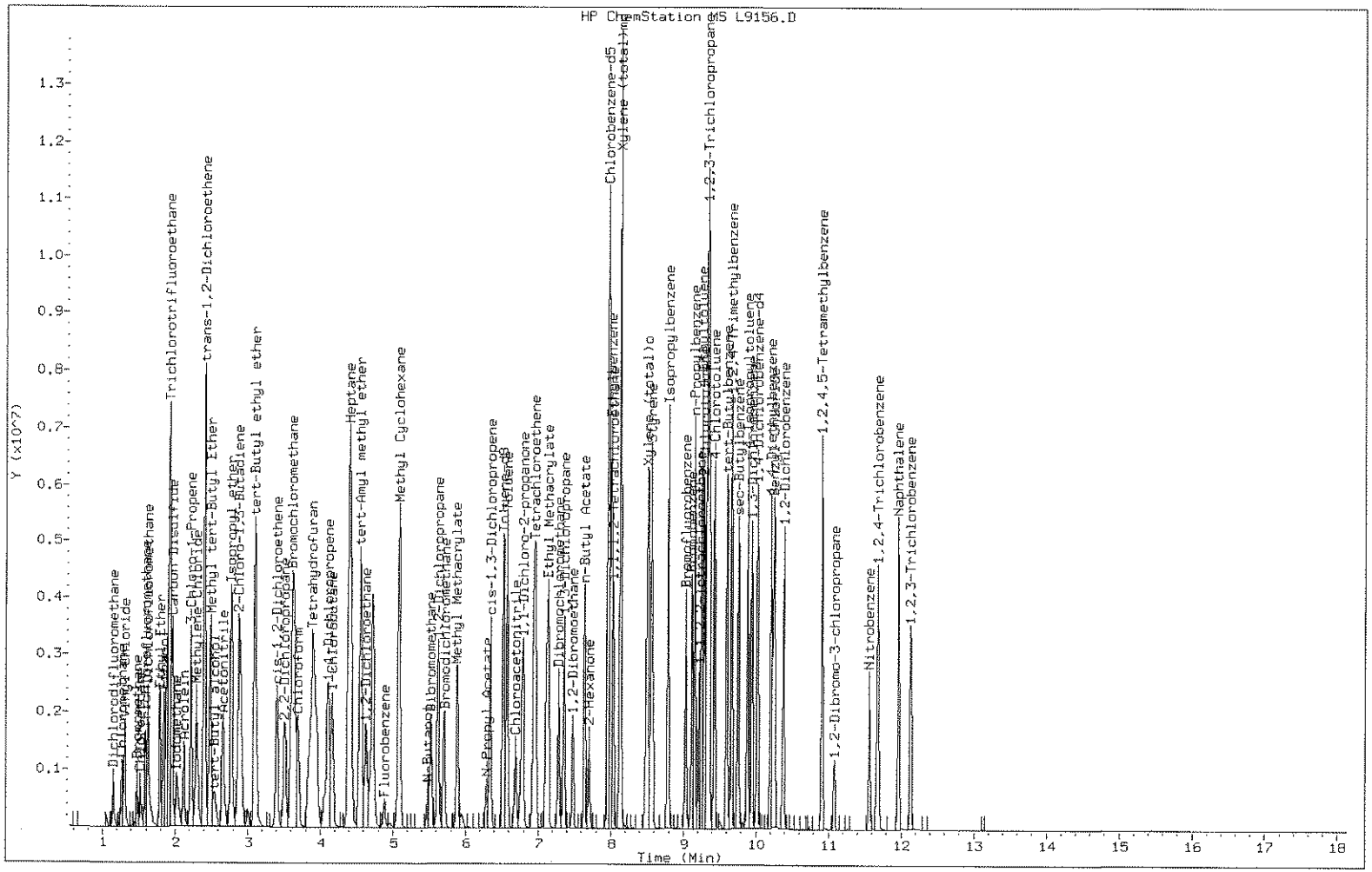
A - Target compound detected but, quantitated amount exceeded maximum amount.
M - Compound response manually integrated.

Data File: L9156.D

Date: 26-JUL-2007 15:11

Client ID: IC;200 Instrument: msl.i

Sample Info: IC;200 Operator: D. HUMBERT



VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Instrument ID: MSL Calibration Date: 07/31/2007 Time: 20:21

Lab File ID: L9310.D Init. Calib. Date(s): 07/26/2007 07/26/2007

Lab Sample ID: CCVIS 220-8307/1 Init. Calib. Time(s): 13:32 15:11

GC Column: RTX-VMS ID: 0.18 (mm) Heated Purge: (Y/N) N Conc. Units: ug/L

Analyte	Curve	Ave RRF	RRF	Min RRF	Calculated Amount	Ccal Amount	% D	Max % D
1,1,1,2-Tetrachloroethane	Ave	0.3128	0.3162		50.5	50.0	1.1	
1,1,1-Trichloroethane	Ave	0.3840	0.4103		53.4	50.0	6.9	
1,1,2,2-Tetrachloroethane	Ave	1.0638	1.0561	0.3000	49.6	50.0	-0.7	
1,1,2-Trichloro-1,2,2-trifluoroethane	Ave	0.2475	0.2706		54.7	50.0	9.4	
1,1,2-Trichloroethane	Ave	0.2502	0.2669		53.3	50.0	6.7	
1,1-Dichloro-1-fluoroethane	Ave	0.4422	0.4703		53.2	50.0	6.4	
1,1-Dichloroacetone	Ave	0.2611	0.2429		233	250	-6.9	
1,1-Dichloroethane	Ave	0.6829	0.7198	0.1000	52.7	50.0	5.4	
1,1-Dichloroethene	Ave	0.2131	0.2306		54.1	50.0	8.2	20.0
1,1-Dichloropropene	Ave	0.4776	0.5060		53.0	50.0	5.9	
1,2,3-Trichlorobenzene	Ave	0.6643	0.7272		54.7	50.0	9.5	
1,2,3-Trichloropropane	Ave	0.3371	0.3565		52.9	50.0	5.8	
1,2,4,5-Tetramethylbenzene	Ave	0.8285	0.9268		55.9	50.0	11.9	
1,2,4-Trichlorobenzene	Ave	0.7280	0.8107		55.7	50.0	11.4	
1,2,4-Trimethylbenzene	Ave	2.4805	2.6112		52.6	50.0	5.3	
1,2-Dibromo-3-Chloropropane	Ave	0.2124	0.2207		52.0	50.0	3.9	
1,2-Dichloro-1,1,2-trifluoroethane	Ave	0.6540			53.9			
1,2-Dichlorobenzene	Ave	1.3400	1.3844		51.7	50.0	3.3	
1,2-Dichloroethane	Ave	0.4746	0.5069		53.4	50.0	6.8	
1,2-Dichloroethane-d4 (Surr)	Ave	0.3932	0.2934		18.7	25.0	-25.4	
1,2-Dichloroethene, Total	Ave	0.2306	0.2910		105	100	26.2	
1,2-Dichloropropane	Ave	0.3960	0.4199		53.0	50.0	6.0	20.0
1,3,5-Trimethylbenzene	Ave	2.4694	2.5755		52.1	50.0	4.3	
1,3-Dichlorobenzene	Ave	1.3616	1.4158		52.0	50.0	4.0	
1,3-Dichloropropane	Ave	0.5774	0.5772		50.0	50.0	0	
1,4-Dichlorobenzene	Ave	1.3897	1.4163		51.0	50.0	1.9	
1,4-Dioxane	Ave	0.0030	0.0033		549	500	9.8	
1-Bromopropane	Ave	0.5150	0.5550		53.9	50.0	7.8	
1-Chlorobutane	Ave	0.6854	0.7342		53.6	50.0	7.1	
1-Chlorohexane	Ave	0.3605	0.3952		54.8	50.0	9.6	
2,2-Dichloropropane	Ave	0.4711	0.4820		51.2	50.0	2.3	
2,4,4-Trimethyl-1-pentene	Ave				0.0			
2-Chloro-1,3-butadiene	Ave	0.1935	0.2054		53.1	50.0	6.2	
2-Chloroethyl vinyl ether	Ave	0.0965	0.1155		59.9	50.0	19.8	
2-Chlorotoluene	Ave	2.2771	2.4558		53.9	50.0	7.8	
2-Hexanone	Ave	0.3432	0.3744		54.5	50.0	9.1	
2-Methyl-2-propanol	Ave	0.0579	0.0636		274	250	9.8	
2-Nitropropane	Ave	0.1112	0.1099		98.9	100	-1.1	
3-Chloro-1-propene	Ave	0.5499	0.5644		51.3	50.0	2.6	
4-Bromofluorobenzene	Ave	0.9379	0.8713		23.2	25.0	-7.1	
4-Chlorotoluene	Ave	2.1467	2.2143		51.6	50.0	3.1	
4-Ethyltoluene	Ave	3.0136	3.1651		52.5	50.0	5.0	

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Instrument ID: MSL Calibration Date: 07/31/2007 Time: 20:21

Lab File ID: L9310.D Init. Calib. Date(s): 07/26/2007 07/26/2007

Lab Sample ID: CCVIS 220-8307/1 Init. Calib. Time(s): 13:32 15:11

GC Column: RTX-VMS ID: 0.18 (mm) Heated Purge: (Y/N) N Conc. Units: ug/L

Analyte	Curve	Ave RRF	RRF	Min RRF	Calculated Amount	Ccal Amount	% D	Max % D
4-Isopropyltoluene	Ave	2.7945	2.8869		51.7	50.0	3.3	
Acetone	Ave	0.1585	0.1838		58.0	50.0	16.0	
Acetonitrile	Ave	0.0661	0.0716		541	500	8.2	
Acrolein	Ave	0.0580	0.0721		311	250	24.3	
Acrylonitrile	Ave	0.1834	0.1880		103	100	2.5	
Benzene	Ave	1.2001	1.2565		52.3	50.0	4.7	
Benzyl chloride	Ave	0.4143	0.4215		50.9	50.0	1.7	
Bromobenzene	Ave	0.7785	0.7987		51.3	50.0	2.6	
Bromoform	Ave	0.2207	0.2136	0.1000	48.4	50.0	-3.2	
Bromomethane	Ave	0.1293	0.1219		47.1	50.0	-5.7	
Carbon disulfide	Ave	1.0400	1.0147		48.8	50.0	-2.4	
Carbon tetrachloride	Ave	0.4301	0.4273		49.7	50.0	-0.7	
Chloroacetonitrile	Ave	0.0162	0.0173		533	500	6.5	
Chlorobenzene	Ave	0.8648	0.9056	0.3000	52.4	50.0	4.7	
Chlorobromomethane	Ave	0.2047	0.2170		53.0	50.0	6.0	
Chlorodibromomethane	Ave	0.4000	0.3958		49.5	50.0	-1.1	
Chloroethane	Ave	0.1962	0.2488		63.4	50.0	26.8	
Chloroform	Ave	0.5224	0.5560		53.2	50.0	6.4	20.0
Chloromethane	Ave	0.3363	0.3326	0.1000	49.5	50.0	-1.1	
cis-1,2-Dichloroethene	Ave	0.2859	0.3032		53.0	50.0	6.0	
cis-1,3-Dichloropropene	Ave	0.5664	0.5831		51.5	50.0	2.9	
Cyclohexane	Ave	0.3652	0.3972		54.4	50.0	8.8	
Dibromofluoromethane	Ave	0.3225	0.2485		19.3	25.0	-23.0	
Dibromomethane	Ave	0.1924	0.2012		52.3	50.0	4.6	
Dichlorobromomethane	Ave	0.3741	0.3803		50.8	50.0	1.7	
Dichlorodifluoromethane	Ave	0.1920	0.1628		42.4	50.0	-15.2	
Dichlorofluoromethane	Ave	0.6540	0.7050		53.9	50.0	7.8	
Ethanol	Ave	0.0178	0.0195		548	500	9.5	
Ethyl acetate	Ave	0.0280	0.0292		104	100	4.5	
Ethyl ether	Ave	0.2255	0.2390		53.0	50.0	6.0	
Ethyl methacrylate	Ave	0.5666	0.5937		52.4	50.0	4.8	
Ethylbenzene	Ave	0.3850	0.4012		52.1	50.0	4.2	20.0
Ethylene Dibromide	Ave	0.3366	0.3433		51.0	50.0	2.0	
Hexachlorobutadiene	Ave	0.2481	0.2421		48.8	50.0	-2.4	
Hexachloroethane	Ave				0.0			
Iodomethane	Ave	0.2764	0.2185		39.5	50.0	-20.9	
Isobutanol	Ave	0.0162	0.0174		539	500	7.8	
Isopropyl acetate	Ave	0.1342	0.1432		107	100	6.8	
Isopropyl alcohol	Ave	0.0574	0.0656		57.2	50.0	14.4	
Isopropyl ether	Ave	1.2074	1.2781		52.9	50.0	5.9	
Isopropylbenzene	Ave	3.0830	3.2023		51.9	50.0	3.9	
m-Xylene & p-Xylene	Ave	0.4812	0.5035		105	100	4.6	
Methacrylonitrile	Ave	0.4469	0.4877		54.6	50.0	9.1	

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Instrument ID: MSL Calibration Date: 07/31/2007 Time: 20:21

Lab File ID: L9310.D Init. Calib. Date(s): 07/26/2007 07/26/2007

Lab Sample ID: CCVIS 220-8307/1 Init. Calib. Time(s): 13:32 15:11

GC Column: RTX-VMS ID: 0.18 (mm) Heated Purge: (Y/N) N Conc. Units: ug/L

Analyte	Curve	Ave RRF	RRF	Min RRF	Calculated Amount	Ccal Amount	% D	Max % D
Methyl acetate	Ave	1.8684	1.9666		52.6	50.0	5.3	
Methyl acrylate	Ave	0.4003	0.4230		52.8	50.0	5.7	
Methyl Ethyl Ketone	Ave	0.2301	0.2494		54.2	50.0	8.4	
methyl isobutyl ketone	Ave	0.4602	0.4889		53.1	50.0	6.2	
Methyl methacrylate	Ave	0.1736	0.3656		105	50.0	111	
Methyl tert-butyl ether	Ave	1.0379	1.0700		51.5	50.0	3.1	
Methylcyclohexane	Ave	0.3498	0.3649		52.2	50.0	4.3	
Methylene Chloride	Ave	0.2929	0.3035		51.8	50.0	3.6	
n-Butanol	Ave	0.0130	0.0145		558	500	11.6	
n-Butyl acetate	Ave	0.2901	0.3144		54.2	50.0	8.4	
n-Butylbenzene	Ave	3.3672	3.4277		50.9	50.0	1.8	
n-Heptane	Ave	0.3090	0.3085		49.9	50.0	-0.2	
n-Propyl acetate	Ave	0.0316	0.0400		126	100	26.3	
N-Propylbenzene	Ave	3.3039	3.4345		52.0	50.0	4.0	
Naphthalene	Ave	2.2068	2.7855		63.1	50.0	26.2	
Nitrobenzene	Ave	0.0473	0.0417		440	500	-11.9	
o-Xylene	Ave	0.4609	0.4936		53.5	50.0	7.1	
p-Diethylbenzene	Ave	0.5445	0.5975		54.9	50.0	9.7	
Pentachloroethane	Ave				0.0			
Propionitrile	Ave	0.0634	0.0694		547	500	9.5	
sec-Butylbenzene	Ave	2.6102	2.6807		51.4	50.0	2.7	
Styrene	Ave	0.7721	0.8121		52.6	50.0	5.2	
Tert-amyl methyl ether	Ave	1.0524	1.1133		52.9	50.0	5.8	
Tert-butyl ethyl ether	Ave	1.2132	1.2972		53.5	50.0	6.9	
tert-Butyl Formate	Ave	0.3499	0.3698		52.8	50.0	5.7	
tert-Butylbenzene	Ave	2.3884	2.4626		51.6	50.0	3.1	
Tetrachloroethene	Ave	0.2100	0.2189		52.1	50.0	4.2	
Tetrahydrofuran	Ave	0.1355	0.1487		110	100	9.8	
Toluene	Ave	1.0645	1.0944		51.4	50.0	2.8	20.0
Toluene-d8 (Surr)	Ave	0.8889	0.7015		19.7	25.0	-21.1	
trans-1,2-Dichloroethene	Ave	0.2676	0.2789		52.1	50.0	4.2	
trans-1,3-Dichloropropene	Ave	0.5141	0.5404		52.6	50.0	5.1	
trans-1,4-Dichloro-2-butene	Ave	0.3461	0.3235		93.5	100	-6.5	
Trichloroethene	Ave	0.3531	0.3884		55.0	50.0	10.0	
Trichlorofluoromethane	Ave	0.3584	0.3917		54.6	50.0	9.3	
Vinyl acetate	Ave	0.9114	0.7874		43.2	50.0	-13.6	
Vinyl chloride	Ave	0.3701	0.3878		52.4	50.0	4.8	20.0
Xylenes, Total	Ave	0.3953	0.5002		158	150	26.5	

Curve Types:
Ave = Average
Lin = Linear
Quad = Quadratic

STL-CT

Volatile Report SW-846 Method 8260B

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079310.b\L9310.D
Lab Smp Id: CCVIS
Inj Date : 31-JUL-2007 20:21 MS Autotune Date: 26-APR-2004 14:21
Operator : D. GAYDA Inst ID: msl.i
Smp Info : CCVIS
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079310.b\L8260BNW.m
Meth Date : 31-Jul-2007 21:02 ctvoa Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 1 Continuing Calibration Sample
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONMSLNT

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	AMOUNTS ON-COL (ug/L)
* 1	Fluorobenzene	96	4.896	4.896	(1.000)	439656	25.0000	
2	Dichlorodifluoromethane	85	1.147	1.147	(0.234)	143111	50.0000	42
3	Chloromethane	50	1.265	1.265	(0.258)	292482	50.0000	49
4	Vinyl Chloride	62	1.304	1.304	(0.267)	341014	50.0000	52
5	Bromomethane	94	1.482	1.482	(0.303)	107180	50.0000	47
6	Chloroethane	64	1.541	1.541	(0.315)	218794	50.0000	63
7	Trichlorofluoromethane	101	1.629	1.629	(0.333)	344396	50.0000	55
8	Dichlorofluoromethane	67	1.649	1.649	(0.337)	619928	50.0000	54
9	Ethyl Ether	45	1.796	1.796	(0.367)	210133	50.0000	53
10	Ethanol	45	1.865	1.865	(0.381)	171112	500.000	550
11	Freon 141	81	1.855	1.855	(0.379)	413505	50.0000	53
12	Freon 123a	67	1.649	1.649	(0.337)	619928	50.0000	54
13	Trichlorotrifluoroethane	101	1.944	1.944	(0.397)	237967	50.0000	55
14	1,1-Dichloroethene	96	1.934	1.934	(0.395)	202750	50.0000	54
15	Carbon Disulfide	76	1.973	1.973	(0.403)	892250	50.0000	49
16	Iodomethane	142	2.033	2.033	(0.415)	192104	50.0000	40
17	Acrolein	56	2.131	2.131	(0.435)	317056	250.000	310
18	2-Propanol	45	2.219	2.219	(0.453)	57715	50.0000	57
19	3-Chloro-1-Propene	41	2.229	2.229	(0.455)	496254	50.0000	51
20	Methylene Chloride	84	2.298	2.298	(0.469)	266873	50.0000	52
21	Acetone	43	2.328	2.328	(0.476)	161654	50.0000	58
22	trans-1,2-Dichloroethene	96	2.416	2.416	(0.494)	245195	50.0000	52

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	AMOUNTS ON-COL (ug/L)
	23 Methyl Acetate	43	2.406	2.406	(0.492)	1729237	50.0000	53
	24 Methyl tert-Butyl Ether	73	2.495	2.495	(0.510)	940842	50.0000	52
	25 tert-Butyl alcohol	59	2.534	2.534	(0.518)	279438	250.000	270
	26 Acetonitrile	41	2.662	2.662	(0.544)	629409	500.000	540
	27 Isopropyl ether	45	2.780	2.780	(0.568)	1123878	50.0000	53
	28 tert-Butyl ethyl ether	59	3.105	3.105	(0.634)	1140619	50.0000	53
	29 2-Chloro-1,3-Butadiene	88	2.888	2.888	(0.590)	180629	50.0000	53
	30 Acrylonitrile	53	2.918	2.918	(0.596)	330643	100.000	100
	31 1,1-Dichloroethane	63	2.898	2.898	(0.592)	632893	50.0000	53
	32 Vinyl Acetate	43	3.105	3.105	(0.634)	692329	50.0000	43
	33 cis-1,2-Dichloroethene	96	3.410	3.410	(0.697)	266583	50.0000	53
	34 2,2-Dichloropropane	77	3.528	3.528	(0.721)	423807	50.0000	51
	35 Bromochloromethane	128	3.626	3.626	(0.741)	190776	50.0000	53
	36 1-Bromopropane	43	3.617	3.617	(0.739)	487991	50.0000	54
	37 Cyclohexane	84	3.656	3.656	(0.747)	349234	50.0000	54
	38 Chloroform	83	3.705	3.705	(0.757)	488870	50.0000	53
	39 Ethyl Acetate	43	3.843	3.843	(0.785)	51369	100.000	100
	40 Methyl Acrylate	55	3.853	3.853	(0.787)	371954	50.0000	53
$	41 Dibromofluoromethane	111	3.922	3.922	(0.801)	109252	25.0000	19
	42 Tetrahydrofuran	42	3.912	3.912	(0.799)	261549	100.000	110
	43 Carbon Tetrachloride	117	3.892	3.892	(0.795)	375743	50.0000	50
	44 1,1,1-Trichloroethane	97	3.961	3.961	(0.809)	360778	50.0000	53
	45 2-Butanone	43	4.069	4.069	(0.831)	219301	50.0000	54
	46 1,1-Dichloropropene	75	4.118	4.118	(0.841)	444936	50.0000	53
	47 tert-Amyl methyl ether	73	4.571	4.571	(0.934)	978907	50.0000	53
	48 tert-Butyl formate	57	3.105	3.105	(0.634)	325145	50.0000	53
	49 1-Chlorobutane	56	4.177	4.177	(0.853)	645592	50.0000	54
	50 Heptane	43	4.394	4.394	(0.898)	271227	50.0000	50
	51 Propionitrile	54	4.394	4.394	(0.898)	610505	500.000	550
	52 Benzene	78	4.414	4.414	(0.902)	1104847	50.0000	52
	53 2-Methyl-2-Propenenitrile	41	4.433	4.433	(0.906)	428854	50.0000	54
	54 Isobutyl alcohol	42	4.679	4.679	(0.956)	153133	500.000	540
$	55 1,2-Dichloroethane-d4	65	4.561	4.561	(0.932)	128977	25.0000	19
	56 1,2-Dichloroethane	62	4.640	4.640	(0.948)	445685	50.0000	53
	59 Methyl Cyclohexane	83	5.092	5.092	(1.040)	320867	50.0000	52
	60 Trichloroethene	130	5.092	5.092	(1.040)	341517	50.0000	55
	61 Isopropyl Acetate	43	4.679	4.679	(0.956)	251899	100.000	110
	62 N-Butanol	56	5.476	5.476	(1.119)	127722	500.000	560
	63 Dibromomethane	93	5.535	5.535	(1.131)	176911	50.0000	52
	64 1,2-Dichloropropane	63	5.634	5.634	(1.151)	369198	50.0000	53
	65 Bromodichloromethane	83	5.712	5.712	(1.167)	334439	50.0000	51
	66 Methyl Methacrylate	69	5.889	5.889	(1.203)	321433	100.000	100
	67 1,4-Dioxane	58	5.929	5.929	(1.211)	28926	500.000	550
	68 N-Propyl Acetate	43	6.303	6.303	(1.287)	70261	100.000	130
	69 2-Chloroethylvinylether	63	6.303	6.303	(1.287)	101597	50.0000	60
	70 cis-1,3-Dichloropropene	75	6.352	6.352	(1.297)	512696	50.0000	51
	71 Chloroacetonitrile	48	6.686	6.686	(1.366)	151771	500.000	530
	72 2-Nitropropane	41	6.765	6.765	(1.382)	193348	100.000	99
	73 trans-1,3-Dichloropropene	75	6.972	6.972	(1.424)	475140	50.0000	52
	74 1,1,2-Trichloroethane	97	7.119	7.119	(1.454)	234684	50.0000	53
*	75 Chlorobenzene-d5	117	7.956	7.956	(1.000)	442233	25.0000	
	76 Toluene	91	6.578	6.578	(0.827)	967975	50.0000	51
$	77 Toluene-d8	98	6.529	6.529	(0.821)	310214	25.0000	20
	78 1,1-Dichloro-2-propanone	43	6.795	6.795	(0.854)	1074275	250.000	230

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	ON-COL (ug/L)
79	4-Methyl-2-Pentanone	43	6.932	6.932	(0.871)	432419	50.0000	53
80	Tetrachloroethene	164	6.952	6.952	(0.874)	193605	50.0000	52
81	Ethyl Methacrylate	69	7.149	7.149	(0.899)	525098	50.0000	52
82	Dibromochloromethane	129	7.287	7.287	(0.916)	350036	50.0000	49
83	1,3-Dichloropropane	76	7.365	7.365	(0.926)	510476	50.0000	50
84	1,2-Dibromoethane	107	7.483	7.483	(0.941)	303604	50.0000	51
85	n-Butyl Acetate	56	7.641	7.641	(0.960)	278075	50.0000	54
86	2-Hexanone	43	7.710	7.710	(0.969)	331164	50.0000	54
87	1-Chlorohexane	91	7.965	7.965	(1.001)	349538	50.0000	55
88	Chlorobenzene	112	7.975	7.975	(1.002)	801005	50.0000	52
89	1,1,1,2-Tetrachloroethane	131	8.034	8.034	(1.010)	279633	50.0000	50
90	Ethylbenzene	106	8.005	8.005	(1.006)	354859	50.0000	52
91	Xylene (total)mp	106	8.143	8.143	(1.023)	890629	100.000	100
92	Xylene (total)o	106	8.516	8.516	(1.070)	436572	50.0000	54
93	Styrene	104	8.566	8.566	(1.077)	718314	50.0000	52
94	Bromoform	173	8.585	8.585	(1.079)	188908	50.0000	48
* 95	1,4-Dichlorobenzene-d4	152	10.012	10.012	(1.000)	160938	25.0000	
96	Isopropylbenzene	105	8.792	8.792	(0.878)	1030749	50.0000	52
97	Bromobenzene	156	9.126	9.126	(0.912)	257079	50.0000	51
98	1,1,2,2-Tetrachloroethane	83	9.225	9.225	(0.921)	339943	50.0000	50
99	4-Ethyltoluene	105	9.254	9.254	(0.924)	1018773	50.0000	52
100	1,2,3-Trichloropropane	110	9.323	9.323	(0.931)	114757	50.0000	53
101	trans-1,4-Dichloro-2-Butene	53	9.363	9.363	(0.935)	208259	100.000	93
102	n-Propylbenzene	91	9.156	9.156	(0.915)	1105483	50.0000	52
103	2-Chlorotoluene	91	9.294	9.294	(0.928)	790446	50.0000	54
104	4-Chlorotoluene	91	9.431	9.431	(0.942)	712713	50.0000	52
105	1,3,5-Trimethylbenzene	105	9.333	9.333	(0.932)	829002	50.0000	52
106	tert-Butylbenzene	119	9.609	9.609	(0.960)	792635	50.0000	52
107	1,2,4-Trimethylbenzene	105	9.668	9.668	(0.966)	840484	50.0000	53
108	sec-Butylbenzene	105	9.766	9.766	(0.975)	862855	50.0000	51
109	4-Isopropyltoluene	119	9.894	9.894	(0.988)	929221	50.0000	52
110	1,3-Dichlorobenzene	146	9.953	9.953	(0.994)	455715	50.0000	52
111	1,4-Dichlorobenzene	146	10.032	10.032	(1.002)	455886	50.0000	51
112	1,2-Dichlorobenzene	146	10.386	10.386	(1.037)	445615	50.0000	52
113	Benzyl Chloride	126	10.238	10.238	(1.023)	135660	50.0000	51
114	1,4-Diethylbenzene	119	10.209	10.209	(2.085)	525346	50.0000	55
115	n-Butylbenzene	91	10.258	10.258	(1.025)	1103294	50.0000	51
118	1,2,4,5-Tetramethylbenzene	119	10.917	10.917	(2.230)	814972	50.0000	56
119	1,2-Dibromo-3-chloropropane	75	11.084	11.084	(1.107)	71043	50.0000	52
120	Nitrobenzene	77	11.566	11.566	(1.155)	134099	500.000	440
121	1,2,4-Trichlorobenzene	180	11.685	11.685	(1.167)	260949	50.0000	56
122	Hexachlorobutadiene	225	11.675	11.675	(1.166)	77928	50.0000	49
123	Naphthalene	128	11.970	11.970	(1.196)	896599	50.0000	63
124	1,2,3-Trichlorobenzene	180	12.137	12.137	(1.212)	234065	50.0000	55
$ 125	Bromofluorobenzene	95	9.038	9.038	(0.903)	140229	25.0000	23
M 126	1,2-Dichloroethene (total)	100				511778	100.000	100
M 127	Xylene (total)	100				1327201	150.000	160

Data File: L9310.D

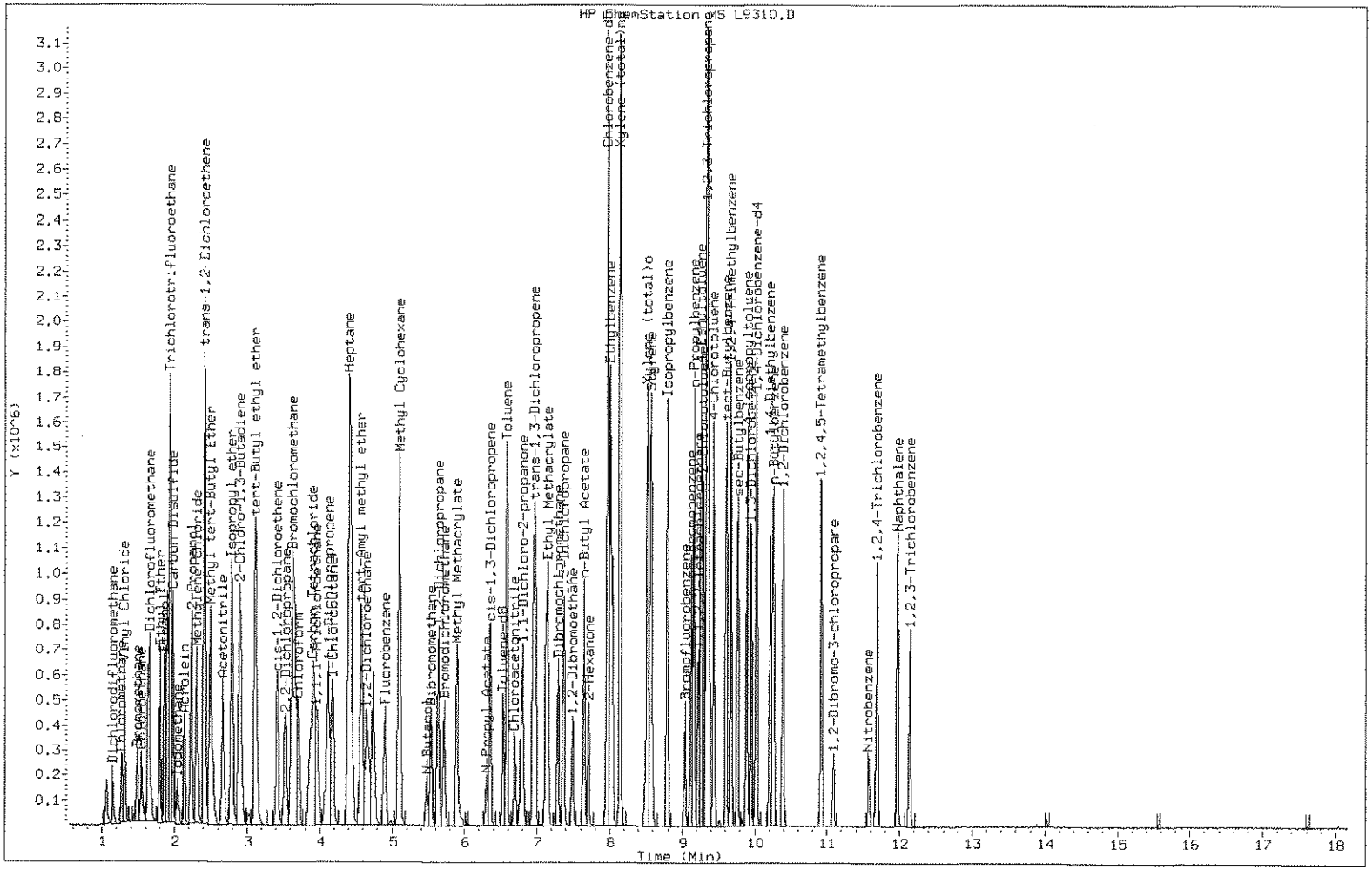
Date: 31-JUL-2007 20:21

Client ID:

Instrument: ms1.i

Sample Info: CCVIS

Operator: D. GAYDA



VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Instrument ID: MSL Calibration Date: 08/01/2007 Time: 09:47

Lab File ID: L9341.D Init. Calib. Date(s): 07/26/2007 07/26/2007

Lab Sample ID: CCVIS 220-8321/1 Init. Calib. Time(s): 13:32 15:11

GC Column: RTX-VMS ID: 0.18 (mm) Heated Purge: (Y/N) N Conc. Units: ug/L

Analyte	Curve	Ave RRF	RRF	Min RRF	Calculated Amount	Ccal Amount	% D	Max % D
1,1,1,2-Tetrachloroethane	Ave	0.3128	0.3125		50.0	50.0	0.1	
1,1,1-Trichloroethane	Ave	0.3840	0.3931		51.2	50.0	2.4	
1,1,2,2-Tetrachloroethane	Ave	1.0638	1.0591	0.3000	49.8	50.0	-0.5	
1,1,2-Trichloro-1,2,2-trifluoroethane	Ave	0.2475	0.2469		49.9	50.0	-0.3	
1,1,2-Trichloroethane	Ave	0.2502	0.2595		51.8	50.0	3.7	
1,1-Dichloro-1-fluoroethane	Ave	0.4422	0.4340		49.1	50.0	-1.8	
1,1-Dichloroacetone	Ave	0.2611	0.2539		243	250	-2.7	
1,1-Dichloroethane	Ave	0.6829	0.6942	0.1000	50.8	50.0	1.6	
1,1-Dichloroethene	Ave	0.2131	0.2038		47.8	50.0	-4.4	20.0
1,1-Dichloropropene	Ave	0.4776	0.4872		51.0	50.0	2.0	
1,2,3-Trichlorobenzene	Ave	0.6643	0.6695		50.4	50.0	0.8	
1,2,3-Trichloropropane	Ave	0.3371	0.3278		48.6	50.0	-2.8	
1,2,4,5-Tetramethylbenzene	Ave	0.8285	0.8525		51.4	50.0	2.9	
1,2,4-Trichlorobenzene	Ave	0.7280	0.7546		51.8	50.0	3.6	
1,2,4-Trimethylbenzene	Ave	2.4805	2.5050		50.5	50.0	1	
1,2-Dibromo-3-Chloropropane	Ave	0.2124	0.2080		49.0	50.0	-2.1	
1,2-Dichloro-1,1,2-trifluoroethane	Ave	0.6540			49.7			
1,2-Dichlorobenzene	Ave	1.3400	1.3208		49.3	50.0	-1.4	
1,2-Dichloroethane	Ave	0.4746	0.4800		50.6	50.0	1.1	
1,2-Dichloroethane-d4 (Surr)	Ave	0.3932	0.2888		18.4	25.0	-26.5	
1,2-Dichloroethene, Total	Ave	0.2306	0.2780		100	100	20.5	
1,2-Dichloropropane	Ave	0.3960	0.4067		51.4	50.0	2.7	20.0
1,3,5-Trimethylbenzene	Ave	2.4694	2.5039		50.7	50.0	1.4	
1,3-Dichlorobenzene	Ave	1.3616	1.3479		49.5	50.0	-1.0	
1,3-Dichloropropane	Ave	0.5774	0.5617		48.6	50.0	-2.7	
1,4-Dichlorobenzene	Ave	1.3897	1.3784		49.6	50.0	-0.8	
1,4-Dioxane	Ave	0.0030	0.0036		597	500	19.4	
1-Bromopropane	Ave	0.5150	0.5337		51.8	50.0	3.6	
1-Chlorobutane	Ave	0.6854	0.7223		52.7	50.0	5.4	
1-Chlorohexane	Ave	0.3605	0.3655		50.7	50.0	1.4	
2,2-Dichloropropane	Ave	0.4711	0.4964		52.7	50.0	5.4	
2,4,4-Trimethyl-1-pentene	Ave				0.0			
2-Chloro-1,3-butadiene	Ave	0.1935	0.1951		50.4	50.0	0.8	
2-Chloroethyl vinyl ether	Ave	0.0965	0.0824		42.7	50.0	-14.6	
2-Chlorotoluene	Ave	2.2771	2.4482		53.8	50.0	7.5	
2-Hexanone	Ave	0.3432	0.3505		51.1	50.0	2.1	
2-Methyl-2-propanol	Ave	0.0579	0.0591		255	250	2.1	
2-Nitropropane	Ave	0.1112	0.1071		96.3	100	-3.7	
3-Chloro-1-propene	Ave	0.5499	0.5375		48.9	50.0	-2.3	
4-Bromofluorobenzene	Ave	0.9379	0.8664		23.1	25.0	-7.6	
4-Chlorotoluene	Ave	2.1467	2.1495		50.1	50.0	0.1	
4-Ethyltoluene	Ave	3.0136	3.0259		50.2	50.0	0.4	

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Instrument ID: MSL Calibration Date: 08/01/2007 Time: 09:47

Lab File ID: L9341.D Init. Calib. Date(s): 07/26/2007 07/26/2007

Lab Sample ID: CCVIS 220-8321/1 Init. Calib. Time(s): 13:32 15:11

GC Column: RTX-VMS ID: 0.18 (mm) Heated Purge: (Y/N) N Conc. Units: ug/L

Analyte	Curve	Ave RRF	RRF	Min RRF	Calculated Amount	Ccal Amount	% D	Max % D
4-Isopropyltoluene	Ave	2.7945	2.7304		48.9	50.0	-2.3	
Acetone	Ave	0.1585	0.1866		58.9	50.0	17.7	
Acetonitrile	Ave	0.0661	0.0681		515	500	3.0	
Acrolein	Ave	0.0580	0.0745		321	250	28.4	
Acrylonitrile	Ave	0.1834	0.1823		99.4	100	-0.6	
Benzene	Ave	1.2001	1.2017		50.1	50.0	0.1	
Benzyl chloride	Ave	0.4143	0.4218		50.9	50.0	1.8	
Bromobenzene	Ave	0.7785	0.7671		49.3	50.0	-1.5	
Bromoform	Ave	0.2207	0.2120	0.1000	48.0	50.0	-3.9	
Bromomethane	Ave	0.1293	0.1143		44.2	50.0	-11.6	
Carbon disulfide	Ave	1.0400	0.8521		41.0	50.0	-18.1	
Carbon tetrachloride	Ave	0.4301	0.4725		54.9	50.0	9.9	
Chloroacetonitrile	Ave	0.0162	0.0156		481	500	-3.8	
Chlorobenzene	Ave	0.8648	0.8892	0.3000	51.4	50.0	2.8	
Chlorobromomethane	Ave	0.2047	0.2078		50.7	50.0	1.5	
Chlorodibromomethane	Ave	0.4000	0.3771		47.1	50.0	-5.7	
Chloroethane	Ave	0.1962	0.2438		62.1	50.0	24.3	
Chloroform	Ave	0.5224	0.5362		51.3	50.0	2.6	20.0
Chloromethane	Ave	0.3363	0.3244	0.1000	48.2	50.0	-3.5	
cis-1,2-Dichloroethene	Ave	0.2859	0.2950		51.6	50.0	3.2	
cis-1,3-Dichloropropene	Ave	0.5664	0.5609		49.5	50.0	-1	
Cyclohexane	Ave	0.3652	0.3675		50.3	50.0	0.6	
Dibromofluoromethane	Ave	0.3225	0.2501		19.4	25.0	-22.4	
Dibromomethane	Ave	0.1924	0.1940		50.4	50.0	0.9	
Dichlorobromomethane	Ave	0.3741	0.3744		50.0	50.0	0.1	
Dichlorodifluoromethane	Ave	0.1920	0.1822		47.5	50.0	-5.1	
Dichlorofluoromethane	Ave	0.6540	0.6496		49.7	50.0	-0.7	
Ethanol	Ave	0.0178	0.0189		532	500	6.5	
Ethyl acetate	Ave	0.0280	0.0272		97.4	100	-2.6	
Ethyl ether	Ave	0.2255	0.2257		50.0	50.0	0.1	
Ethyl methacrylate	Ave	0.5666	0.5687		50.2	50.0	0.4	
Ethylbenzene	Ave	0.3850	0.4033		52.4	50.0	4.8	20.0
Ethylene Dibromide	Ave	0.3366	0.3383		50.3	50.0	0.5	
Hexachlorobutadiene	Ave	0.2481	0.2362		47.6	50.0	-4.8	
Hexachloroethane	Ave				0.0			
Iodomethane	Ave	0.2764	0.2585		46.8	50.0	-6.5	
Isobutanol	Ave	0.0162	0.0154		477	500	-4.6	
Isopropyl acetate	Ave	0.1342	0.1303		97.1	100	-2.9	
Isopropyl alcohol	Ave	0.0574	0.0626		54.5	50.0	9.0	
Isopropyl ether	Ave	1.2074	1.2414		51.4	50.0	2.8	
Isopropylbenzene	Ave	3.0830	3.0984		50.2	50.0	0.5	
m-Xylene & p-Xylene	Ave	0.4812	0.4979		103	100	3.5	
Methacrylonitrile	Ave	0.4469	0.4642		51.9	50.0	3.9	

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Instrument ID: MSL Calibration Date: 08/01/2007 Time: 09:47

Lab File ID: L9341.D Init. Calib. Date(s): 07/26/2007 07/26/2007

Lab Sample ID: CCVIS 220-8321/1 Init. Calib. Time(s): 13:32 15:11

GC Column: RTX-VMS ID: 0.18 (mm) Heated Purge: (Y/N) N Conc. Units: ug/L

Analyte	Curve	Ave RRF	RRF	Min RRF	Calculated Amount	Ccal Amount	% D	Max % D
Methyl acetate	Ave	1.8684	1.7774		47.6	50.0	-4.9	
Methyl acrylate	Ave	0.4003	0.3971		49.6	50.0	-0.8	
Methyl Ethyl Ketone	Ave	0.2301	0.2476		53.8	50.0	7.6	
methyl isobutyl ketone	Ave	0.4602	0.4537		49.3	50.0	-1.4	
Methyl methacrylate	Ave	0.1736	0.3471		100.0	50.0	99.9	
Methyl tert-butyl ether	Ave	1.0379	1.0341		49.8	50.0	-0.4	
Methylcyclohexane	Ave	0.3498	0.3397		48.6	50.0	-2.9	
Methylene Chloride	Ave	0.2929	0.2870		49.0	50.0	-2.0	
n-Butanol	Ave	0.0130	0.0142		547	500	9.4	
n-Butyl acetate	Ave	0.2901	0.2985		51.4	50.0	2.9	
n-Butylbenzene	Ave	3.3672	3.6452		54.1	50.0	8.3	
n-Heptane	Ave	0.3090	0.2987		48.3	50.0	-3.3	
n-Propyl acetate	Ave	0.0316	0.0288		91.2	100	-8.8	
N-Propylbenzene	Ave	3.3039	3.2975		49.9	50.0	-0.2	
Naphthalene	Ave	2.2068	2.4163		54.7	50.0	9.5	
Nitrobenzene	Ave	0.0473	0.0367		388	500	-22.3	
o-Xylene	Ave	0.4609	0.4725		51.3	50.0	2.5	
p-Diethylbenzene	Ave	0.5445	0.5544		50.9	50.0	1.8	
Pentachloroethane	Ave				0.0			
Propionitrile	Ave	0.0634	0.0633		499	500	-0.3	
sec-Butylbenzene	Ave	2.6102	2.5595		49.0	50.0	-1.9	
Styrene	Ave	0.7721	0.7829		50.7	50.0	1.4	
Tert-amyl methyl ether	Ave	1.0524	1.0534		50.0	50.0	0.1	
Tert-butyl ethyl ether	Ave	1.2132	1.2529		51.6	50.0	3.3	
tert-Butyl Formate	Ave	0.3499	0.3562		50.9	50.0	1.8	
tert-Butylbenzene	Ave	2.3884	2.3244		48.7	50.0	-2.7	
Tetrachloroethene	Ave	0.2100	0.2128		50.7	50.0	1.3	
Tetrahydrofuran	Ave	0.1355	0.1320		97.5	100	-2.5	
Toluene	Ave	1.0645	1.0747		50.5	50.0	1	20.0
Toluene-d8 (Surr)	Ave	0.8889	0.7195		20.2	25.0	-19.1	
trans-1,2-Dichloroethene	Ave	0.2676	0.2611		48.8	50.0	-2.4	
trans-1,3-Dichloropropene	Ave	0.5141	0.5174		50.3	50.0	0.7	
trans-1,4-Dichloro-2-butene	Ave	0.3461	0.2887		83.4	100	-16.6	
Trichloroethene	Ave	0.3531	0.3486		49.4	50.0	-1.3	
Trichlorofluoromethane	Ave	0.3584	0.3736		52.1	50.0	4.2	
Vinyl acetate	Ave	0.9114	0.9212		50.5	50.0	1.1	
Vinyl chloride	Ave	0.3701	0.3865		52.2	50.0	4.4	20.0
Xylenes, Total	Ave	0.3953	0.4894		155	150	23.8	

Curve Types:
Ave = Average
Lin = Linear
Quad = Quadratic

STL-CT

Volatile Report SW-846 Method 8260B

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079341.b\L9341.D
Lab Smp Id: CCVIS Client Smp ID: CCVIS
Inj Date : 01-AUG-2007 09:47 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : CCVIS
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079341.b\L8260BNW.m
Meth Date : 02-Aug-2007 09:12 target Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 31 Continuing Calibration Sample
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONMSLNT

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	ON-COL (ug/L)
* 1	Fluorobenzene	96	4.908	4.908	(1.000)	451284	25.0000	
2	Dichlorodifluoromethane	85	1.149	1.149	(0.234)	164433	50.0000	47
3	Chloromethane	50	1.267	1.267	(0.258)	292749	50.0000	48
4	Vinyl Chloride	62	1.307	1.307	(0.266)	348879	50.0000	52
5	Bromomethane	94	1.474	1.474	(0.300)	103170	50.0000	44
6	Chloroethane	64	1.543	1.543	(0.314)	220023	50.0000	62
7	Trichlorofluoromethane	101	1.631	1.631	(0.332)	337170	50.0000	52
8	Dichlorofluoromethane	67	1.651	1.651	(0.336)	586313	50.0000	50
9	Ethyl Ether	45	1.799	1.799	(0.367)	203686	50.0000	50
10	Ethanol	45	1.858	1.858	(0.379)	170740	500.000	530
11	Freon 141	81	1.858	1.858	(0.379)	391747	50.0000	49
12	Freon 123a	67	1.651	1.651	(0.336)	586313	50.0000	50
13	Trichlorotrifluoroethane	101	1.946	1.946	(0.397)	222811	50.0000	50
14	1,1-Dichloroethene	96	1.936	1.936	(0.395)	183919	50.0000	48
15	Carbon Disulfide	76	1.976	1.976	(0.403)	769109	50.0000	41
16	Iodomethane	142	2.035	2.035	(0.415)	233315	50.0000	47
17	Acrolein	56	2.133	2.133	(0.435)	336214	250.000	320
18	2-Propanol	45	2.222	2.222	(0.453)	56457	50.0000	54
19	3-Chloro-1-Propene	41	2.232	2.232	(0.455)	485152	50.0000	49
20	Methylene Chloride	84	2.300	2.300	(0.469)	259059	50.0000	49
21	Acetone	43	2.330	2.330	(0.475)	168437	50.0000	59
22	trans-1,2-Dichloroethene	96	2.428	2.428	(0.495)	235611	50.0000	49

			QUANT SIG					AMOUNTS	
		Compounds	MASS	RT	EXP RT	REL RT	RESPONSE	CAL-AMT (ug/L)	ON-COL (ug/L)
	23	Methyl Acetate	43	2.409	2.409	(0.491)	1604236	50.0000	48
	24	Methyl tert-Butyl Ether	73	2.497	2.497	(0.509)	933314	50.0000	50
	25	tert-Butyl alcohol	59	2.537	2.537	(0.517)	266718	250.000	260
	26	Acetonitrile	41	2.665	2.665	(0.543)	614521	500.000	510
	27	Isopropyl ether	45	2.783	2.783	(0.567)	1120446	50.0000	51
	28	tert-Butyl ethyl ether	59	3.117	3.117	(0.635)	1130831	50.0000	52
	29	2-Chloro-1,3-Butadiene	88	2.891	2.891	(0.589)	176099	50.0000	50
	30	Acrylonitrile	53	2.920	2.920	(0.595)	329096	100.000	99
	31	1,1-Dichloroethane	63	2.901	2.901	(0.591)	626567	50.0000	51
	32	Vinyl Acetate	43	3.107	3.107	(0.633)	831481	50.0000	50
	33	cis-1,2-Dichloroethene	96	3.422	3.422	(0.697)	266265	50.0000	52
	34	2,2-Dichloropropane	77	3.530	3.530	(0.719)	448046	50.0000	53
	35	Bromochloromethane	128	3.629	3.629	(0.739)	187504	50.0000	51
	36	1-Bromopropane	43	3.619	3.619	(0.737)	481733	50.0000	52
	37	Cyclohexane	84	3.658	3.658	(0.745)	331698	50.0000	50
	38	Chloroform	83	3.707	3.707	(0.755)	483971	50.0000	51
	39	Ethyl Acetate	43	3.855	3.855	(0.786)	49132	100.000	97
	40	Methyl Acrylate	55	3.865	3.865	(0.788)	358386	50.0000	50
$	41	Dibromofluoromethane	111	3.924	3.924	(0.800)	112880	25.0000	19
	42	Tetrahydrofuran	42	3.924	3.924	(0.800)	238327	100.000	97
	43	Carbon Tetrachloride	117	3.904	3.904	(0.796)	426471	50.0000	55
	44	1,1,1-Trichloroethane	97	3.973	3.973	(0.810)	354754	50.0000	51
	45	2-Butanone	43	4.071	4.071	(0.830)	223443	50.0000	54
	46	1,1-Dichloropropene	75	4.131	4.131	(0.842)	439697	50.0000	51
	47	tert-Amyl methyl ether	73	4.573	4.573	(0.932)	950752	50.0000	50
	48	tert-Butyl formate	57	3.117	3.117	(0.635)	321466	50.0000	51
	49	1-Chlorobutane	56	4.180	4.180	(0.852)	651964	50.0000	53
	50	Heptane	43	4.406	4.406	(0.898)	269617	50.0000	48
	51	Propionitrile	54	4.406	4.406	(0.898)	570906	500.000	500
	52	Benzene	78	4.426	4.426	(0.902)	1084581	50.0000	50
	53	2-Methyl-2-Propenenitrile	41	4.436	4.436	(0.904)	418989	50.0000	52
	54	Isobutyl alcohol	42	4.682	4.682	(0.954)	139124	500.000	480
$	55	1,2-Dichloroethane-d4	65	4.563	4.563	(0.930)	130340	25.0000	18
	56	1,2-Dichloroethane	62	4.652	4.652	(0.948)	433194	50.0000	50
	58	2,4,4-Trimethyl 2-Pentene	55	1.809	1.809	(0.369)	3113	50.0000	
	59	Methyl Cyclohexane	83	5.095	5.095	(1.038)	306641	50.0000	48
	60	Trichloroethene	130	5.105	5.105	(1.040)	314660	50.0000	49
	61	Isopropyl Acetate	43	4.682	4.682	(0.954)	235168	100.000	97
	62	N-Butanol	56	5.488	5.488	(1.118)	128515	500.000	550
	63	Dibromomethane	93	5.538	5.538	(1.128)	175138	50.0000	50
	64	1,2-Dichloropropane	63	5.636	5.636	(1.148)	367092	50.0000	51
	65	Bromodichloromethane	83	5.724	5.724	(1.166)	337926	50.0000	50
	66	Methyl Methacrylate	69	5.902	5.902	(1.202)	313244	100.000	100
	67	1,4-Dioxane	58	5.941	5.941	(1.210)	32302	500.000	600
	68	N-Propyl Acetate	43	6.305	6.305	(1.285)	52060	100.000	91
	69	2-Chloroethylvinylether	63	6.305	6.305	(1.285)	74393	50.0000	43
	70	cis-1,3-Dichloropropene	75	6.354	6.354	(1.295)	506205	50.0000	50
	71	Chloroacetonitrile	48	6.698	6.698	(1.365)	140643	500.000	480
	72	2-Nitropropane	41	6.777	6.777	(1.381)	193237	100.000	96
	73	trans-1,3-Dichloropropene	75	6.984	6.984	(1.423)	467023	50.0000	50
	74	1,1,2-Trichloroethane	97	7.122	7.122	(1.451)	234171	50.0000	52
*	75	Chlorobenzene-d5	117	7.968	7.968	(1.000)	444659	25.0000	
	76	Toluene	91	6.580	6.580	(0.826)	955775	50.0000	50
$	77	Toluene-d8	98	6.541	6.541	(0.821)	319943	25.0000	20

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	AMOUNTS ON-COL (ug/L)
78	1,1-Dichloro-2-propanone	43	6.797	6.797	(0.853)	1128926	250.000	240
79	4-Methyl-2-Pentanone	43	6.944	6.944	(0.872)	403461	50.0000	49
80	Tetrachloroethene	164	6.964	6.964	(0.874)	189251	50.0000	51
81	Ethyl Methacrylate	69	7.151	7.151	(0.898)	505707	50.0000	50
82	Dibromochloromethane	129	7.289	7.289	(0.915)	335349	50.0000	47
83	1,3-Dichloropropane	76	7.368	7.368	(0.925)	499527	50.0000	49
84	1,2-Dibromoethane	107	7.495	7.495	(0.941)	300850	50.0000	50
85	n-Butyl Acetate	56	7.653	7.653	(0.960)	265432	50.0000	51
86	2-Hexanone	43	7.712	7.712	(0.968)	311693	50.0000	51
87	1-Chlorohexane	91	7.968	7.968	(1.000)	325084	50.0000	51
88	Chlorobenzene	112	7.978	7.978	(1.001)	790765	50.0000	51
89	1,1,1,2-Tetrachloroethane	131	8.037	8.037	(1.009)	277869	50.0000	50
90	Ethylbenzene	106	8.007	8.007	(1.005)	358693	50.0000	52
91	Xylene (total)mp	106	8.145	8.145	(1.022)	885500	100.000	100
92	Xylene (total)o	106	8.519	8.519	(1.069)	420241	50.0000	51
93	Styrene	104	8.568	8.568	(1.075)	696271	50.0000	51
94	Bromoform	173	8.588	8.588	(1.078)	188563	50.0000	48
* 95	1,4-Dichlorobenzene-d4	152	10.014	10.014	(1.000)	164048	25.0000	
96	Isopropylbenzene	105	8.804	8.804	(0.879)	1016570	50.0000	50
97	Bromobenzene	156	9.129	9.129	(0.912)	251668	50.0000	49
98	1,1,2,2-Tetrachloroethane	83	9.227	9.227	(0.921)	347471	50.0000	50
99	4-Ethyltoluene	105	9.257	9.257	(0.924)	992791	50.0000	50
100	1,2,3-Trichloropropane	110	9.335	9.335	(0.932)	107539	50.0000	49
101	trans-1,4-Dichloro-2-Butene	53	9.375	9.375	(0.936)	189442	100.000	83
102	n-Propylbenzene	91	9.158	9.158	(0.915)	1081910	50.0000	50
103	2-Chlorotoluene	91	9.296	9.296	(0.928)	803250	50.0000	54
104	4-Chlorotoluene	91	9.434	9.434	(0.942)	705255	50.0000	50
105	1,3,5-Trimethylbenzene	105	9.335	9.335	(0.932)	821534	50.0000	51
106	tert-Butylbenzene	119	9.611	9.611	(0.960)	762614	50.0000	49
107	1,2,4-Trimethylbenzene	105	9.670	9.670	(0.966)	821886	50.0000	50
108	sec-Butylbenzene	105	9.768	9.768	(0.975)	839764	50.0000	49
109	4-Isopropyltoluene	119	9.896	9.896	(0.988)	895824	50.0000	49
110	1,3-Dichlorobenzene	146	9.955	9.955	(0.994)	442238	50.0000	49
111	1,4-Dichlorobenzene	146	10.034	10.034	(1.002)	452234	50.0000	50
112	1,2-Dichlorobenzene	146	10.388	10.388	(1.037)	433353	50.0000	49
113	Benzyl Chloride	126	10.241	10.241	(1.023)	138390	50.0000	51
114	1,4-Diethylbenzene	119	10.211	10.211	(2.080)	500395	50.0000	51
115	n-Butylbenzene	91	10.260	10.260	(1.025)	1195984	50.0000	54
118	1,2,4,5-Tetramethylbenzene	119	10.919	10.919	(2.225)	769413	50.0000	51
119	1,2-Dibromo-3-chloropropane	75	11.087	11.087	(1.107)	68241	50.0000	49
120	Nitrobenzene	77	11.579	11.579	(1.156)	120494	500.000	390
121	1,2,4-Trichlorobenzene	180	11.697	11.697	(1.168)	247569	50.0000	52
122	Hexachlorobutadiene	225	11.677	11.677	(1.166)	77501	50.0000	48
123	Naphthalene	128	11.972	11.972	(1.196)	792770	50.0000	55
124	1,2,3-Trichlorobenzene	180	12.139	12.139	(1.212)	219654	50.0000	50
$ 125	Bromofluorobenzene	95	9.040	9.040	(0.903)	142127	25.0000	23
M 126	1,2-Dichloroethene (total)	100				501876	100.000	100
M 127	Xylene (total)	100				1305741	150.000	150

Data File: L9341.D

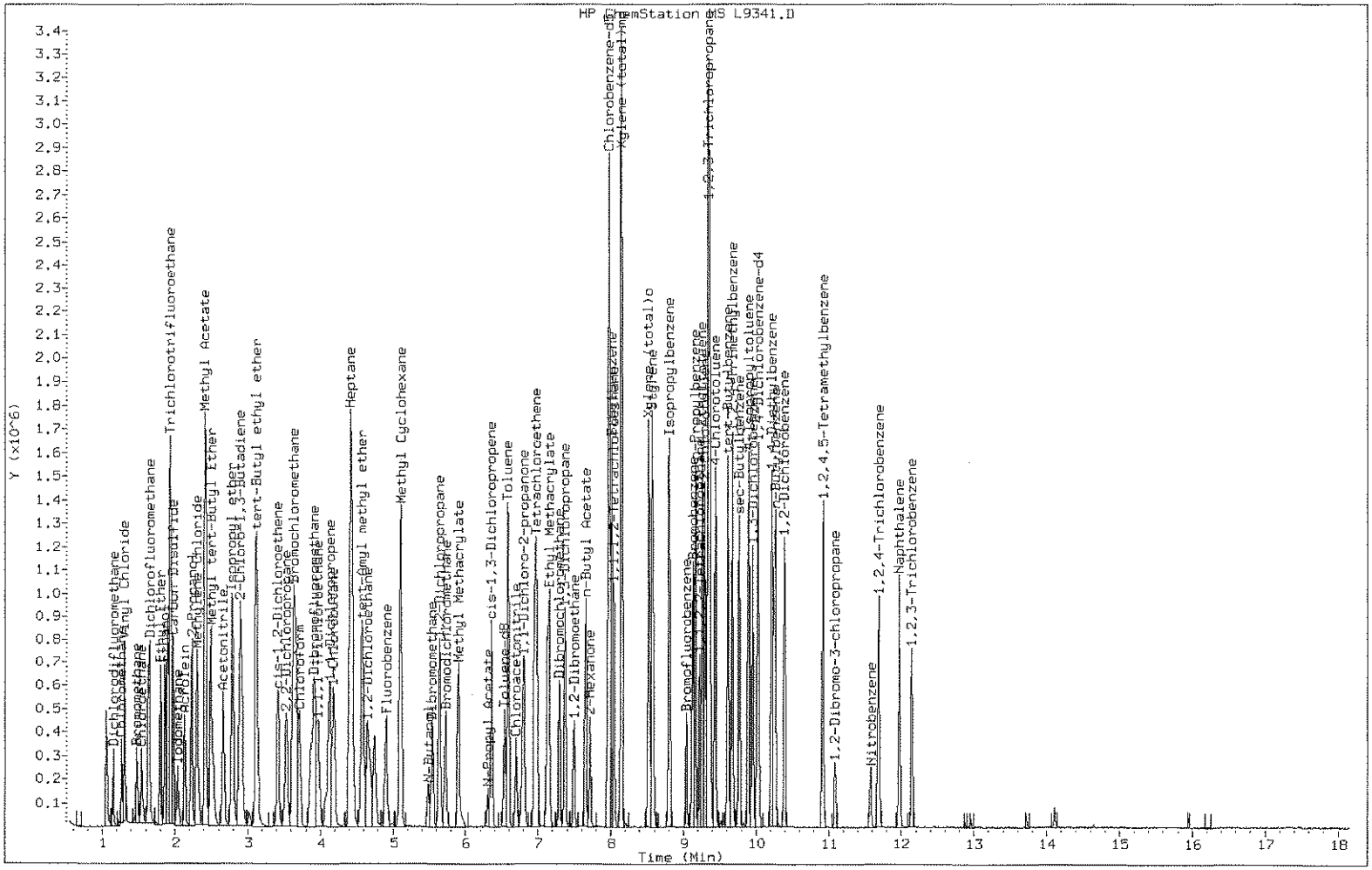
Date: 01-AUG-2007 09:47

Client ID: CCVIS

Instrument: msl.i

Sample Info: CCVIS

Operator: D. HUMBERT



08/09/2007

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Instrument ID: MSL Calibration Date: 08/02/2007 Time: 09:18

Lab File ID: L9371.D Init. Calib. Date(s): 07/26/2007 07/26/2007

Lab Sample ID: CCVIS 220-8356/1 Init. Calib. Time(s): 13:32 15:11

GC Column: RTX-VMS ID: 0.18 (mm) Heated Purge: (Y/N) N Conc. Units: ug/L

Analyte	Curve	Ave RRF	RRF	Min RRF	Calculated Amount	Ccal Amount	% D	Max % D
1,1,1,2-Tetrachloroethane	Ave	0.3128	0.3049		48.7	50.0	-2.5	
1,1,1-Trichloroethane	Ave	0.3840	0.4136		53.9	50.0	7.7	
1,1,2,2-Tetrachloroethane	Ave	1.0638	1.1000	0.3000	51.7	50.0	3.4	
1,1,2-Trichloro-1,2,2-trifluoroethane	Ave	0.2475	0.2754		55.6	50.0	11.3	
1,1,2-Trichloroethane	Ave	0.2502	0.2668		53.3	50.0	6.6	
1,1-Dichloro-1-fluoroethane	Ave	0.4422	0.4654		52.6	50.0	5.2	
1,1-Dichloroacetone	Ave	0.2611	0.2529		242	250	-3.1	
1,1-Dichloroethane	Ave	0.6829	0.7360	0.1000	53.9	50.0	7.8	
1,1-Dichloroethene	Ave	0.2131	0.2315		54.3	50.0	8.6	20.0
1,1-Dichloropropene	Ave	0.4776	0.5055		52.9	50.0	5.8	
1,2,3-Trichlorobenzene	Ave	0.6643	0.7071		53.2	50.0	6.4	
1,2,3-Trichloropropane	Ave	0.3371	0.3409		50.6	50.0	1.1	
1,2,4,5-Tetramethylbenzene	Ave	0.8285	0.9070		54.7	50.0	9.5	
1,2,4-Trichlorobenzene	Ave	0.7280	0.7871		54.1	50.0	8.1	
1,2,4-Trimethylbenzene	Ave	2.4805	2.6177		52.8	50.0	5.5	
1,2-Dibromo-3-Chloropropane	Ave	0.2124	0.2118		49.8	50.0	-0.3	
1,2-Dichloro-1,1,2-trifluoroethane	Ave	0.6540			54.4			
1,2-Dichlorobenzene	Ave	1.3400	1.3780		51.4	50.0	2.8	
1,2-Dichloroethane	Ave	0.4746	0.5078		53.5	50.0	7.0	
1,2-Dichloroethane-d4 (Surr)	Ave	0.3932	0.2854		18.1	25.0	-27.4	
1,2-Dichloroethene, Total	Ave	0.2306	0.2985		108	100	29.4	
1,2-Dichloropropane	Ave	0.3960	0.4264		53.8	50.0	7.7	20.0
1,3,5-Trimethylbenzene	Ave	2.4694	2.6361		53.4	50.0	6.7	
1,3-Dichlorobenzene	Ave	1.3616	1.4330		52.6	50.0	5.2	
1,3-Dichloropropane	Ave	0.5774	0.5761		49.9	50.0	-0.2	
1,4-Dichlorobenzene	Ave	1.3897	1.4355		51.6	50.0	3.3	
1,4-Dioxane	Ave	0.0030	0.0033		545	500	9.1	
1-Bromopropane	Ave	0.5150	0.5573		54.1	50.0	8.2	
1-Chlorobutane	Ave	0.6854	0.7565		55.2	50.0	10.4	
1-Chlorohexane	Ave	0.3605	0.3531		49.0	50.0	-2.1	
2,2-Dichloropropane	Ave	0.4711	0.5186		55.0	50.0	10.1	
2,4,4-Trimethyl-1-pentene	Ave				0.0			
2-Chloro-1,3-butadiene	Ave	0.1935	0.2073		53.6	50.0	7.1	
2-Chloroethyl vinyl ether	Ave	0.0965	0.1470		76.2	50.0	52.4	
2-Chlorotoluene	Ave	2.2771	2.2325		49.0	50.0	-2.0	
2-Hexanone	Ave	0.3432	0.3641		53.0	50.0	6.1	
2-Methyl-2-propanol	Ave	0.0579	0.0569		246	250	-1.7	
2-Nitropropane	Ave	0.1112	0.1130		102	100	1.6	
3-Chloro-1-propene	Ave	0.5499	0.6042		54.9	50.0	9.9	
4-Bromofluorobenzene	Ave	0.9379	0.8817		23.5	25.0	-6.0	
4-Chlorotoluene	Ave	2.1467	2.2785		53.1	50.0	6.1	
4-Ethyltoluene	Ave	3.0136	3.1924		53.0	50.0	5.9	

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Instrument ID: MSL Calibration Date: 08/02/2007 Time: 09:18

Lab File ID: L9371.D Init. Calib. Date(s): 07/26/2007 07/26/2007

Lab Sample ID: CCVIS 220-8356/1 Init. Calib. Time(s): 13:32 15:11

GC Column: RTX-VMS ID: 0.18 (mm) Heated Purge: (Y/N) N Conc. Units: ug/L

Analyte	Curve	Ave RRF	RRF	Min RRF	Calculated Amount	Ccal Amount	% D	Max % D
4-Isopropyltoluene	Ave	2.7945	2.9041		52.0	50.0	3.9	
Acetone	Ave	0.1585	0.1939		61.1	50.0	22.3	
Acetonitrile	Ave	0.0661	0.0692		523	500	4.6	
Acrolein	Ave	0.0580	0.0677		292	250	16.7	
Acrylonitrile	Ave	0.1834	0.2031		111	100	10.7	
Benzene	Ave	1.2001	1.2830		53.5	50.0	6.9	
Benzyl chloride	Ave	0.4143	0.4812		58.1	50.0	16.1	
Bromobenzene	Ave	0.7785	0.8117		52.1	50.0	4.3	
Bromoform	Ave	0.2207	0.2205	0.1000	50.0	50.0	0.1	
Bromomethane	Ave	0.1293	0.1371		53.0	50.0	6.0	
Carbon disulfide	Ave	1.0400	1.1001		52.9	50.0	5.8	
Carbon tetrachloride	Ave	0.4301	0.5017		58.3	50.0	16.6	
Chloroacetonitrile	Ave	0.0162	0.0171		527	500	5.4	
Chlorobenzene	Ave	0.8648	0.8952	0.3000	51.8	50.0	3.5	
Chlorobromomethane	Ave	0.2047	0.2189		53.5	50.0	6.9	
Chlorodibromomethane	Ave	0.4000	0.3909		48.9	50.0	-2.3	
Chloroethane	Ave	0.1962	0.2427		61.9	50.0	23.7	
Chloroform	Ave	0.5224	0.5508		52.7	50.0	5.4	20.0
Chloromethane	Ave	0.3363	0.3329	0.1000	49.5	50.0	-1	
cis-1,2-Dichloroethene	Ave	0.2859	0.3041		53.2	50.0	6.3	
cis-1,3-Dichloropropene	Ave	0.5664	0.6044		53.4	50.0	6.7	
Cyclohexane	Ave	0.3652	0.3987		54.6	50.0	9.2	
Dibromofluoromethane	Ave	0.3225	0.2476		19.2	25.0	-23.2	
Dibromomethane	Ave	0.1924	0.2084		54.2	50.0	8.3	
Dichlorobromomethane	Ave	0.3741	0.3938		52.6	50.0	5.3	
Dichlorodifluoromethane	Ave	0.1920	0.1780		46.4	50.0	-7.3	
Dichlorofluoromethane	Ave	0.6540	0.7121		54.4	50.0	8.9	
Ethanol	Ave	0.0178	0.0195		549	500	9.9	
Ethyl acetate	Ave	0.0280	0.0290		104	100	3.8	
Ethyl ether	Ave	0.2255	0.2477		54.9	50.0	9.8	
Ethyl methacrylate	Ave	0.5666	0.5738		50.6	50.0	1.3	
Ethylbenzene	Ave	0.3850	0.4125		53.6	50.0	7.2	20.0
Ethylene Dibromide	Ave	0.3366	0.3483		51.7	50.0	3.5	
Hexachlorobutadiene	Ave	0.2481	0.2496		50.3	50.0	0.6	
Hexachloroethane	Ave				0.0			
Iodomethane	Ave	0.2764	0.2140		38.7	50.0	-22.6	
Isobutanol	Ave	0.0162	0.0163		506	500	1.1	
Isopropyl acetate	Ave	0.1342	0.1328		99.0	100	-1.0	
Isopropyl alcohol	Ave	0.0574	0.0596		51.9	50.0	3.9	
Isopropyl ether	Ave	1.2074	1.3068		54.1	50.0	8.2	
Isopropylbenzene	Ave	3.0830	3.2747		53.1	50.0	6.2	
m-Xylene & p-Xylene	Ave	0.4812	0.4997		104	100	3.9	
Methacrylonitrile	Ave	0.4469	0.4995		55.9	50.0	11.8	

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Instrument ID: MSL Calibration Date: 08/02/2007 Time: 09:18

Lab File ID: L9371.D Init. Calib. Date(s): 07/26/2007 07/26/2007

Lab Sample ID: CCVIS 220-8356/1 Init. Calib. Time(s): 13:32 15:11

GC Column: RTX-VMS ID: 0.18 (mm) Heated Purge: (Y/N) N Conc. Units: ug/L

Analyte	Curve	Ave RRF	RRF	Min RRF	Calculated Amount	Ccal Amount	% D	Max % D
Methyl acetate	Ave	1.8684	1.9307		51.7	50.0	3.3	
Methyl acrylate	Ave	0.4003	0.4314		53.9	50.0	7.8	
Methyl Ethyl Ketone	Ave	0.2301	0.2470		53.7	50.0	7.3	
methyl isobutyl ketone	Ave	0.4602	0.4649		50.5	50.0	1.0	
Methyl methacrylate	Ave	0.1736	0.3622		104	50.0	109	
Methyl tert-butyl ether	Ave	1.0379	1.0680		51.4	50.0	2.9	
Methylcyclohexane	Ave	0.3498	0.3628		51.9	50.0	3.7	
Methylene Chloride	Ave	0.2929	0.3088		52.7	50.0	5.4	
n-Butanol	Ave	0.0130	0.0135		518	500	3.5	
n-Butyl acetate	Ave	0.2901	0.2977		51.3	50.0	2.6	
n-Butylbenzene	Ave	3.3672	3.9512		58.7	50.0	17.3	
n-Heptane	Ave	0.3090	0.3306		53.5	50.0	7.0	
n-Propyl acetate	Ave	0.0316	0.0495		157	100	56.6	
N-Propylbenzene	Ave	3.3039	3.4341		52.0	50.0	3.9	
Naphthalene	Ave	2.2068	2.3706		53.7	50.0	7.4	
Nitrobenzene	Ave	0.0473	0.0397		420	500	-16.0	
o-Xylene	Ave	0.4609	0.4918		53.4	50.0	6.7	
p-Diethylbenzene	Ave	0.5445	0.5796		53.2	50.0	6.4	
Pentachloroethane	Ave				0.0			
Propionitrile	Ave	0.0634	0.0665		524	500	4.8	
sec-Butylbenzene	Ave	2.6102	2.6991		51.7	50.0	3.4	
Styrene	Ave	0.7721	0.8190		53.0	50.0	6.1	
Tert-amyl methyl ether	Ave	1.0524	1.1024		52.4	50.0	4.8	
Tert-butyl ethyl ether	Ave	1.2132	1.2997		53.6	50.0	7.1	
tert-Butyl Formate	Ave	0.3499	0.3652		52.2	50.0	4.4	
tert-Butylbenzene	Ave	2.3884	2.3903		50.0	50.0	0.1	
Tetrachloroethene	Ave	0.2100	0.2141		51.0	50.0	1.9	
Tetrahydrofuran	Ave	0.1355	0.1395		103	100	3.0	
Toluene	Ave	1.0645	1.0897		51.2	50.0	2.4	20.0
Toluene-d8 (Surr)	Ave	0.8889	0.6960		19.6	25.0	-21.7	
trans-1,2-Dichloroethene	Ave	0.2676	0.2930		54.7	50.0	9.5	
trans-1,3-Dichloropropene	Ave	0.5141	0.5575		54.2	50.0	8.4	
trans-1,4-Dichloro-2-butene	Ave	0.3461	0.3398		98.2	100	-1.8	
Trichloroethene	Ave	0.3531	0.3758		53.2	50.0	6.5	
Trichlorofluoromethane	Ave	0.3584	0.3727		52.0	50.0	4.0	
Vinyl acetate	Ave	0.9114	1.0031		55.0	50.0	10.1	
Vinyl chloride	Ave	0.3701	0.3675		49.6	50.0	-0.7	20.0
Xylenes, Total	Ave	0.3953	0.4971		157	150	25.7	

Curve Types:
Ave = Average
Lin = Linear
Quad = Quadratic

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079370.b\L9371.D
Lab Smp Id: CCVIS Client Smp ID: CCVIS
Inj Date : 02-AUG-2007 09:18 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : CCVIS
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 56 Continuing Calibration Sample
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONMSLNT

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	ON-COL (ug/L)
* 1	Fluorobenzene	96	4.901	4.901	(1.000)	436789	25.0000	
2	Dichlorodifluoromethane	85	1.143	1.143	(0.233)	155467	50.0000	46
3	Chloromethane	50	1.261	1.261	(0.257)	290845	50.0000	50
4	Vinyl Chloride	62	1.300	1.300	(0.265)	321051	50.0000	50
5	Bromomethane	94	1.477	1.477	(0.301)	119749	50.0000	53
6	Chloroethane	64	1.546	1.546	(0.316)	212012	50.0000	62
7	Trichlorofluoromethane	101	1.625	1.625	(0.332)	325550	50.0000	52
8	Dichlorofluoromethane	67	1.645	1.645	(0.336)	622033	50.0000	54
9	Ethyl Ether	45	1.792	1.792	(0.366)	216367	50.0000	55
10	Ethanol	45	1.861	1.861	(0.380)	170532	500.000	550
11	Freon 141	81	1.861	1.861	(0.380)	406535	50.0000	53
12	Freon 123a	67	1.645	1.645	(0.336)	622033	50.0000	54
13	Trichlorotrifluoroethane	101	1.950	1.950	(0.398)	240577	50.0000	56
14	1,1-Dichloroethene	96	1.930	1.930	(0.394)	202206	50.0000	54
15	Carbon Disulfide	76	1.969	1.969	(0.402)	961046	50.0000	53
16	Iodomethane	142	2.038	2.038	(0.416)	186963	50.0000	39
17	Acrolein	56	2.127	2.127	(0.434)	295881	250.000	290
18	2-Propanol	45	2.215	2.215	(0.452)	52066	50.0000	52
19	3-Chloro-1-Propene	41	2.225	2.225	(0.454)	527810	50.0000	55
20	Methylene Chloride	84	2.304	2.304	(0.470)	269741	50.0000	53
21	Acetone	43	2.323	2.323	(0.474)	169354	50.0000	61
22	trans-1,2-Dichloroethene	96	2.422	2.422	(0.494)	255914	50.0000	55

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	AMOUNTS ON-COL (ug/L)
	23 Methyl Acetate	43	2.402	2.402	(0.490)	1686599	50.0000	52
	24 Methyl tert-Butyl Ether	73	2.491	2.491	(0.508)	932996	50.0000	51
	25 tert-Butyl alcohol	59	2.530	2.530	(0.516)	248520	250.000	240
	26 Acetonitrile	41	2.658	2.658	(0.542)	604181	500.000	520
	27 Isopropyl ether	45	2.786	2.786	(0.568)	1141587	50.0000	54
	28 tert-Butyl ethyl ether	59	3.111	3.111	(0.635)	1135391	50.0000	54
	29 2-Chloro-1,3-Butadiene	88	2.884	2.884	(0.589)	181059	50.0000	54
	30 Acrylonitrile	53	2.924	2.924	(0.597)	354776	100.000	110
	31 1,1-Dichloroethane	63	2.904	2.904	(0.593)	642966	50.0000	54
	32 Vinyl Acetate	43	3.111	3.111	(0.635)	876246	50.0000	55
	33 cis-1,2-Dichloroethene	96	3.416	3.416	(0.697)	265634	50.0000	53
	34 2,2-Dichloropropane	77	3.524	3.524	(0.719)	453044	50.0000	55
	35 Bromochloromethane	128	3.622	3.622	(0.739)	191262	50.0000	53
	36 1-Bromopropane	43	3.612	3.612	(0.737)	486864	50.0000	54
	37 Cyclohexane	84	3.652	3.652	(0.745)	348248	50.0000	54
	38 Chloroform	83	3.701	3.701	(0.755)	481135	50.0000	53
	39 Ethyl Acetate	43	3.858	3.858	(0.787)	50716	100.000	100
	40 Methyl Acrylate	55	3.858	3.858	(0.787)	376818	50.0000	54
$	41 Dibromofluoromethane	111	3.927	3.927	(0.801)	108164	25.0000	19
	42 Tetrahydrofuran	42	3.917	3.917	(0.799)	243667	100.000	100
	43 Carbon Tetrachloride	117	3.898	3.898	(0.795)	438237	50.0000	58
	44 1,1,1-Trichloroethane	97	3.967	3.967	(0.809)	361304	50.0000	54
	45 2-Butanone	43	4.065	4.065	(0.829)	215746	50.0000	54
	46 1,1-Dichloropropene	75	4.124	4.124	(0.841)	441547	50.0000	53
	47 tert-Amyl methyl ether	73	4.567	4.567	(0.932)	963067	50.0000	52
	48 tert-Butyl formate	57	3.111	3.111	(0.635)	319024	50.0000	52
	49 1-Chlorobutane	56	4.173	4.173	(0.851)	660820	50.0000	55
	50 Heptane	43	4.399	4.399	(0.898)	288789	50.0000	53
	51 Propionitrile	54	4.399	4.399	(0.898)	580650	500.000	520
	52 Benzene	78	4.419	4.419	(0.902)	1120770	50.0000	53
	53 2-Methyl-2-Propenenitrile	41	4.439	4.439	(0.906)	436326	50.0000	56
	54 Isobutyl alcohol	42	4.685	4.685	(0.956)	142742	500.000	500
$	55 1,2-Dichloroethane-d4	65	4.567	4.567	(0.932)	124675	25.0000	18
	56 1,2-Dichloroethane	62	4.645	4.645	(0.948)	443604	50.0000	53
	59 Methyl Cyclohexane	83	5.098	5.098	(1.040)	316971	50.0000	52
	60 Trichloroethene	130	5.098	5.098	(1.040)	328316	50.0000	53
	61 Isopropyl Acetate	43	4.685	4.685	(0.956)	232034	100.000	99
	62 N-Butanol	56	5.482	5.482	(1.118)	117771	500.000	520
	63 Dibromomethane	93	5.531	5.531	(1.128)	182066	50.0000	54
	64 1,2-Dichloropropane	63	5.629	5.629	(1.149)	372482	50.0000	54
	65 Bromodichloromethane	83	5.718	5.718	(1.167)	344037	50.0000	53
	66 Methyl Methacrylate	69	5.895	5.895	(1.203)	316392	100.000	100
	67 1,4-Dioxane	58	5.934	5.934	(1.211)	28552	500.000	540
	68 N-Propyl Acetate	43	6.298	6.298	(1.285)	86508	100.000	160
	69 2-Chloroethylvinylether	63	6.298	6.298	(1.285)	128432	50.0000	76
	70 cis-1,3-Dichloropropene	75	6.348	6.348	(1.295)	527995	50.0000	53
	71 Chloroacetonitrile	48	6.692	6.692	(1.365)	149131	500.000	530
	72 2-Nitropropane	41	6.771	6.771	(1.381)	197367	100.000	100
	73 trans-1,3-Dichloropropene	75	6.977	6.977	(1.424)	487016	50.0000	54
	74 1,1,2-Trichloroethane	97	7.125	7.125	(1.454)	233051	50.0000	53
*	75 Chlorobenzene-d5	117	7.961	7.961	(1.000)	448655	25.0000	
	76 Toluene	91	6.584	6.584	(0.827)	977836	50.0000	51
$	77 Toluene-d8	98	6.535	6.535	(0.821)	312259	25.0000	20
	78 1,1-Dichloro-2-propanone	43	6.800	6.800	(0.854)	1134417	250.000	240

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (ug/L)	ON-COL (ug/L)
79	4-Methyl-2-Pentanone	43	6.938	6.938	(0.871)	417120	50.0000	50
80	Tetrachloroethene	164	6.958	6.958	(0.874)	192125	50.0000	51
81	Ethyl Methacrylate	69	7.145	7.145	(0.897)	514838	50.0000	51
82	Dibromochloromethane	129	7.292	7.292	(0.916)	350753	50.0000	49
83	1,3-Dichloropropane	76	7.361	7.361	(0.925)	516940	50.0000	50
84	1,2-Dibromoethane	107	7.489	7.489	(0.941)	312563	50.0000	52
85	n-Butyl Acetate	56	7.646	7.646	(0.960)	267166	50.0000	51
86	2-Hexanone	43	7.705	7.705	(0.968)	326714	50.0000	53
87	1-Chlorohexane	91	7.971	7.971	(1.001)	316792	50.0000	49
88	Chlorobenzene	112	7.971	7.971	(1.001)	803262	50.0000	52
89	1,1,1,2-Tetrachloroethane	131	8.040	8.040	(1.010)	273621	50.0000	49
90	Ethylbenzene	106	8.010	8.010	(1.006)	370130	50.0000	54
91	Xylene (total)mp	106	8.138	8.138	(1.022)	896773	100.000	100
92	Xylene (total)o	106	8.512	8.512	(1.069)	441304	50.0000	53
93	Styrene	104	8.561	8.561	(1.075)	734851	50.0000	53
94	Bromoform	173	8.581	8.581	(1.078)	197842	50.0000	50
* 95	1,4-Dichlorobenzene-d4	152	10.018	10.018	(1.000)	157752	25.0000	
96	Isopropylbenzene	105	8.797	8.797	(0.878)	1033187	50.0000	53
97	Bromobenzene	156	9.122	9.122	(0.911)	256098	50.0000	52
98	1,1,2,2-Tetrachloroethane	83	9.221	9.221	(0.920)	347066	50.0000	52
99	4-Ethyltoluene	105	9.260	9.260	(0.924)	1007220	50.0000	53
100	1,2,3-Trichloropropane	110	9.329	9.329	(0.931)	107566	50.0000	50
101	trans-1,4-Dichloro-2-Butene	53	9.368	9.368	(0.935)	214424	100.000	98
102	n-Propylbenzene	91	9.162	9.162	(0.915)	1083456	50.0000	52
103	2-Chlorotoluene	91	9.289	9.289	(0.927)	704363	50.0000	49
104	4-Chlorotoluene	91	9.437	9.437	(0.942)	718861	50.0000	53
105	1,3,5-Trimethylbenzene	105	9.339	9.339	(0.932)	831688	50.0000	53
106	tert-Butylbenzene	119	9.604	9.604	(0.959)	754140	50.0000	50
107	1,2,4-Trimethylbenzene	105	9.673	9.673	(0.966)	825899	50.0000	53
108	sec-Butylbenzene	105	9.762	9.762	(0.974)	851572	50.0000	52
109	4-Isopropyltoluene	119	9.890	9.890	(0.987)	916267	50.0000	52
110	1,3-Dichlorobenzene	146	9.949	9.949	(0.993)	452126	50.0000	53
111	1,4-Dichlorobenzene	146	10.027	10.027	(1.001)	452914	50.0000	52
112	1,2-Dichlorobenzene	146	10.391	10.391	(1.037)	434751	50.0000	51
113	Benzyl Chloride	126	10.234	10.234	(1.022)	151810	50.0000	58
114	1,4-Diethylbenzene	119	10.204	10.204	(2.082)	506336	50.0000	53
115	n-Butylbenzene	91	10.254	10.254	(1.024)	1246625	50.0000	59
118	1,2,4,5-Tetramethylbenzene	119	10.913	10.913	(2.226)	792315	50.0000	55
119	1,2-Dibromo-3-chloropropane	75	11.080	11.080	(1.106)	66810	50.0000	50
120	Nitrobenzene	77	11.572	11.572	(1.155)	125298	500.000	420
121	1,2,4-Trichlorobenzene	180	11.690	11.690	(1.167)	248340	50.0000	54
122	Hexachlorobutadiene	225	11.670	11.670	(1.165)	78745	50.0000	50
123	Naphthalene	128	11.966	11.966	(1.194)	747935	50.0000	54
124	1,2,3-Trichlorobenzene	180	12.133	12.133	(1.211)	223076	50.0000	53
$ 125	Bromofluorobenzene	95	9.043	9.043	(0.903)	139090	25.0000	24
M 126	1,2-Dichloroethene (total)	100				521548	100.000	110
M 127	Xylene (total)	100				1338077	150.000	160

Data File: L9371.D

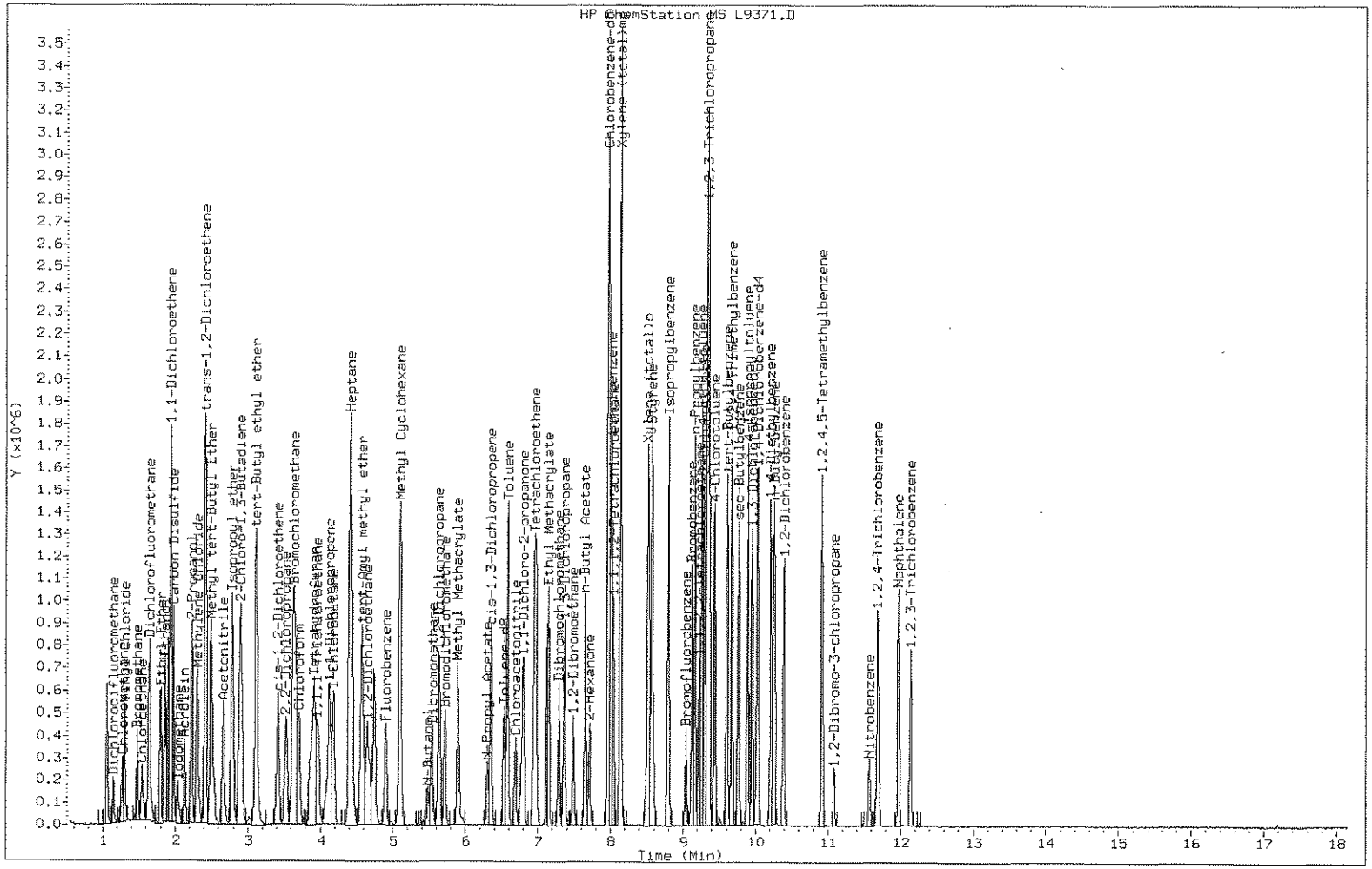
Date: 02-AUG-2007 09:18

Client ID: CCVIS

Instrument: msl.i

Sample Info: CCVIS

Operator: D. HUMBERT



08/09/2007

STL-CT

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079145.b\LB417.D
Lab Smp Id: BFB Client Smp ID: BFB
Inj Date : 26-JUL-2007 13:09 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : BFB
Misc Info : : ;;; BFB ; 8260B ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079145.b\LBFBNCLP.M
Meth Date : 16-May-2007 12:03 pattym Quant Type: ISTD
Cal Date : Cal File:
Als bottle: 29 QC Sample: BFB
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14 Sample Matrix: WATER
Processing Host: CONMSLNT

Concentration Formula: Amt * DF * Uf * Vf * VI * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vf	1.000	Volumetric correction factor
VI	2.000	Injection Volume
Cpnd Variable		Local Compound Variable

RT	EXP RT	REL RT	MASS	RESPONSE	CONCENTRATIONS ON-COL (ug/L)	CONCENTRATIONS FINAL (ug/L)	TARGET RANGE	RATIO
1 bfb						CAS #: 460-00-4		
3.269	3.400	(0.000)	95	78760			0.00- 100.00	100.00
3.269	3.400	(0.000)	50	15646			15.00- 40.00	19.87
3.269	3.400	(0.000)	75	43208			30.00- 60.00	54.86
3.269	3.400	(0.000)	96	4903			5.00- 9.00	6.23
3.269	3.400	(0.000)	173	0	0.0	0.0	0.00- 2.00	0.00
3.269	3.400	(0.000)	174	68864			50.00- 100.00	87.44
3.269	3.400	(0.000)	175	4851			5.00- 9.00	7.04
3.269	3.400	(0.000)	176	69264			95.00- 101.00	100.58
3.269	3.400	(0.000)	177	4313			5.00- 9.00	6.23

Data File: LB417.D

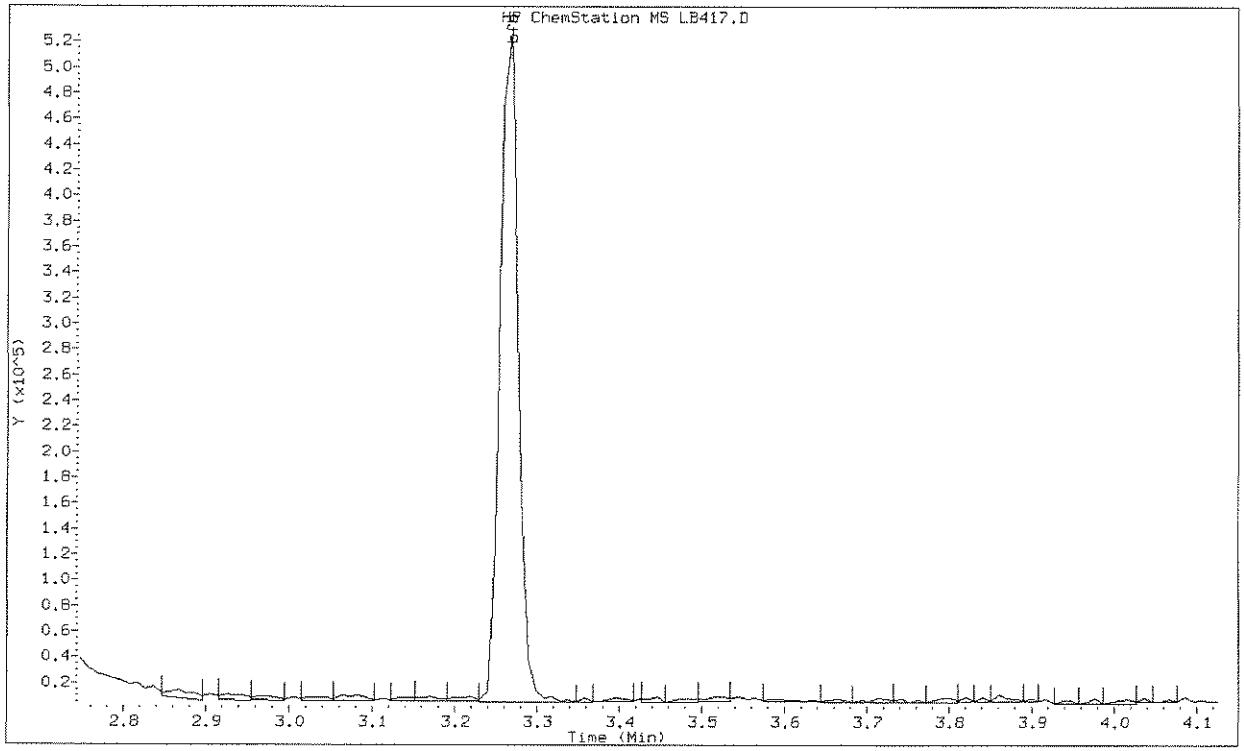
Date: 26-JUL-2007 13:09

Client ID: BFB

Instrument: msl.i

Sample Info: BFB

Operator: D. HUMBERT



Data File: LB417.D

Date: 26-JUL-2007 13:09

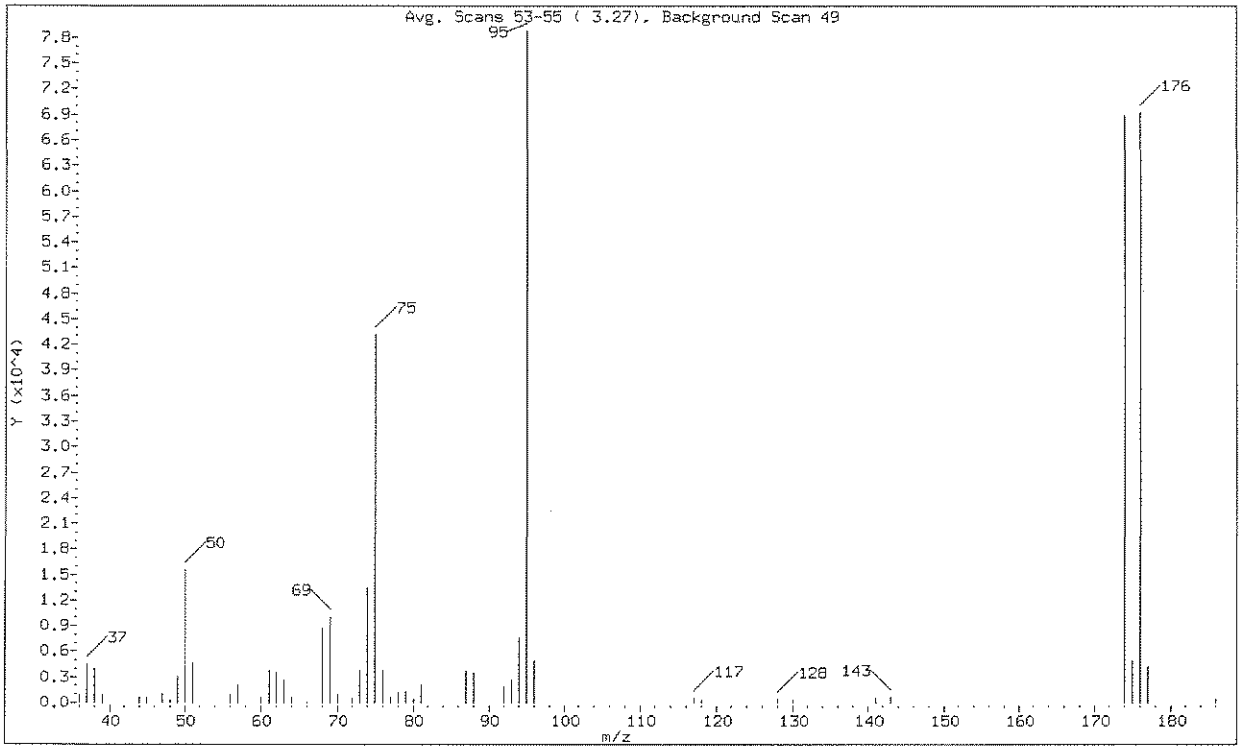
Client ID: BFB

Instrument: msl.i

Sample Info: BFB

Operator: D. HUMBERT

1 bfb



m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
95	Base Peak, 100% relative abundance	100.00
50	15.00 - 40.00% of mass 95	19.87
75	30.00 - 60.00% of mass 95	54.86
96	5.00 - 9.00% of mass 95	6.23
173	Less than 2.00% of mass 174	0.00 (0.00)
174	50.00 - 100.00% of mass 95	87.44
175	5.00 - 9.00% of mass 174	6.16 (7.04)
176	95.00 - 101.00% of mass 174	87.94 (100.58)
177	5.00 - 9.00% of mass 176	5.48 (6.23)

Data File: LB417.D

Date: 26-JUL-2007 13:09

Client ID: BFB

Instrument: msl.i

Sample Info: BFB

Operator: D. HUMBERT

Data File: \\target1_ct\files\chem\VOA\msl.i\L079145.b\LB417.D
Spectrum: Avg. Scans 53-55 (3.27), Background Scan 49
Location of Maximum: 95.00
Number of points: 50

m/z	Y	m/z	Y	m/z	Y	m/z	Y
36.00	903	60.00	692	76.00	3709	117.00	485
37.00	4539	61.00	3689	77.00	583	118.00	237
38.00	3936	62.00	3435	78.00	1101	128.00	408
39.00	928	63.00	2528	79.00	1344	130.00	214
44.00	682	64.00	637	80.00	427	141.00	552
45.00	533	66.00	190	81.00	2095	143.00	606
47.00	1056	68.00	8803	87.00	3785	174.00	68864
48.00	199	69.00	10024	88.00	3472	175.00	4851
49.00	3036	70.00	851	92.00	1806	176.00	69264
50.00	15646	72.00	526	93.00	2606	177.00	4313
51.00	4682	73.00	3768	94.00	7586	186.00	371
56.00	940	74.00	13365	95.00	78760		
57.00	2073	75.00	43208	96.00	4903		

STL-CT

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079310.b\LB425.D
Lab Smp Id: BFB Client Smp ID: BFB
Inj Date : 31-JUL-2007 20:13 MS Autotune Date: 26-APR-2004 14:21
Operator : D. GAYDA Inst ID: msl.i
Smp Info : BFB
Misc Info : : ;;; BFB ; 8260B ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079310.b\LBFBNCLP.M
Meth Date : 16-May-2007 12:03 pattym Quant Type: ISTD
Cal Date : Cal File:
Als bottle: 33 QC Sample: BFB
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14 Sample Matrix: WATER
Processing Host: CONMSLNT

Concentration Formula: Amt * DF * Uf * Vf * VI * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vf	1.000	Volumetric correction factor
VI	2.000	Injection Volume
Cpnd Variable		Local Compound Variable

RT	EXP RT	REL RT	MASS	RESPONSE	CONCENTRATIONS ON-COL (ug/L)	CONCENTRATIONS FINAL (ug/L)	TARGET RANGE	RATIO
1 bfb						CAS #: 460-00-4		
3.254	3.400	(0.000)	95	82768			0.00- 100.00	100.00
3.254	3.400	(0.000)	50	17536			15.00- 40.00	21.19
3.254	3.400	(0.000)	75	45632			30.00- 60.00	55.13
3.254	3.400	(0.000)	96	6082			5.00- 9.00	7.35
3.254	3.400	(0.000)	173	232			0.00- 2.00	0.31
3.254	3.400	(0.000)	174	73952			50.00- 100.00	89.35
3.254	3.400	(0.000)	175	5631			5.00- 9.00	7.61
3.254	3.400	(0.000)	176	73912			95.00- 101.00	99.95
3.254	3.400	(0.000)	177	4687			5.00- 9.00	6.34

Data File: LB425.D

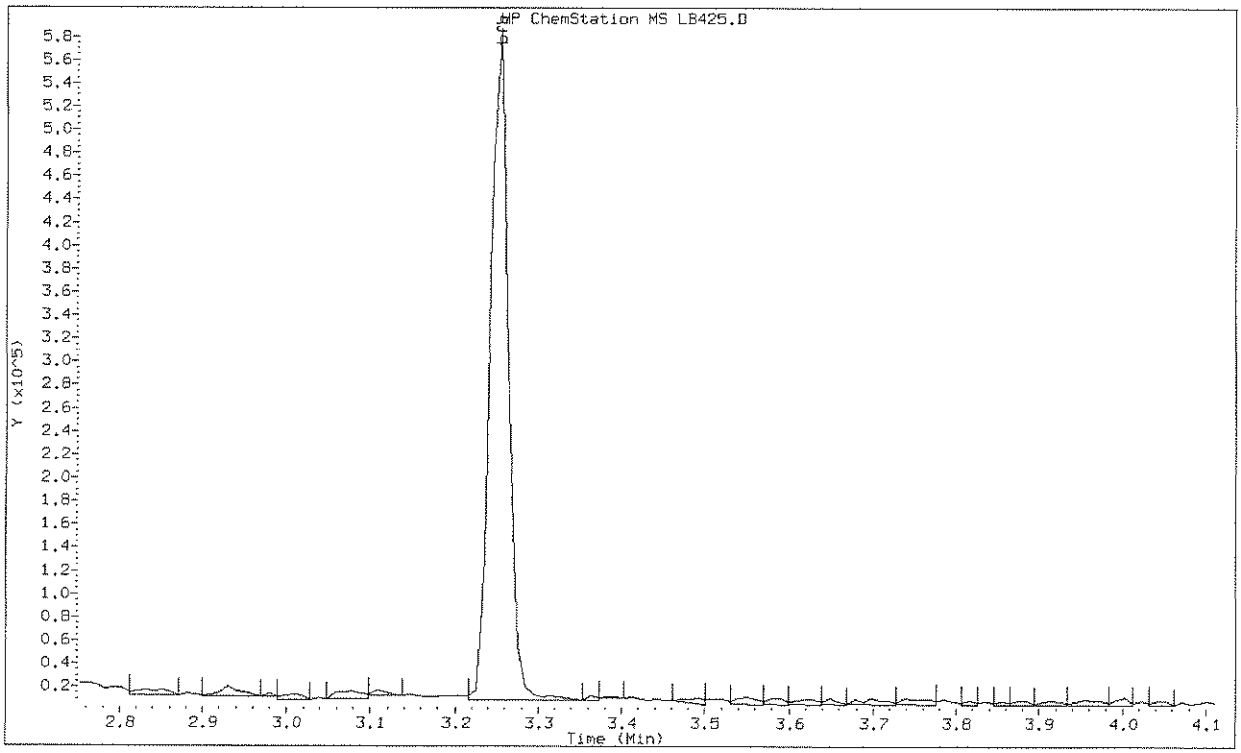
Date: 31-JUL-2007 20:13

Client ID: BFB

Instrument: msl.i

Sample Info: BFB

Operator: D. GAYDA



Data File: LB425.D

Date: 31-JUL-2007 20:13

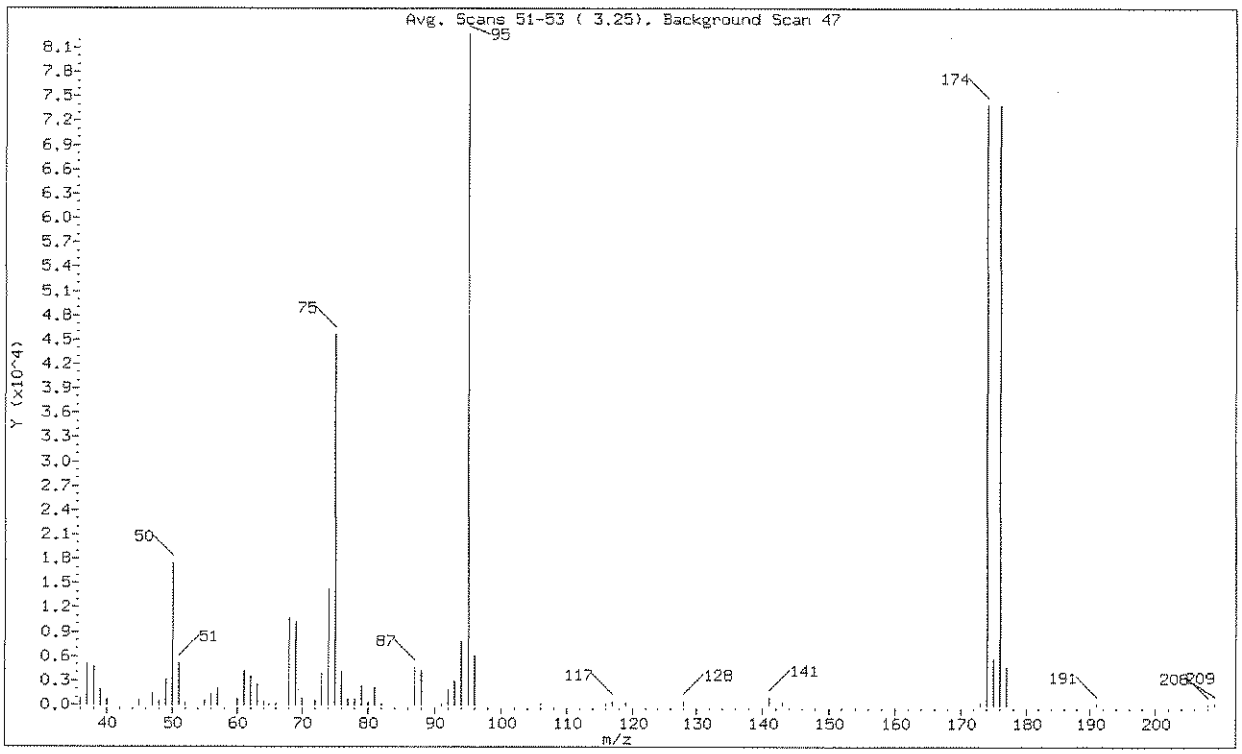
Client ID: BFB

Instrument: msl.i

Sample Info: BFB

Operator: D. GAYDA

1 bfb



m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
95	Base Peak, 100% relative abundance	100.00
50	15.00 - 40.00% of mass 95	21.19
75	30.00 - 60.00% of mass 95	55.13
96	5.00 - 9.00% of mass 95	7.35
173	Less than 2.00% of mass 174	0.28 (0.31)
174	50.00 - 100.00% of mass 95	89.35
175	5.00 - 9.00% of mass 174	6.80 (7.61)
176	95.00 - 101.00% of mass 174	89.30 (99.95)
177	5.00 - 9.00% of mass 176	5.66 (6.34)

Data File: LB425.D

Date: 31-JUL-2007 20:13

Client ID: BFB

Instrument: msl.i

Sample Info: BFB

Operator: D. GAYDA

Data File: \\target1_ct\files\chem\VOA\msl.i\L079310.b\LB425.D
Spectrum: Avg. Scans 51-53 (3.25), Background Scan 47
Location of Maximum: 95.00
Number of points: 59

m/z	Y	m/z	Y	m/z	Y	m/z	Y
36.00	850	60.00	849	77.00	640	116.00	169
37.00	5061	61.00	4228	78.00	655	117.00	387
38.00	4792	62.00	3507	79.00	2254	119.00	217
39.00	1857	63.00	2410	80.00	266	128.00	357
40.00	661	64.00	435	81.00	2108	141.00	978
45.00	623	65.00	22	82.00	172	143.00	385
47.00	1415	66.00	192	87.00	4539	173.00	232
48.00	441	68.00	10733	88.00	4138	174.00	73952
49.00	3067	69.00	10184	91.00	175	175.00	5631
50.00	17536	70.00	844	92.00	1960	176.00	73912
51.00	5173	72.00	554	93.00	2943	177.00	4687
52.00	228	73.00	3954	94.00	7890	191.00	172
55.00	530	74.00	14260	95.00	82768	208.00	30
56.00	1290	75.00	45632	96.00	6082	209.00	192
57.00	2015	76.00	4091	106.00	180		

STL-CT

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079341.b\LB426.D
Lab Smp Id: BFB Client Smp ID: BFB
Inj Date : 01-AUG-2007 09:13 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : BFB
Misc Info : : ;;; BFB ; 8260B ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079341.b\LBFBNCLP.M
Meth Date : 16-May-2007 12:03 pattym Quant Type: ISTD
Cal Date : Cal File:
Als bottle: 33 QC Sample: BFB
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14 Sample Matrix: WATER
Processing Host: CONMSLNT

Concentration Formula: Amt * DF * Uf * Vf * VI * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vf	1.000	Volumetric correction factor
VI	2.000	Injection Volume
Cpnd Variable		Local Compound Variable

RT	EXP RT	REL RT	MASS	RESPONSE	CONCENTRATIONS ON-COL (ug/L)	CONCENTRATIONS FINAL (ug/L)	TARGET RANGE	RATIO
1 bfb						CAS #: 460-00-4		
3.264	3.400	(0.000)	95	102928			0.00- 100.00	100.00
3.264	3.400	(0.000)	50	21160			15.00- 40.00	20.56
3.264	3.400	(0.000)	75	55200			30.00- 60.00	53.63
3.264	3.400	(0.000)	96	6703			5.00- 9.00	6.51
3.264	3.400	(0.000)	173	535			0.00- 2.00	0.58
3.264	3.400	(0.000)	174	92384			50.00- 100.00	89.76
3.264	3.400	(0.000)	175	6947			5.00- 9.00	7.52
3.264	3.400	(0.000)	176	91376			95.00- 101.00	98.91
3.264	3.400	(0.000)	177	5983			5.00- 9.00	6.55

Data File: LB426.D

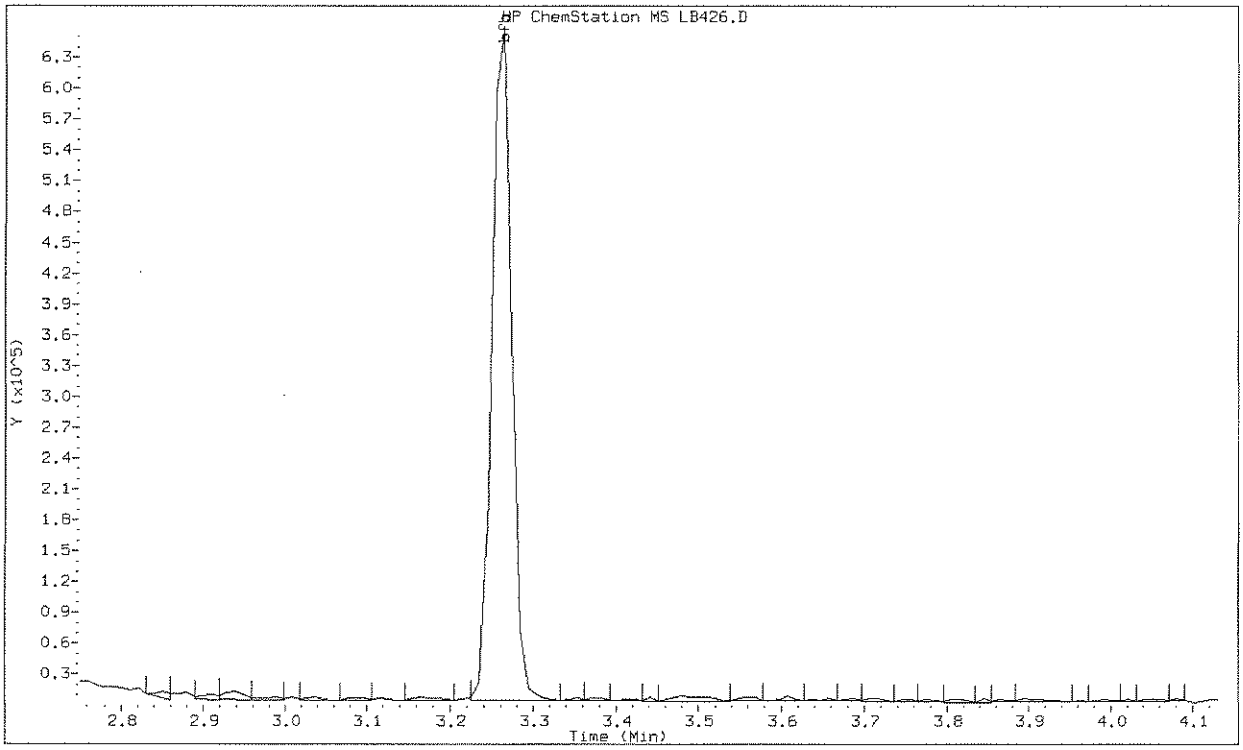
Date: 01-AUG-2007 09:13

Client ID: BFB

Instrument: msl.i

Sample Info: BFB

Operator: D. HUMBERT



Data File: LB426.D

Date: 01-AUG-2007 09:13

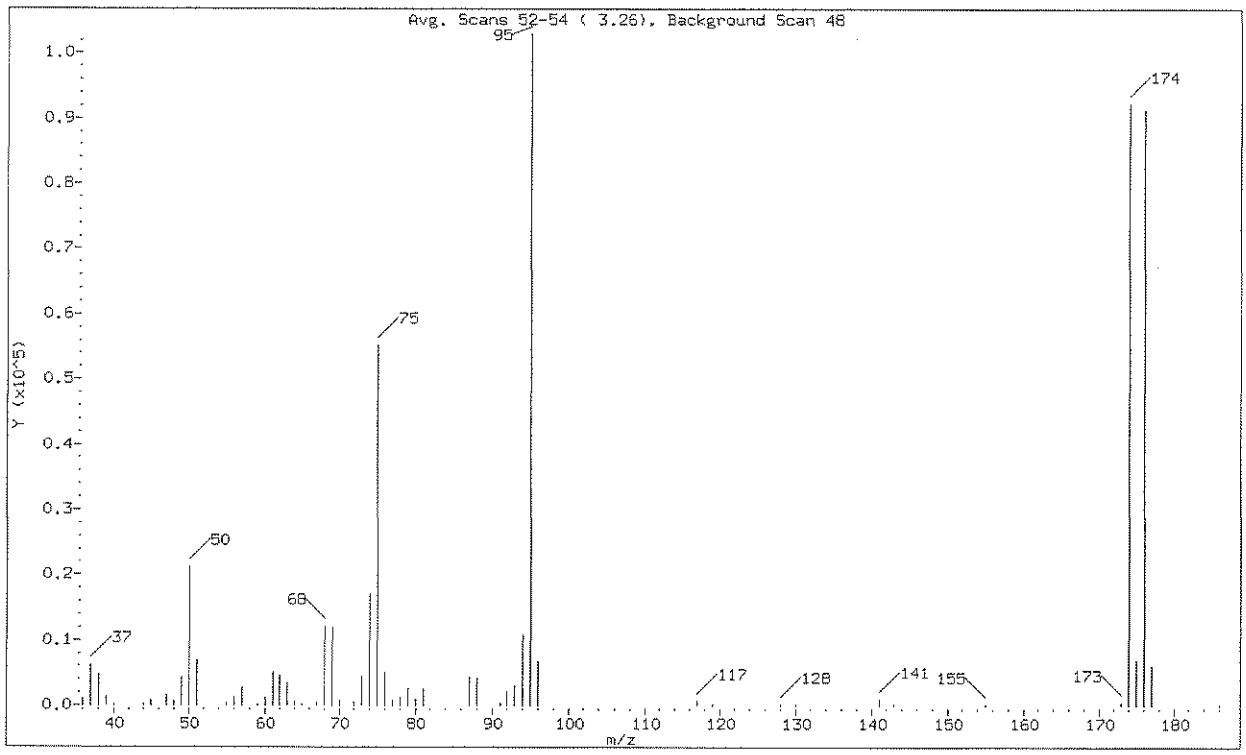
Client ID: BFB

Instrument: msl.i

Sample Info: BFB

Operator: D. HUMBERT

1 bfb



m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
95	Base Peak, 100% relative abundance	100.00
50	15.00 - 40.00% of mass 95	20.56
75	30.00 - 60.00% of mass 95	53.63
96	5.00 - 9.00% of mass 95	6.51
173	Less than 2.00% of mass 174	0.52 (0.58)
174	50.00 - 100.00% of mass 95	89.76
175	5.00 - 9.00% of mass 174	6.75 (7.52)
176	95.00 - 101.00% of mass 174	88.78 (98.91)
177	5.00 - 9.00% of mass 176	5.81 (6.55)

Data File: LB426.D

Date: 01-AUG-2007 09:13

Client ID: BFB

Instrument: msl.i

Sample Info: BFB

Operator: D. HUMBERT

Data File: \\target1_ct\files\chem\VOA\msl.i\L079341.b\LB426.D
Spectrum: Avg. Scans 52-54 (3.26), Background Scan 48
Location of Maximum: 95.00
Number of points: 57

m/z	Y	m/z	Y	m/z	Y	m/z	Y
36.00	1049	59.00	220	76.00	4989	119.00	180
37.00	6260	60.00	1177	77.00	750	128.00	191
38.00	4694	61.00	5099	78.00	1240	130.00	169
39.00	1333	62.00	4515	79.00	2576	141.00	1064
40.00	190	63.00	3367	80.00	780	143.00	229
44.00	233	64.00	429	81.00	2459	155.00	175
45.00	829	65.00	181	87.00	4355	173.00	535
47.00	1502	67.00	495	88.00	4210	174.00	92384
48.00	704	68.00	12193	91.00	415	175.00	6947
49.00	4399	69.00	12021	92.00	2114	176.00	91376
50.00	21160	70.00	749	93.00	3106	177.00	5983
51.00	6852	72.00	539	94.00	10725	186.00	221
55.00	358	73.00	4427	95.00	102928		
56.00	1333	74.00	16968	96.00	6703		
57.00	2727	75.00	55200	117.00	719		

STL-CT

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\LB428.D
Lab Smp Id: BFB Client Smp ID: BFB
Inj Date : 02-AUG-2007 08:50 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : BFB
Misc Info : : ;;; BFB ; 8260B ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\LBFBNCLP.M
Meth Date : 16-May-2007 12:03 pattym Quant Type: ISTD
Cal Date : Cal File:
Als bottle: 33 QC Sample: BFB
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14 Sample Matrix: WATER
Processing Host: CONMSLNT

Concentration Formula: Amt * DF * Uf * Vf * VI * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vf	1.000	Volumetric correction factor
VI	2.000	Injection Volume
Cpnd Variable		Local Compound Variable

RT	EXP RT	REL RT	MASS	RESPONSE	CONCENTRATIONS ON-COL (ug/L)	CONCENTRATIONS FINAL (ug/L)	TARGET RANGE	RATIO
1 bfb						CAS #: 460-00-4		
3.265	3.400	(0.000)	95	89720			0.00- 100.00	100.00
3.265	3.400	(0.000)	50	19704			15.00- 40.00	21.96
3.265	3.400	(0.000)	75	50720			30.00- 60.00	56.53
3.265	3.400	(0.000)	96	6030			5.00- 9.00	6.72
3.265	3.400	(0.000)	173	235			0.00- 2.00	0.28
3.265	3.400	(0.000)	174	83792			50.00- 100.00	93.39
3.265	3.400	(0.000)	175	6395			5.00- 9.00	7.63
3.265	3.400	(0.000)	176	81168			95.00- 101.00	96.87
3.265	3.400	(0.000)	177	5101			5.00- 9.00	6.28

Data File: LB428.D

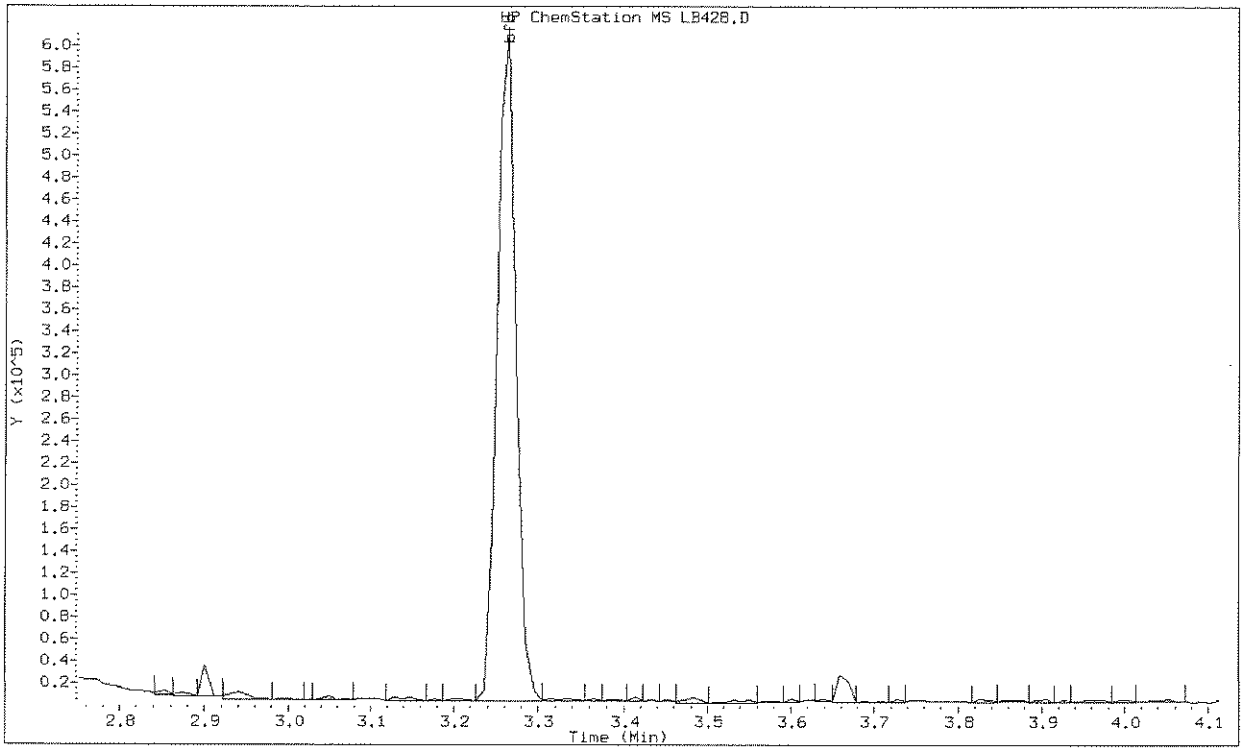
Date: 02-AUG-2007 08:50

Client ID: BFB

Instrument: msl.i

Sample Info: BFB

Operator: D. HUMBERT



Data File: LB428.D

Date: 02-AUG-2007 08:50

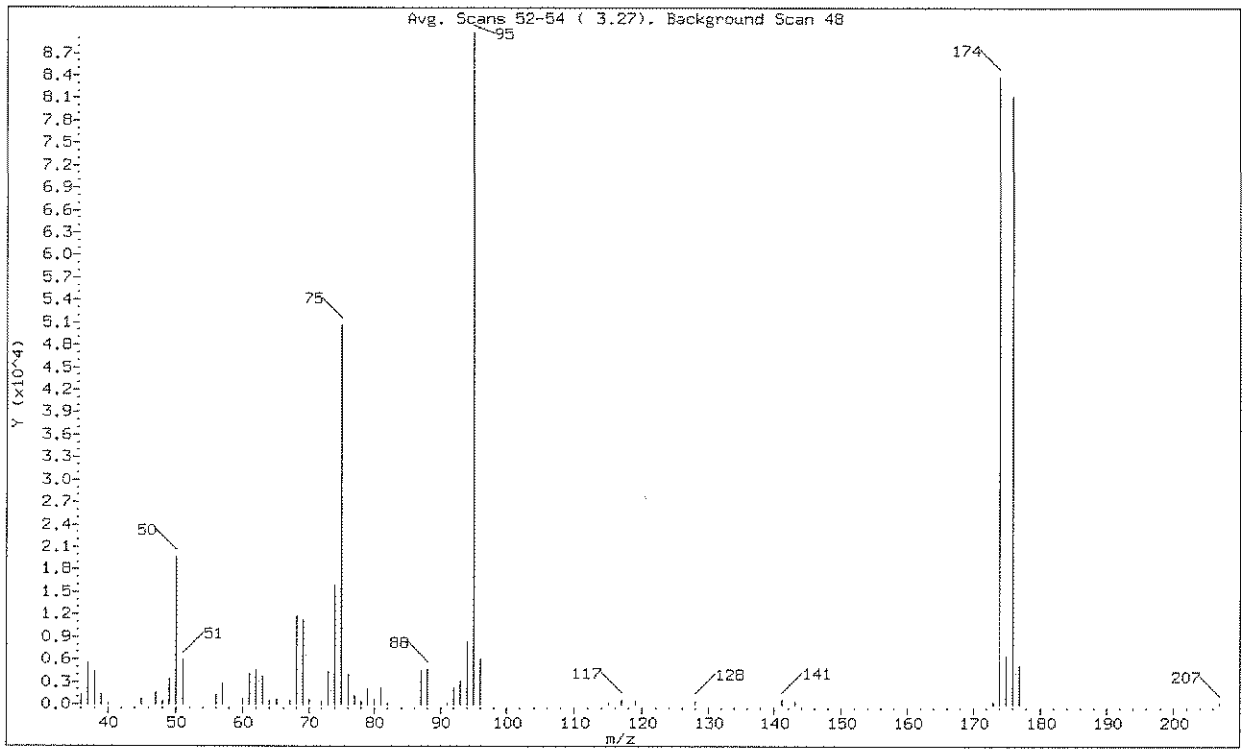
Client ID: BFB

Instrument: msl.i

Sample Info: BFB

Operator: D. HUMBERT

1 bfb



m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
95	Base Peak, 100% relative abundance	100.00
50	15.00 - 40.00% of mass 95	21.96
75	30.00 - 60.00% of mass 95	56.53
96	5.00 - 9.00% of mass 95	6.72
173	Less than 2.00% of mass 174	0.26 (0.28)
174	50.00 - 100.00% of mass 95	93.39
175	5.00 - 9.00% of mass 174	7.13 (7.63)
176	95.00 - 101.00% of mass 174	90.47 (96.87)
177	5.00 - 9.00% of mass 176	5.69 (6.28)

Data File: LB428.D

Date: 02-AUG-2007 08:50

Client ID: BFB

Instrument: msl.i

Sample Info: BFB

Operator: D. HUMBERT

Data File: \\target1_ct\files\chem\VOA\msl.i\L079370.b\LB428.D
Spectrum: Avg. Scans 52-54 (3.27), Background Scan 48
Location of Maximum: 95.00
Number of points: 54

m/z	Y	m/z	Y	m/z	Y	m/z	Y
36.00	1320	61.00	4147	77.00	1122	115.00	203
37.00	5653	62.00	4600	78.00	487	117.00	578
38.00	4403	63.00	3652	79.00	2244	119.00	409
39.00	1319	64.00	400	80.00	595	128.00	447
40.00	171	65.00	586	81.00	2239	141.00	563
45.00	771	67.00	433	82.00	189	143.00	410
47.00	1648	68.00	11769	87.00	4500	173.00	235
48.00	402	69.00	11291	88.00	4716	174.00	83792
49.00	3436	70.00	572	91.00	180	175.00	6395
50.00	19704	72.00	486	92.00	2199	176.00	81168
51.00	5820	73.00	4386	93.00	3062	177.00	5101
56.00	1357	74.00	15880	94.00	8409	207.00	187
57.00	2762	75.00	50720	95.00	89720		
60.00	763	76.00	3998	96.00	6030		

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: | Lab Sample ID: MB 220-8307/3

Matrix: Water | Lab File ID: L9313.D

Analysis Method: 8260B | Date Received:

Sample wt/vol: 5 (mL) | Date Analyzed: 07/31/2007 21:36

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8307 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	5.0	U	5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-CT

Volatile Report SW-846 Method 8260B

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079310.b\L9313.D
Lab Smp Id: MB
Inj Date : 31-JUL-2007 21:36 MS Autotune Date: 26-APR-2004 14:21
Operator : D. GAYDA Inst ID: msl.i
Smp Info : MB
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079310.b\L8260BNW.m
Meth Date : 31-Jul-2007 21:07 target Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 4 QC Sample: BLANK
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONMSLNT

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1 Fluorobenzene	96	4.897	4.896	(1.000)	450138	25.0000	
$ 41 Dibromofluoromethane	111	3.923	3.922	(0.801)	112197	19.3203	19
$ 55 1,2-Dichloroethane-d4	65	4.562	4.561	(0.932)	134039	18.9318	19
* 75 Chlorobenzene-d5	117	7.957	7.956	(1.000)	433216	25.0000	
$ 77 Toluene-d8	98	6.530	6.529	(0.821)	320395	20.7993	21
* 95 1,4-Dichlorobenzene-d4	152	10.013	10.012	(1.000)	130327	25.0000	
$ 125 Bromofluorobenzene	95	9.039	9.038	(0.903)	132069	27.0118	27

Data File: L9313.D

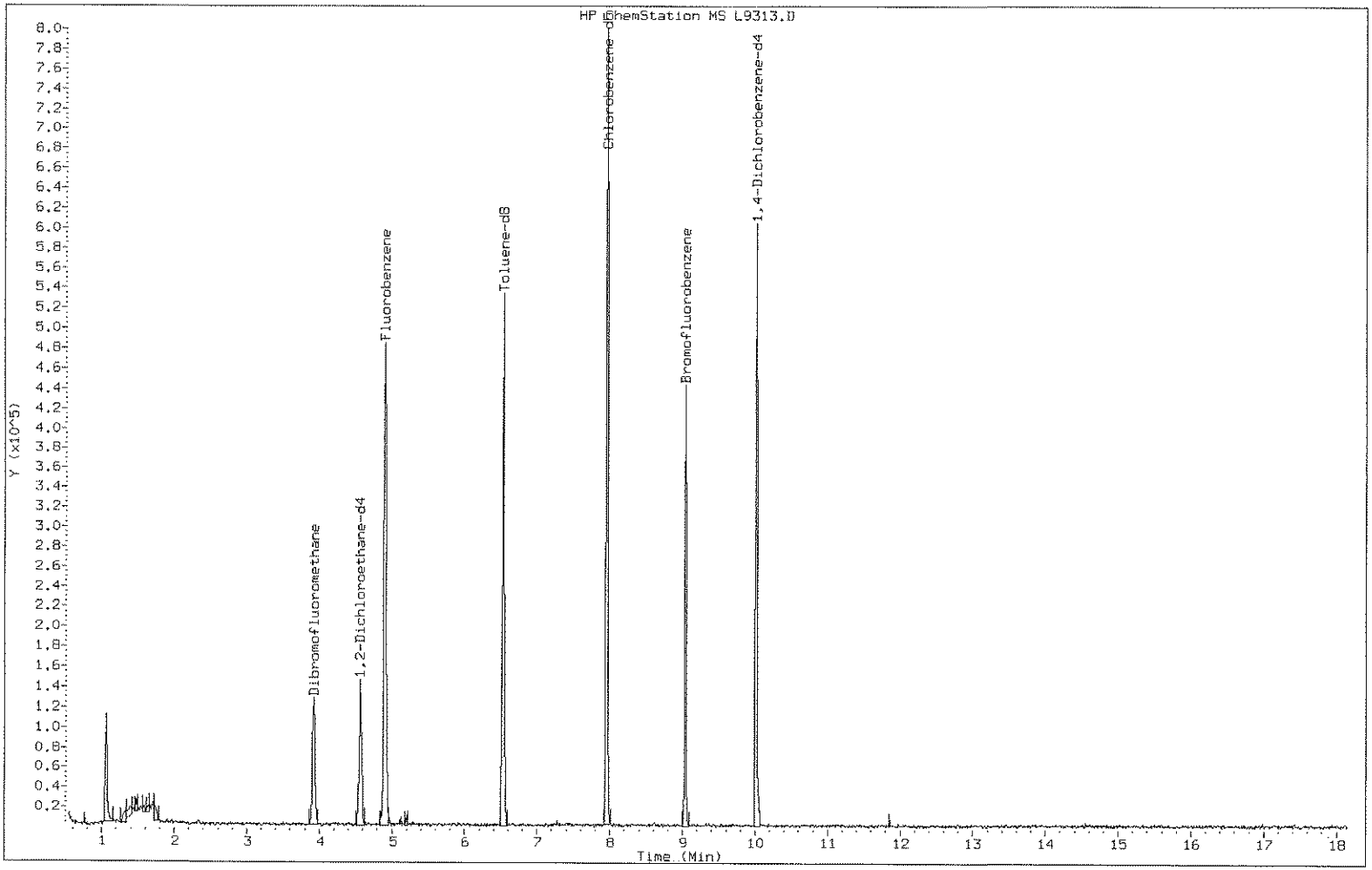
Date: 31-JUL-2007 21:36

Client ID: MB

Instrument: msl.i

Sample Info: MB

Operator: D. GAYDA



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: Lab Sample ID: MB 220-8321/3

Matrix: Water Lab File ID: L9344.D

Analysis Method: 8260B Date Received:

Sample wt/vol: 5 (mL) Date Analyzed: 08/01/2007 11:41

Level: (low/med) Low Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) Soil Aliquot:

Soil Extract Vol.: % Moisture:

Analy. Batch No.: 8321 Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	3.1	J	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	0.56	J	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	5.0	U	5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-CT

Volatile Report SW-846 Method 8260B

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079341.b\L9344.D
Lab Smp Id: MB Client Smp ID: MB
Inj Date : 01-AUG-2007 11:41 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : MB
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079341.b\L8260BNW.m
Meth Date : 02-Aug-2007 09:12 target Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 32 QC Sample: BLANK
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1 Fluorobenzene	96	4.896	4.908	(1.000)	449562	25.0000	
20 Methylene Chloride	84	2.298	2.300	(0.469)	2931	0.55647	0.6
21 Acetone	43	2.328	2.330	(0.476)	8948	3.13849	3
$ 41 Dibromofluoromethane	111	3.922	3.924	(0.801)	110095	18.9826	19
$ 55 1,2-Dichloroethane-d4	65	4.561	4.563	(0.932)	131294	18.5679	18
* 75 Chlorobenzene-d5	117	7.955	7.968	(1.000)	431702	25.0000	
$ 77 Toluene-d8	98	6.529	6.541	(0.821)	314635	20.4970	20
* 95 1,4-Dichlorobenzene-d4	152	10.012	10.014	(1.000)	124435	25.0000	
$ 125 Bromofluorobenzene	95	9.038	9.040	(0.903)	131324	28.1313	28

Data File: L9344.D

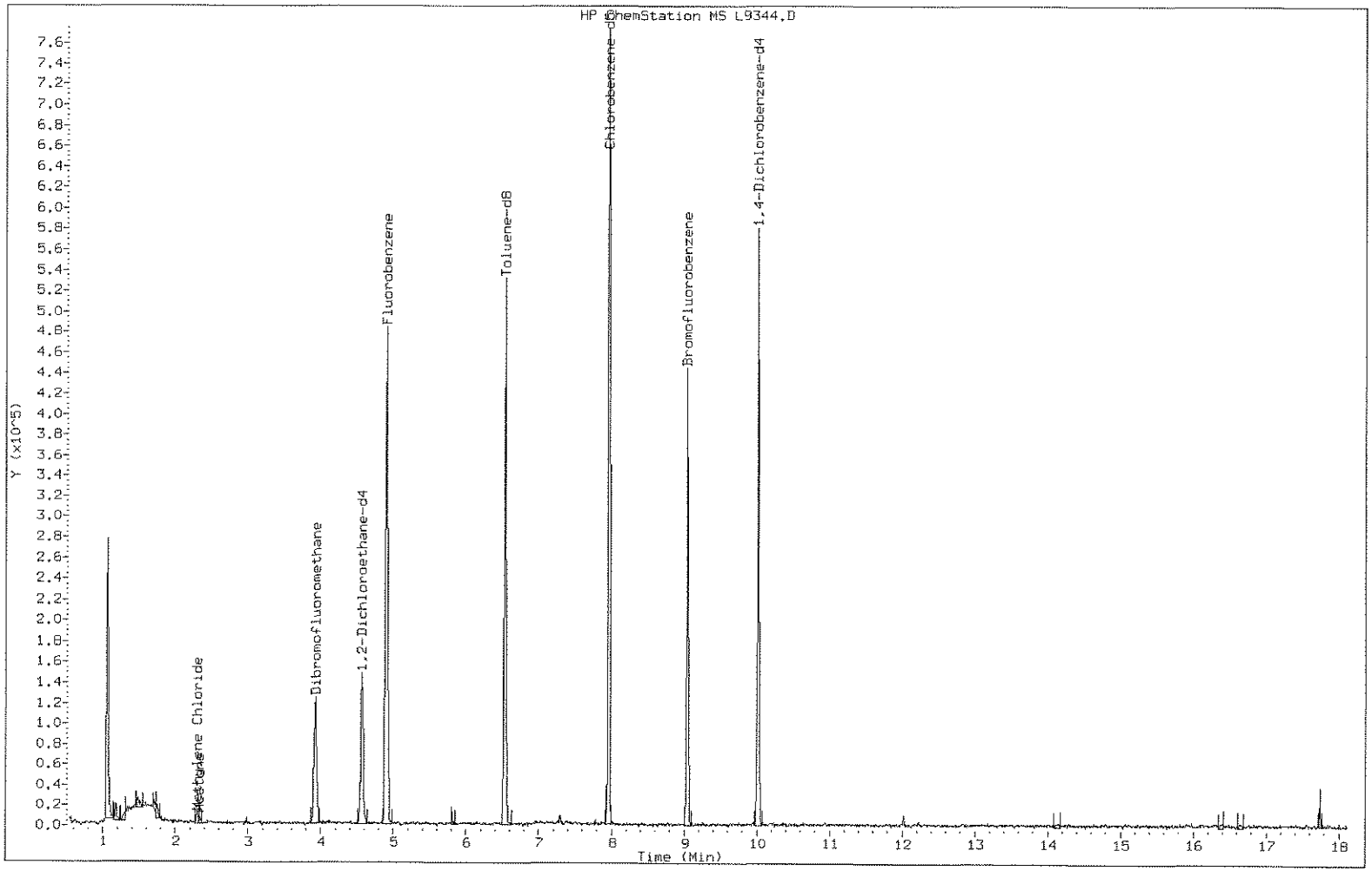
Date: 01-AUG-2007 11:41

Client ID: MB

Instrument: ms1.i

Sample Info: MB

Operator: D. HUMBERT



Data File: L9344.D

Date: 01-AUG-2007 11:41

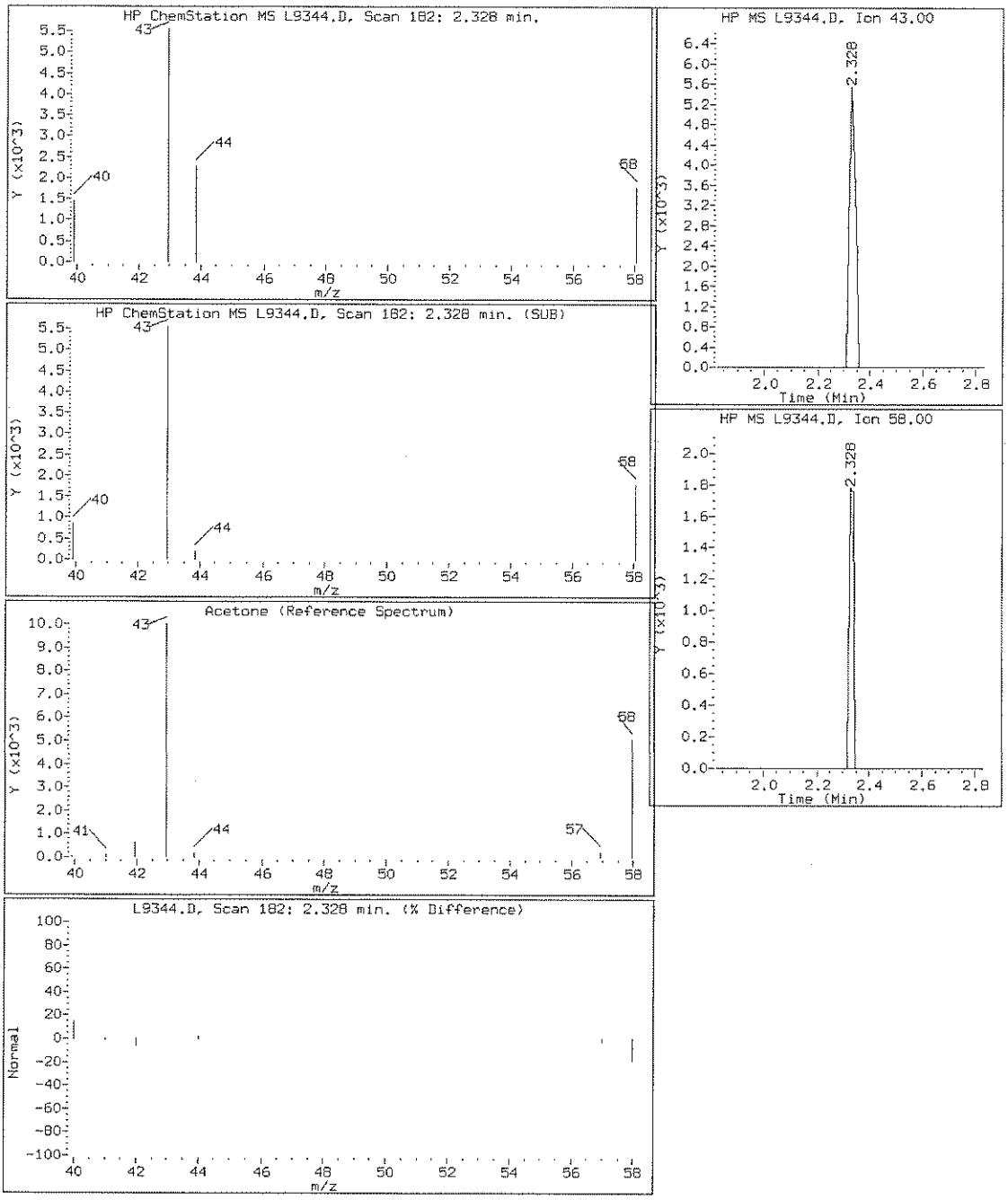
Client ID: MB

Instrument: msl.i

Sample Info: MB

Operator: D. HUMBERT

21 Acetone



Data File: L9344.D

Date: 01-AUG-2007 11:41

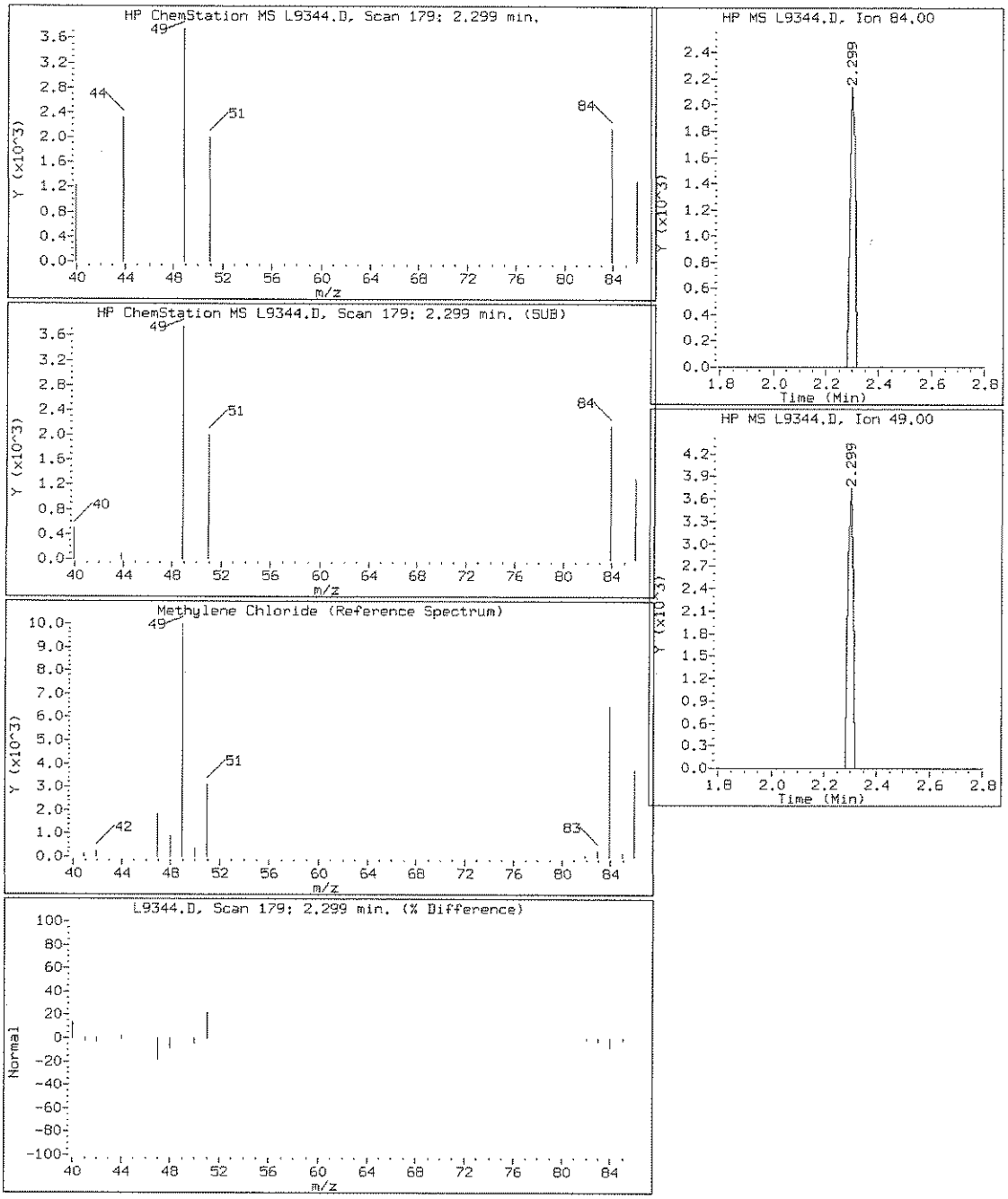
Client ID: MB

Instrument: msl.i

Sample Info: MB

Operator: D. HUMBERT

20 Methylene Chloride



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: | Lab Sample ID: MB 220-8356/4

Matrix: Water | Lab File ID: L9375.D

Analysis Method: 8260B | Date Received:

Sample wt/vol: 5 (mL) | Date Analyzed: 08/02/2007 11:30

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8356 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	10	U	10	1.4
71-43-2	Benzene	5.0	U	5.0	0.40
75-27-4	Bromodichloromethane	5.0	U	5.0	0.40
75-25-2	Bromoform	5.0	U	5.0	0.80
74-83-9	Bromomethane	5.0	U	5.0	1.2
78-93-3	2-Butanone (MEK)	10	U	10	1.2
75-15-0	Carbon disulfide	5.0	U	5.0	0.90
56-23-5	Carbon tetrachloride	5.0	U	5.0	1.0
108-90-7	Chlorobenzene	5.0	U	5.0	0.40
75-00-3	Chloroethane	5.0	U	5.0	0.80
67-66-3	Chloroform	5.0	U	5.0	0.70
74-87-3	Chloromethane	5.0	U	5.0	0.50
124-48-1	Dibromochloromethane	5.0	U	5.0	0.50
75-34-3	1,1-Dichloroethane	5.0	U	5.0	0.60
107-06-2	1,2-Dichloroethane	5.0	U	5.0	0.60
75-35-4	1,1-Dichloroethene	5.0	U	5.0	0.70
78-87-5	1,2-Dichloropropane	5.0	U	5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	5.0	U	5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	5.0	U	5.0	0.80
100-41-4	Ethylbenzene	5.0	U	5.0	1.0
591-78-6	2-Hexanone	10	U	10	0.80
75-09-2	Methylene Chloride	5.0	U	5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	10	U	10	0.70
100-42-5	Styrene	5.0	U	5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	5.0	0.40
127-18-4	Tetrachloroethene	5.0	U	5.0	0.50
108-88-3	Toluene	5.0	U	5.0	0.30
71-55-6	1,1,1-Trichloroethane	5.0	U	5.0	0.40
79-00-5	1,1,2-Trichloroethane	5.0	U	5.0	0.60
79-01-6	Trichloroethene	5.0	U	5.0	0.70
75-01-4	Vinyl chloride	5.0	U	5.0	0.80
1330-20-7	Xylenes, Total	5.0	U	5.0	1.0
156-59-2	cis-1,2-Dichloroethene	5.0	U	5.0	0.60
156-60-5	trans-1,2-Dichloroethene	5.0	U	5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079370.b\L9375.D
Lab Smp Id: MB Client Smp ID: MB
Inj Date : 02-AUG-2007 11:30 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : MB
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\target1_ct\Files\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 11:58 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 58 QC Sample: BLANK
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONMSW

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1 Fluorobenzene	96	4.895	4.901	(1.000)	438738	25.0000	
$ 41 Dibromofluoromethane	111	3.912	3.927	(0.799)	103827	18.3435	18
$ 55 1,2-Dichloroethane-d4	65	4.561	4.567	(0.932)	128164	18.5724	18
* 75 Chlorobenzene-d5	117	7.955	7.961	(1.000)	427020	25.0000	
$ 77 Toluene-d8	98	6.529	6.535	(0.821)	305247	20.1035	20
* 95 1,4-Dichlorobenzene-d4	152	10.012	10.018	(1.000)	128132	25.0000	
$ 125 Bromofluorobenzene	95	9.038	9.043	(0.903)	131006	27.2534	27

Data File: L9375.D

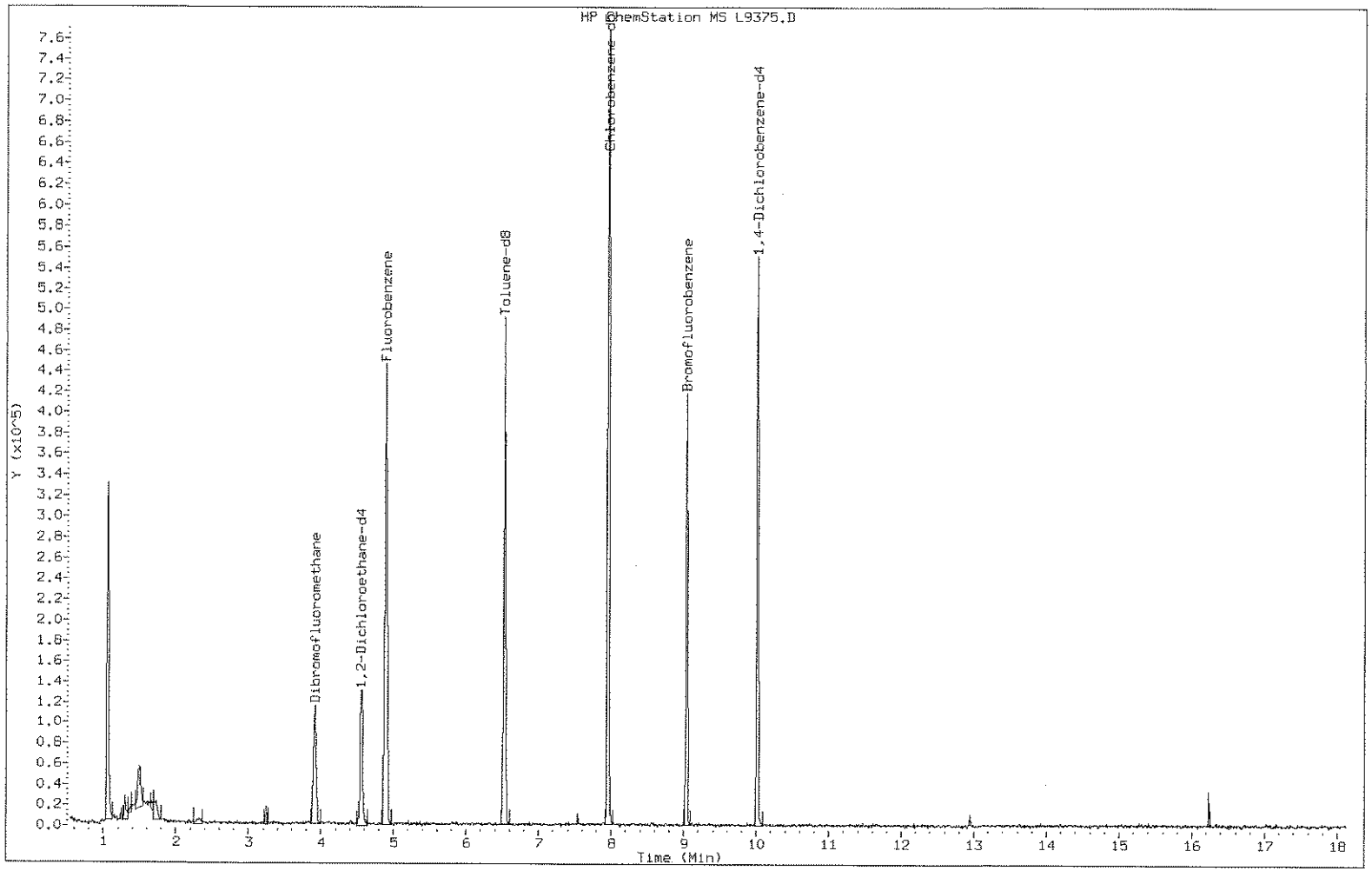
Date: 02-AUG-2007 11:30

Client ID: MB

Instrument: msl.i

Sample Info: MB

Operator: D. HUMBERT



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: MW-112 MS | Lab Sample ID: 220-2277-22 MS

Matrix: Water | Lab File ID: L9393.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/02/2007 19:42

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8356 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	39.8		10	1.4
71-43-2	Benzene	49.4		5.0	0.40
75-27-4	Bromodichloromethane	47.8		5.0	0.40
75-25-2	Bromoform	44.8		5.0	0.80
74-83-9	Bromomethane	41.7		5.0	1.2
78-93-3	2-Butanone (MEK)	50.3		10	1.2
75-15-0	Carbon disulfide	45.2		5.0	0.90
56-23-5	Carbon tetrachloride	46.1		5.0	1.0
108-90-7	Chlorobenzene	49.3		5.0	0.40
75-00-3	Chloroethane	60.9		5.0	0.80
67-66-3	Chloroform	49.1		5.0	0.70
74-87-3	Chloromethane	46.2		5.0	0.50
124-48-1	Dibromochloromethane	44.7		5.0	0.50
75-34-3	1,1-Dichloroethane	52.0		5.0	0.60
107-06-2	1,2-Dichloroethane	49.3		5.0	0.60
75-35-4	1,1-Dichloroethene	52.2		5.0	0.70
78-87-5	1,2-Dichloropropane	50.6		5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	46.9		5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	46.9		5.0	0.80
100-41-4	Ethylbenzene	49.4		5.0	1.0
591-78-6	2-Hexanone	48.9		10	0.80
75-09-2	Methylene Chloride	49.3		5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	50.5		10	0.70
100-42-5	Styrene	48.3		5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	49.2		5.0	0.40
127-18-4	Tetrachloroethene	49.9		5.0	0.50
108-88-3	Toluene	49.7		5.0	0.30
71-55-6	1,1,1-Trichloroethane	50.2		5.0	0.40
79-00-5	1,1,2-Trichloroethane	51.8		5.0	0.60
79-01-6	Trichloroethene	49.0		5.0	0.70
75-01-4	Vinyl chloride	49.8		5.0	0.80
1330-20-7	Xylenes, Total	149		5.0	1.0
156-59-2	cis-1,2-Dichloroethene	49.8		5.0	0.60
156-60-5	trans-1,2-Dichloroethene	49.7		5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079370.b\L9393.D
Lab Smp Id: 220-2277-C-22 MS
Inj Date : 02-AUG-2007 19:42 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-C-22 MS
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 75 QC Sample: MS
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONMSLNT

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1	Fluorobenzene	96	4.892	4.901	(1.000)	440538	25.0000	
2	Dichlorodifluoromethane	85	1.144	1.143	(0.234)	150001	44.3448	44
3	Chloromethane	50	1.262	1.261	(0.258)	273614	46.1775	46
4	Vinyl Chloride	62	1.301	1.300	(0.266)	324796	49.7965	50
5	Bromomethane	94	1.478	1.477	(0.302)	94971	41.6946	42
6	Chloroethane	64	1.537	1.546	(0.314)	210519	60.8982	61
7	Trichlorofluoromethane	101	1.626	1.625	(0.332)	313750	49.6803	50
8	Dichlorofluoromethane	67	1.645	1.645	(0.336)	597730	51.8641	52
9	Ethyl Ether	45	1.793	1.792	(0.367)	202349	50.9261	51
10	Ethanol	45	1.862	1.861	(0.381)	158794	507.126	510
11	Freon 141	81	1.852	1.861	(0.379)	387587	49.7436	50
12	Freon 123a	67	1.645	1.645	(0.336)	597730	51.8641	52
13	Trichlorotrifluoroethane	101	1.941	1.950	(0.397)	226016	51.8272	52
14	1,1-Dichloroethene	96	1.931	1.930	(0.395)	195834	52.1589	52
15	Carbon Disulfide	76	1.970	1.969	(0.403)	829154	45.2418	45
16	Iodomethane	142	2.029	2.038	(0.415)	250511	51.4404	51
17	Acrolein	56	2.127	2.127	(0.435)	360229	352.255	350
18	2-Propanol	45	2.216	2.215	(0.453)	55236	54.6389	55
19	3-Chloro-1-Propene	41	2.226	2.225	(0.455)	465054	47.9889	48
20	Methylene Chloride	84	2.295	2.304	(0.469)	254572	49.3226	49
21	Acetone	43	2.324	2.323	(0.475)	111293	39.8353	40
22	trans-1,2-Dichloroethene	96	2.413	2.422	(0.493)	234385	49.7063	50

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
	23 Methyl Acetate	43	2.403	2.402	(0.491)	1566397	47.5763	48
	24 Methyl tert-Butyl Ether	73	2.482	2.491	(0.507)	905181	49.4904	49
	25 tert-Butyl alcohol	59	2.531	2.530	(0.517)	268024	262.680	260
	26 Acetonitrile	41	2.659	2.658	(0.543)	616985	529.396	530
	27 Isopropyl ether	45	2.777	2.786	(0.568)	1078103	50.6728	51
	28 tert-Butyl ethyl ether	59	3.111	3.111	(0.636)	1083762	50.6956	51
	29 2-Chloro-1,3-Butadiene	88	2.885	2.884	(0.590)	169392	49.6791	50
	30 Acrylonitrile	53	2.915	2.924	(0.596)	314692	97.3711	97
	31 1,1-Dichloroethane	63	2.895	2.904	(0.592)	625681	51.9915	52
	32 Vinyl Acetate	43	3.102	3.111	(0.634)	795094	49.5091	50
	33 cis-1,2-Dichloroethene	96	3.407	3.416	(0.696)	250870	49.7885	50
	34 2,2-Dichloropropane	77	3.515	3.524	(0.718)	371345	44.7357	45
	35 Bromochloromethane	128	3.623	3.622	(0.741)	176585	48.9499	49
	36 1-Bromopropane	43	3.613	3.612	(0.739)	473919	52.2207	52
	37 Cyclohexane	84	3.653	3.652	(0.747)	323461	50.2673	50
	38 Chloroform	83	3.702	3.701	(0.757)	452112	49.1147	49
	39 Ethyl Acetate	43	3.849	3.858	(0.787)	50553	102.610	100
	40 Methyl Acrylate	55	3.849	3.858	(0.787)	365425	51.8028	52
$	41 Dibromofluoromethane	111	3.918	3.927	(0.801)	108023	19.0069	19
	42 Tetrahydrofuran	42	3.908	3.917	(0.799)	249522	104.536	100
	43 Carbon Tetrachloride	117	3.889	3.898	(0.795)	349515	46.1155	46
	44 1,1,1-Trichloroethane	97	3.958	3.967	(0.809)	339975	50.2481	50
	45 2-Butanone	43	4.056	4.065	(0.829)	203960	50.2921	50
	46 1,1-Dichloropropene	75	4.115	4.124	(0.841)	420034	49.9059	50
	47 tert-Amyl methyl ether	73	4.568	4.567	(0.934)	914541	49.3144	49
	48 tert-Butyl formate	57	3.111	3.111	(0.636)	317752	51.5356	52
	49 1-Chlorobutane	56	4.174	4.173	(0.853)	615599	50.9661	51
	50 Heptane	43	4.390	4.399	(0.897)	242808	44.5913	44
	51 Propionitrile	54	4.390	4.399	(0.897)	576709	515.956	520
	52 Benzene	78	4.410	4.419	(0.901)	1045502	49.4374	49
	53 2-Methyl-2-Propenenitrile	41	4.430	4.439	(0.905)	399448	50.7256	51
	54 Isobutyl alcohol	42	4.676	4.685	(0.956)	127526	447.808	450
$	55 1,2-Dichloroethane-d4	65	4.558	4.567	(0.932)	127050	18.3357	18
	56 1,2-Dichloroethane	62	4.636	4.645	(0.948)	412618	49.3337	49
	59 Methyl Cyclohexane	83	5.089	5.098	(1.040)	291653	47.3089	47
	60 Trichloroethene	130	5.089	5.098	(1.040)	304696	48.9764	49
	61 Isopropyl Acetate	43	4.676	4.685	(0.956)	211897	89.6228	90
	62 N-Butanol	56	5.473	5.482	(1.119)	126035	549.279	550
	63 Dibromomethane	93	5.532	5.531	(1.131)	168625	49.7473	50
	64 1,2-Dichloropropane	63	5.630	5.629	(1.151)	352828	50.5617	50
	65 Bromodichloromethane	83	5.709	5.718	(1.167)	315366	47.8372	48
	66 Methyl Methacrylate	69	5.886	5.895	(1.203)	303831	99.3231	99
	67 1,4-Dioxane	58	5.935	5.934	(1.213)	28175	533.381	530
	68 N-Propyl Acetate	43	6.348	6.298	(1.298)	12693	22.7779	23
	69 2-Chloroethylvinylether	63	6.348	6.298	(1.298)	1091	0.64170	0.6(M)
	70 cis-1,3-Dichloropropene	75	6.348	6.348	(1.298)	468256	46.9176	47
	71 Chloroacetonitrile	48	6.693	6.692	(1.368)	138983	486.812	490
	72 2-Nitropropane	41	6.771	6.771	(1.384)	191362	97.6592	98
	73 trans-1,3-Dichloropropene	75	6.968	6.977	(1.424)	424572	46.8673	47
	74 1,1,2-Trichloroethane	97	7.116	7.125	(1.454)	228395	51.8025	52
*	75 Chlorobenzene-d5	117	7.952	7.961	(1.000)	438582	25.0000	
	76 Toluene	91	6.575	6.584	(0.827)	927993	49.6913	50
$	77 Toluene-d8	98	6.525	6.535	(0.821)	309453	19.8432	20
	78 1,1-Dichloro-2-propanone	43	6.791	6.800	(0.854)	1160099	253.312	250

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
79	4-Methyl-2-Pentanone	43	6.929	6.938	(0.871)	408059	50.5412	50
80	Tetrachloroethene	164	6.949	6.958	(0.874)	183831	49.8870	50
81	Ethyl Methacrylate	69	7.145	7.145	(0.899)	479547	48.2403	48
82	Dibromochloromethane	129	7.283	7.292	(0.916)	313805	44.7204	45
83	1,3-Dichloropropane	76	7.362	7.361	(0.926)	482370	47.6233	48
84	1,2-Dibromoethane	107	7.480	7.489	(0.941)	295227	49.9966	50
85	n-Butyl Acetate	56	7.637	7.646	(0.960)	241101	47.3712	47
86	2-Hexanone	43	7.706	7.705	(0.969)	294468	48.9051	49
87	1-Chlorohexane	91	7.962	7.971	(1.001)	309020	48.8563	49
88	Chlorobenzene	112	7.972	7.971	(1.002)	748181	49.3161	49
89	1,1,1,2-Tetrachloroethane	131	8.031	8.040	(1.010)	259073	47.2171	47
90	Ethylbenzene	106	8.001	8.010	(1.006)	333340	49.3594	49
91	Xylene (total)mp	106	8.139	8.138	(1.024)	832581	98.6344	99
92	Xylene (total)o	106	8.513	8.512	(1.071)	404854	50.0687	50
93	Styrene	104	8.562	8.561	(1.077)	654115	48.2901	48
94	Bromoform	173	8.582	8.581	(1.079)	173290	44.7601	45
* 95	1,4-Dichlorobenzene-d4	152	10.008	10.018	(1.000)	160374	25.0000	
96	Isopropylbenzene	105	8.798	8.797	(0.879)	975973	49.3482	49
97	Bromobenzene	156	9.123	9.122	(0.912)	241853	48.4306	48
98	1,1,2,2-Tetrachloroethane	83	9.221	9.221	(0.921)	335540	49.1666	49
99	4-Ethyltoluene	105	9.251	9.260	(0.924)	935360	48.3834	48
100	1,2,3-Trichloropropane	110	9.330	9.329	(0.932)	107313	49.6312	50
101	trans-1,4-Dichloro-2-Butene	53	9.369	9.368	(0.936)	111405	50.1706	50
102	n-Propylbenzene	91	9.153	9.162	(0.914)	1008986	47.6065	48
103	2-Chlorotoluene	91	9.290	9.289	(0.928)	741793	50.7812	51
104	4-Chlorotoluene	91	9.428	9.437	(0.942)	672278	48.8195	49
105	1,3,5-Trimethylbenzene	105	9.330	9.339	(0.932)	782948	49.4254	49
106	tert-Butylbenzene	119	9.605	9.604	(0.960)	725540	47.3549	47
107	1,2,4-Trimethylbenzene	105	9.664	9.673	(0.966)	777296	48.8495	49
108	sec-Butylbenzene	105	9.763	9.762	(0.975)	781920	46.6973	47
109	4-Isopropyltoluene	119	9.890	9.890	(0.988)	848473	47.3309	47
110	1,3-Dichlorobenzene	146	9.949	9.949	(0.994)	414338	47.4352	47
111	1,4-Dichlorobenzene	146	10.028	10.027	(1.002)	423313	47.4833	47
112	1,2-Dichlorobenzene	146	10.382	10.391	(1.037)	415498	48.3354	48
113	Benzyl Chloride	126	10.235	10.234	(1.023)	116442	43.8106	44
114	1,4-Diethylbenzene	119	10.205	10.204	(2.086)	464317	48.3887	48
115	n-Butylbenzene	91	10.254	10.254	(1.025)	964206	44.6388	45
118	1,2,4,5-Tetramethylbenzene	119	10.914	10.913	(2.231)	727966	49.8633	50
119	1,2-Dibromo-3-chloropropane	75	11.081	11.080	(1.107)	66116	48.5227	48
120	Nitrobenzene	77	11.563	11.572	(1.155)	71793	236.654	240
121	1,2,4-Trichlorobenzene	180	11.681	11.690	(1.167)	229516	49.1458	49
122	Hexachlorobutadiene	225	11.671	11.670	(1.166)	68554	43.0809	43
123	Naphthalene	128	11.966	11.966	(1.196)	694366	49.0496	49
124	1,2,3-Trichlorobenzene	180	12.134	12.133	(1.212)	207287	48.6459	49
$ 125	Bromofluorobenzene	95	9.034	9.043	(0.903)	140862	23.4125	23
M 126	1,2-Dichloroethene (total)	100				485255	99.4948	99
M 127	Xylene (total)	100				1237435	148.703	150

QC Flag Legend

M - Compound response manually integrated.

Data File: L9393.D

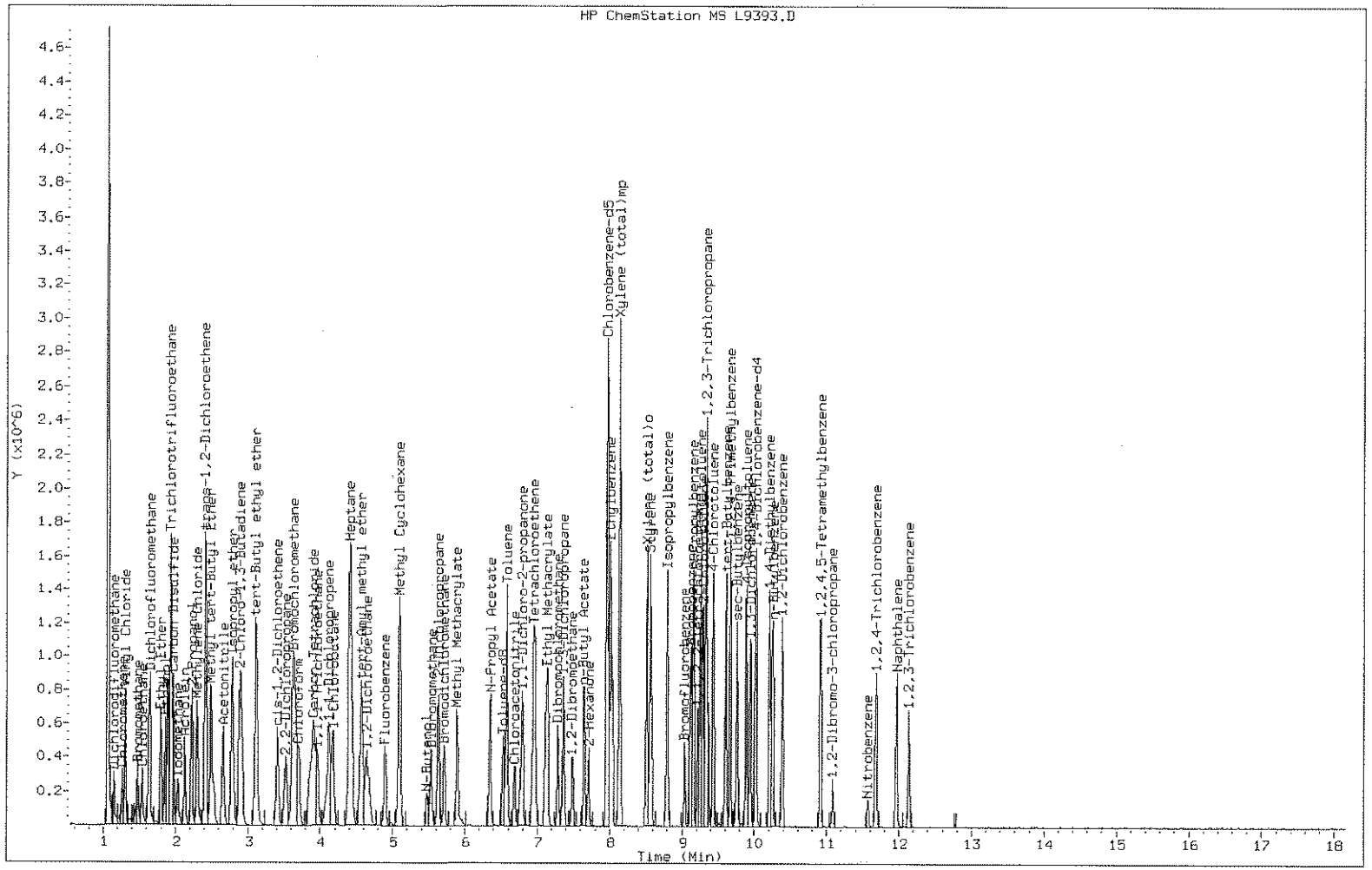
Date: 02-AUG-2007 19:42

Client ID:

Instrument: msl.i

Sample Info: 220-2277-C-22 MS

Operator: D. HUMBERT



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: MW-112 MSD | Lab Sample ID: 220-2277-22 MSD

Matrix: Water | Lab File ID: L9394.D

Analysis Method: 8260B | Date Received: 07/28/2007 10:00

Sample wt/vol: 5 (mL) | Date Analyzed: 08/02/2007 20:07

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8356 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	44.4		10	1.4
71-43-2	Benzene	50.8		5.0	0.40
75-27-4	Bromodichloromethane	49.3		5.0	0.40
75-25-2	Bromoform	47.5		5.0	0.80
74-83-9	Bromomethane	47.2		5.0	1.2
78-93-3	2-Butanone (MEK)	49.4		10	1.2
75-15-0	Carbon disulfide	47.2		5.0	0.90
56-23-5	Carbon tetrachloride	47.8		5.0	1.0
108-90-7	Chlorobenzene	51.3		5.0	0.40
75-00-3	Chloroethane	60.3		5.0	0.80
67-66-3	Chloroform	50.7		5.0	0.70
74-87-3	Chloromethane	48.2		5.0	0.50
124-48-1	Dibromochloromethane	47.2		5.0	0.50
75-34-3	1,1-Dichloroethane	51.6		5.0	0.60
107-06-2	1,2-Dichloroethane	52.1		5.0	0.60
75-35-4	1,1-Dichloroethene	52.6		5.0	0.70
78-87-5	1,2-Dichloropropane	51.6		5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	48.2		5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	48.1		5.0	0.80
100-41-4	Ethylbenzene	52.4		5.0	1.0
591-78-6	2-Hexanone	51.7		10	0.80
75-09-2	Methylene Chloride	50.6		5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	54.4		10	0.70
100-42-5	Styrene	51.7		5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	52.8		5.0	0.40
127-18-4	Tetrachloroethene	56.1		5.0	0.50
108-88-3	Toluene	51.3		5.0	0.30
71-55-6	1,1,1-Trichloroethane	51.4		5.0	0.40
79-00-5	1,1,2-Trichloroethane	52.1		5.0	0.60
79-01-6	Trichloroethene	50.4		5.0	0.70
75-01-4	Vinyl chloride	49.9		5.0	0.80
1330-20-7	Xylenes, Total	156		5.0	1.0
156-59-2	cis-1,2-Dichloroethene	52.3		5.0	0.60
156-60-5	trans-1,2-Dichloroethene	51.2		5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079370.b\L9394.D
Lab Smp Id: 220-2277-C-22 MSD
Inj Date : 02-AUG-2007 20:07 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : 220-2277-C-22 MSD
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 76 QC Sample: MSD
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONMSLNT

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1	Fluorobenzene	96	4.891	4.901	(1.000)	455584	25.0000	
2	Dichlorodifluoromethane	85	1.142	1.143	(0.234)	156701	44.7956	45
3	Chloromethane	50	1.260	1.261	(0.258)	295618	48.2434	48
4	Vinyl Chloride	62	1.300	1.300	(0.266)	336907	49.9475	50
5	Bromomethane	94	1.477	1.477	(0.302)	111088	47.1597	47
6	Chloroethane	64	1.536	1.546	(0.314)	215533	60.2896	60
7	Trichlorofluoromethane	101	1.625	1.625	(0.332)	327239	50.1049	50
8	Dichlorofluoromethane	67	1.644	1.645	(0.336)	616637	51.7376	52
9	Ethyl Ether	45	1.792	1.792	(0.366)	215344	52.4067	52
10	Ethanol	45	1.861	1.861	(0.380)	176553	545.220	540
11	Freon 141	81	1.861	1.861	(0.380)	411929	51.1217	51
12	Freon 123a	67	1.644	1.645	(0.336)	616637	51.7376	52
13	Trichlorotrifluoroethane	101	1.939	1.950	(0.397)	241798	53.6150	54
14	1,1-Dichloroethene	96	1.930	1.930	(0.395)	204427	52.6494	53
15	Carbon Disulfide	76	1.969	1.969	(0.403)	894858	47.2143	47
16	Iodomethane	142	2.028	2.038	(0.415)	306820	60.9222	61
17	Acrolein	56	2.126	2.127	(0.435)	384217	363.304	360
18	2-Propanol	45	2.215	2.215	(0.453)	60398	57.7720	58
19	3-Chloro-1-Propene	41	2.225	2.225	(0.455)	486278	48.5218	48
20	Methylene Chloride	84	2.294	2.304	(0.469)	270330	50.6460	51
21	Acetone	43	2.323	2.323	(0.475)	128424	44.4489	44
22	trans-1,2-Dichloroethene	96	2.421	2.422	(0.495)	249740	51.2135	51

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
23	Methyl Acetate	43	2.402	2.402	(0.491)	1669702	49.0392	49
24	Methyl tert-Butyl Ether	73	2.490	2.491	(0.509)	963436	50.9358	51
25	tert-Butyl alcohol	59	2.530	2.530	(0.517)	288354	273.271	270
26	Acetonitrile	41	2.658	2.658	(0.543)	664138	551.035	550
27	Isopropyl ether	45	2.776	2.786	(0.568)	1137591	51.7030	52
28	tert-Butyl ethyl ether	59	3.110	3.111	(0.636)	1154435	52.2181	52
29	2-Chloro-1,3-Butadiene	88	2.884	2.884	(0.590)	180974	51.3230	51
30	Acrylonitrile	53	2.913	2.924	(0.596)	344752	103.149	100
31	1,1-Dichloroethane	63	2.894	2.904	(0.592)	642671	51.6396	52
32	Vinyl Acetate	43	3.100	3.111	(0.634)	851409	51.2649	51
33	cis-1,2-Dichloroethene	96	3.405	3.416	(0.696)	272387	52.2735	52
34	2,2-Dichloropropane	77	3.523	3.524	(0.720)	391997	45.6640	46
35	Bromochloromethane	128	3.622	3.622	(0.741)	190406	51.0379	51
36	1-Bromopropane	43	3.612	3.612	(0.739)	494031	52.6390	53
37	Cyclohexane	84	3.651	3.652	(0.747)	346266	52.0342	52
38	Chloroform	83	3.701	3.701	(0.757)	482476	50.6823	51
39	Ethyl Acetate	43	3.848	3.858	(0.787)	51369	100.823	100
40	Methyl Acrylate	55	3.848	3.858	(0.787)	388963	53.3186	53
$ 41	Dibromofluoromethane	111	3.907	3.927	(0.799)	110463	18.7943	19
42	Tetrahydrofuran	42	3.907	3.917	(0.799)	269190	109.051	110
43	Carbon Tetrachloride	117	3.887	3.898	(0.795)	374976	47.8409	48
44	1,1,1-Trichloroethane	97	3.956	3.967	(0.809)	359842	51.4279	51
45	2-Butanone	43	4.055	4.065	(0.829)	207041	49.3658	49
46	1,1-Dichloropropene	75	4.114	4.124	(0.841)	448228	51.4970	51
47	tert-Amyl methyl ether	73	4.566	4.567	(0.934)	968548	50.5018	50
48	tert-Butyl formate	57	3.110	3.111	(0.636)	338799	53.1344	53
49	1-Chlorobutane	56	4.173	4.173	(0.853)	653808	52.3418	52
50	Heptane	43	4.399	4.399	(0.899)	258124	45.8385	46
51	Propionitrile	54	4.389	4.399	(0.897)	618452	535.028	540
52	Benzene	78	4.409	4.419	(0.901)	1110688	50.7853	51
53	2-Methyl-2-Propenenitrile	41	4.429	4.439	(0.905)	424429	52.1179	52
54	Isobutyl alcohol	42	4.675	4.685	(0.956)	135751	460.947	460
$ 55	1,2-Dichloroethane-d4	65	4.557	4.567	(0.932)	130372	18.1938	18
56	1,2-Dichloroethane	62	4.635	4.645	(0.948)	450376	52.0698	52
59	Methyl Cyclohexane	83	5.088	5.098	(1.040)	320397	50.2550	50
60	Trichloroethene	130	5.098	5.098	(1.042)	324032	50.3643	50
61	Isopropyl Acetate	43	4.675	4.685	(0.956)	228853	93.5977	94
62	N-Butanol	56	5.472	5.482	(1.119)	138969	585.645	580
63	Dibromomethane	93	5.531	5.531	(1.131)	181163	51.6811	52
64	1,2-Dichloropropane	63	5.629	5.629	(1.151)	372283	51.5878	52
65	Bromodichloromethane	83	5.708	5.718	(1.167)	336210	49.3147	49
66	Methyl Methacrylate	69	5.885	5.895	(1.203)	323169	102.156	100
67	1,4-Dioxane	58	5.924	5.934	(1.211)	31519	576.980	580
68	N-Propyl Acetate	43	6.347	6.298	(1.298)	13510	23.4433	23
69	2-Chloroethylvinylether	63	6.347	6.298	(1.298)	3213	1.82740	2(M)
70	cis-1,3-Dichloropropene	75	6.347	6.348	(1.298)	497412	48.1929	48
71	Chloroacetonitrile	48	6.692	6.692	(1.368)	145601	493.150	490
72	2-Nitropropane	41	6.770	6.771	(1.384)	205901	101.609	100
73	trans-1,3-Dichloropropene	75	6.967	6.977	(1.424)	450568	48.0943	48
74	1,1,2-Trichloroethane	97	7.115	7.125	(1.455)	237360	52.0579	52
* 75	Chlorobenzene-d5	117	7.951	7.961	(1.000)	444075	25.0000	
76	Toluene	91	6.574	6.584	(0.827)	970690	51.3347	51
$ 77	Toluene-d8	98	6.524	6.535	(0.821)	313316	19.8424	20
78	1,1-Dichloro-2-propanone	43	6.790	6.800	(0.854)	1239840	267.375	270

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
79	4-Methyl-2-Pentanone	43	6.938	6.938	(0.873)	444921	54.4252	54
80	Tetrachloroethene	164	6.947	6.958	(0.874)	209349	56.1092	56
81	Ethyl Methacrylate	69	7.144	7.145	(0.899)	517214	51.3858	51
82	Dibromochloromethane	129	7.282	7.292	(0.916)	335425	47.2102	47
83	1,3-Dichloropropane	76	7.361	7.361	(0.926)	518126	50.5207	50
84	1,2-Dibromoethane	107	7.489	7.489	(0.942)	306028	51.1847	51
85	n-Butyl Acetate	56	7.646	7.646	(0.962)	255925	49.6618	50
86	2-Hexanone	43	7.705	7.705	(0.969)	315059	51.6776	52
87	1-Chlorohexane	91	7.961	7.971	(1.001)	334311	52.2010	52
88	Chlorobenzene	112	7.971	7.971	(1.002)	788125	51.3064	51
89	1,1,1,2-Tetrachloroethane	131	8.030	8.040	(1.010)	272611	49.0699	49
90	Ethylbenzene	106	8.000	8.010	(1.006)	358359	52.4077	52
91	Xylene (total)mp	106	8.138	8.138	(1.024)	886384	103.709	100
92	Xylene (total)o	106	8.512	8.512	(1.071)	430018	52.5230	52
93	Styrene	104	8.561	8.561	(1.077)	709079	51.7004	52
94	Bromoform	173	8.581	8.581	(1.079)	186117	47.4786	47
* 95	1,4-Dichlorobenzene-d4	152	10.007	10.018	(1.000)	164899	25.0000	
96	Isopropylbenzene	105	8.797	8.797	(0.879)	1013774	49.8529	50
97	Bromobenzene	156	9.122	9.122	(0.912)	254092	49.4852	49
98	1,1,2,2-Tetrachloroethane	83	9.220	9.221	(0.921)	370476	52.7961	53
99	4-Ethyltoluene	105	9.250	9.260	(0.924)	1000998	50.3578	50
100	1,2,3-Trichloropropane	110	9.328	9.329	(0.932)	117092	52.6678	53
101	trans-1,4-Dichloro-2-Butene	53	9.368	9.368	(0.936)	129311	56.6365	57
102	n-Propylbenzene	91	9.161	9.162	(0.915)	1085642	49.8177	50
103	2-Chlorotoluene	91	9.289	9.289	(0.928)	800476	53.2947	53
104	4-Chlorotoluene	91	9.427	9.437	(0.942)	721940	50.9872	51
105	1,3,5-Trimethylbenzene	105	9.328	9.339	(0.932)	825773	50.6983	51
106	tert-Butylbenzene	119	9.604	9.604	(0.960)	756259	48.0054	48
107	1,2,4-Trimethylbenzene	105	9.663	9.673	(0.966)	818689	50.0390	50
108	sec-Butylbenzene	105	9.761	9.762	(0.975)	852832	49.5346	50
109	4-Isopropyltoluene	119	9.889	9.890	(0.988)	888841	48.2222	48
110	1,3-Dichlorobenzene	146	9.948	9.949	(0.994)	443890	49.4239	49
111	1,4-Dichlorobenzene	146	10.027	10.027	(1.002)	455461	49.6874	50
112	1,2-Dichlorobenzene	146	10.381	10.391	(1.037)	435372	49.2576	49
113	Benzyl Chloride	126	10.234	10.234	(1.023)	121512	44.4636	44
114	1,4-Diethylbenzene	119	10.204	10.204	(2.086)	508206	51.2134	51
115	n-Butylbenzene	91	10.253	10.254	(1.025)	1124903	50.6494	51
118	1,2,4,5-Tetramethylbenzene	119	10.913	10.913	(2.231)	772939	51.1953	51
119	1,2-Dibromo-3-chloropropane	75	11.080	11.080	(1.107)	67646	48.2833	48
120	Nitrobenzene	77	11.572	11.572	(1.156)	84756	271.718	270
121	1,2,4-Trichlorobenzene	180	11.690	11.690	(1.168)	247682	51.5803	52
122	Hexachlorobutadiene	225	11.670	11.670	(1.166)	78393	47.9120	48
123	Naphthalene	128	11.965	11.966	(1.196)	751392	51.6214	52
124	1,2,3-Trichlorobenzene	180	12.133	12.133	(1.212)	224000	51.1256	51
$ 125	Bromofluorobenzene	95	9.033	9.043	(0.903)	145032	23.4441	23
M 126	1,2-Dichloroethene (total)	100				522127	103.487	100
M 127	Xylene (total)	100				1316402	156.232	160

QC Flag Legend

M - Compound response manually integrated.

Data File: L9394.D

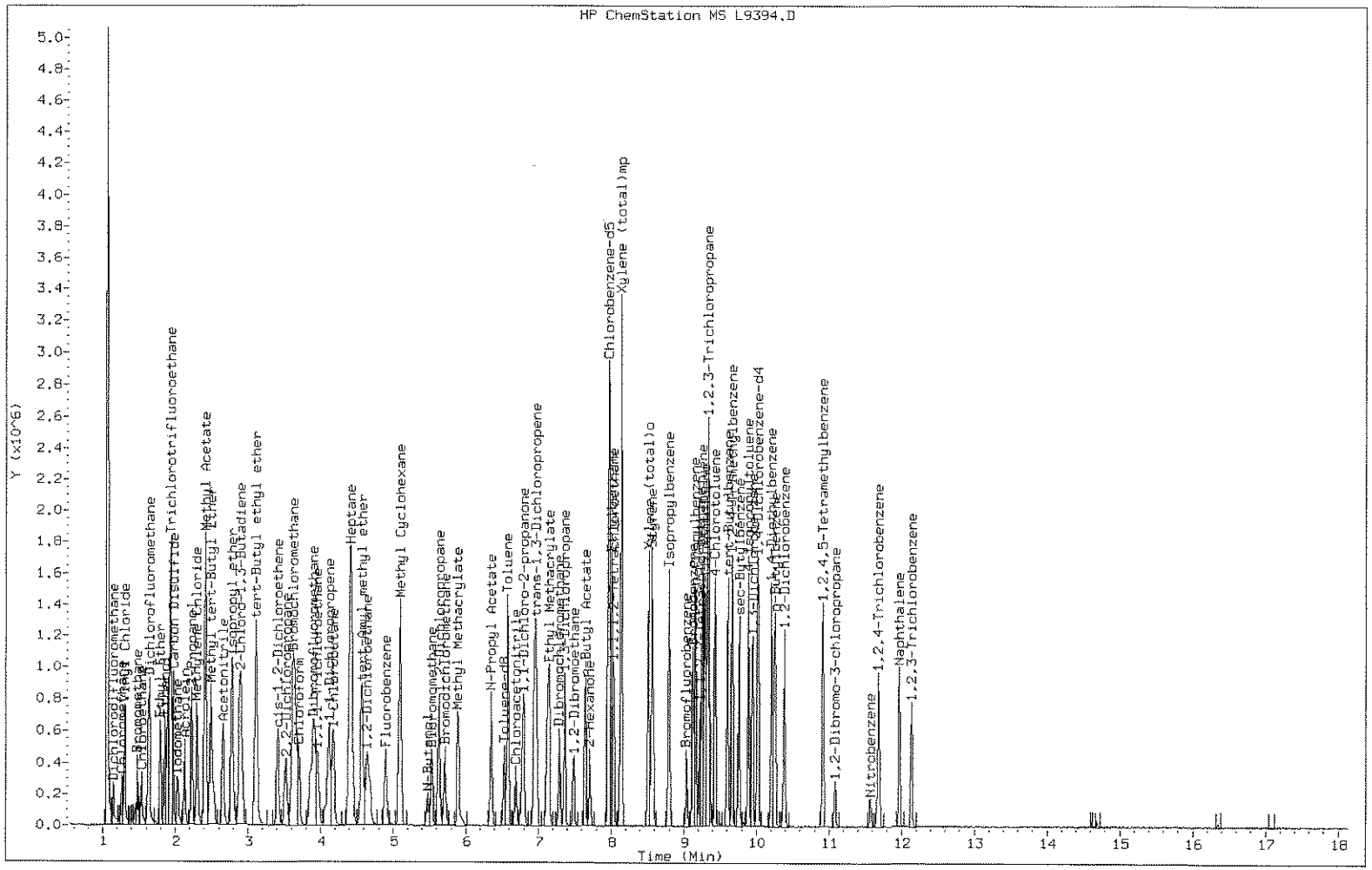
Date: 02-AUG-2007 20:07

Client ID:

Instrument: msl.i

Sample Info: 220-2277-C-22 MSD

Operator: D. HUMBERT



FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: Lab Sample ID: LCS 220-8307/2

Matrix: Water Lab File ID: L9311.D

Analysis Method: 8260B Date Received:

Sample wt/vol: 5 (mL) Date Analyzed: 07/31/2007 20:46

Level: (low/med) Low Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) Soil Aliquot:

Soil Extract Vol.: % Moisture:

Analy. Batch No.: 8307 Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	32.5		10	1.4
71-43-2	Benzene	20.0		5.0	0.40
75-27-4	Bromodichloromethane	19.6		5.0	0.40
75-25-2	Bromoform	17.8		5.0	0.80
74-83-9	Bromomethane	22.3		5.0	1.2
78-93-3	2-Butanone (MEK)	24.9		10	1.2
75-15-0	Carbon disulfide	17.8		5.0	0.90
56-23-5	Carbon tetrachloride	21.7		5.0	1.0
108-90-7	Chlorobenzene	20.9		5.0	0.40
75-00-3	Chloroethane	27.3		5.0	0.80
67-66-3	Chloroform	21.1		5.0	0.70
74-87-3	Chloromethane	22.5		5.0	0.50
124-48-1	Dibromochloromethane	18.4		5.0	0.50
75-34-3	1,1-Dichloroethane	20.7		5.0	0.60
107-06-2	1,2-Dichloroethane	20.7		5.0	0.60
75-35-4	1,1-Dichloroethene	21.7		5.0	0.70
78-87-5	1,2-Dichloropropane	20.9		5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	19.1		5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	19.2		5.0	0.80
100-41-4	Ethylbenzene	21.0		5.0	1.0
591-78-6	2-Hexanone	22.2		10	0.80
75-09-2	Methylene Chloride	19.8		5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	19.7		10	0.70
100-42-5	Styrene	18.6		5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	20.1		5.0	0.40
127-18-4	Tetrachloroethene	20.4		5.0	0.50
108-88-3	Toluene	20.7		5.0	0.30
71-55-6	1,1,1-Trichloroethane	20.1		5.0	0.40
79-00-5	1,1,2-Trichloroethane	20.0		5.0	0.60
79-01-6	Trichloroethene	21.0		5.0	0.70
75-01-4	Vinyl chloride	22.3		5.0	0.80
1330-20-7	Xylenes, Total	61.9		5.0	1.0
156-59-2	cis-1,2-Dichloroethene	20.9		5.0	0.60
156-60-5	trans-1,2-Dichloroethene	20.3		5.0	0.50

STL-CT

Volatile Report SW-846 Method 8260B

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079310.b\L9311.D
Lab Smp Id: LCS
Inj Date : 31-JUL-2007 20:46 MS Autotune Date: 26-APR-2004 14:21
Operator : D. GAYDA Inst ID: msl.i
Smp Info : LCS
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079310.b\L8260BNW.m
Meth Date : 31-Jul-2007 21:02 ctvoa Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 2 QC Sample: LCS
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONMSLNT

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1	Fluorobenzene	96	4.902	4.896	(1.000)	440213	25.0000	
2	Dichlorodifluoromethane	85	1.153	1.147	(0.235)	89095	26.3586	26
3	Chloromethane	50	1.271	1.265	(0.259)	133185	22.4941	22
4	Vinyl Chloride	62	1.301	1.304	(0.265)	145260	22.2872	22
5	Bromomethane	94	1.478	1.482	(0.302)	50783	22.3115	22
6	Chloroethane	64	1.547	1.541	(0.316)	94373	27.3201	27
7	Trichlorofluoromethane	101	1.635	1.629	(0.334)	137903	21.8522	22
8	Dichlorofluoromethane	67	1.930	1.649	(0.394)	34260	2.97488	3
9	Ethyl Ether	45	1.802	1.796	(0.368)	76691	19.3154	19
10	Ethanol	45	1.861	1.865	(0.380)	47700	152.448	150
11	Freon 141	81	1.861	1.855	(0.380)	151526	19.4615	19
12	Freon 123a	67	1.930	1.649	(0.394)	34260	2.97488	3
13	Trichlorotrifluoroethane	101	1.950	1.944	(0.398)	92182	21.1537	21
14	1,1-Dichloroethene	96	1.930	1.934	(0.394)	81489	21.7200	22
15	Carbon Disulfide	76	1.970	1.973	(0.402)	326842	17.8469	18
16	Iodomethane	142	2.039	2.033	(0.416)	97388	20.0126	20
17	Acrolein	56	2.501	2.131	(0.510)	13876	13.5789	14
18	2-Propanol	45	2.403	2.219	(0.490)	3063	3.03212	3
19	3-Chloro-1-Propene	41	2.226	2.229	(0.454)	184632	19.0662	19
20	Methylene Chloride	84	2.304	2.298	(0.470)	102258	19.8268	20
21	Acetone	43	2.324	2.328	(0.474)	90817	32.5303	32
22	trans-1,2-Dichloroethene	96	2.422	2.416	(0.494)	95622	20.2936	20

			QUANT SIG					CONCENTRATIONS	
		Compounds	MASS	RT	EXP RT	REL RT	RESPONSE	ON-COLUMN (ug/L)	FINAL (ug/L)
	23	Methyl Acetate	43	2.412	2.406	(0.492)	306819	9.32593	9
	24	Methyl tert-Butyl Ether	73	2.491	2.495	(0.508)	347611	19.0195	19
	25	tert-Butyl alcohol	59	2.531	2.534	(0.516)	75844	74.3866	74
	26	Acetonitrile	41	2.678	2.662	(0.546)	8867	7.61383	8
	30	Acrylonitrile	53	2.924	2.918	(0.597)	106692	33.0367	33
	31	1,1-Dichloroethane	63	2.904	2.898	(0.593)	248689	20.6803	21
	33	cis-1,2-Dichloroethene	96	3.416	3.410	(0.697)	105446	20.9426	21
	34	2,2-Dichloropropane	77	3.534	3.528	(0.721)	168454	20.3085	20
	35	Bromochloromethane	128	3.632	3.626	(0.741)	72542	20.1237	20
	36	1-Bromopropane	43	3.662	3.617	(0.747)	21703	2.39320	2
	37	Cyclohexane	84	3.652	3.656	(0.745)	129986	20.2153	20
	38	Chloroform	83	3.701	3.705	(0.755)	194334	21.1269	21
	39	Ethyl Acetate	43	3.859	3.843	(0.787)	7987	16.2236	16
	40	Methyl Acrylate	55	3.859	3.853	(0.787)	133495	18.9383	19
$	41	Dibromofluoromethane	111	3.918	3.922	(0.799)	108409	19.0889	19
	42	Tetrahydrofuran	42	3.928	3.912	(0.801)	87010	36.4793	36
	43	Carbon Tetrachloride	117	3.898	3.892	(0.795)	164277	21.6909	22
	44	1,1,1-Trichloroethane	97	3.967	3.961	(0.809)	135691	20.0698	20
	45	2-Butanone	43	4.075	4.069	(0.831)	100783	24.8693	25
	46	1,1-Dichloropropene	75	4.115	4.118	(0.839)	168927	20.0857	20
	49	1-Chlorobutane	56	4.174	4.177	(0.851)	247222	20.4829	20
	50	Heptane	43	4.410	4.394	(0.900)	9778	1.79704	2
	51	Propionitrile	54	4.400	4.394	(0.898)	204246	182.865	180
	52	Benzene	78	4.420	4.414	(0.902)	422277	19.9824	20
	53	2-Methyl-2-Propenenitrile	41	4.439	4.433	(0.906)	104538	13.2850	13
	54	Isobutyl alcohol	42	4.705	4.679	(0.960)	2268	7.96997	8
$	55	1,2-Dichloroethane-d4	65	4.557	4.561	(0.930)	127824	18.4610	18
	56	1,2-Dichloroethane	62	4.646	4.640	(0.948)	172780	20.6733	21
	59	Methyl Cyclohexane	83	5.089	5.092	(1.038)	128262	20.8207	21
	60	Trichloroethene	130	5.098	5.092	(1.040)	130824	21.0440	21
	61	Isopropyl Acetate	43	4.410	4.679	(0.900)	9778	4.13870	4
	62	N-Butanol	56	5.502	5.476	(1.122)	4366	19.0417	19
	63	Dibromomethane	93	5.531	5.535	(1.128)	67302	19.8699	20
	64	1,2-Dichloropropane	63	5.630	5.634	(1.149)	145412	20.8535	21
	65	Bromodichloromethane	83	5.718	5.712	(1.167)	128909	19.5684	20
	66	Methyl Methacrylate	69	5.895	5.889	(1.203)	221590	72.4918	72
	68	N-Propyl Acetate	43	6.584	6.303	(1.343)	3397	6.10049	6
	70	cis-1,3-Dichloropropene	75	6.348	6.352	(1.295)	190131	19.0645	19
	71	Chloroacetonitrile	48	6.692	6.686	(1.365)	13440	47.1107	47
	72	2-Nitropropane	41	6.771	6.765	(1.381)	63644	32.5039	32
	73	trans-1,3-Dichloropropene	75	6.978	6.972	(1.423)	173803	19.1998	19
	74	1,1,2-Trichloroethane	97	7.115	7.119	(1.452)	88312	20.0449	20
*	75	Chlorobenzene-d5	117	7.962	7.956	(1.000)	428646	25.0000	
	76	Toluene	91	6.584	6.578	(0.827)	377516	20.6835	21
$	77	Toluene-d8	98	6.535	6.529	(0.821)	325810	21.3763	21
	78	1,1-Dichloro-2-propanone	43	6.801	6.795	(0.854)	346292	77.3669	77
	79	4-Methyl-2-Pentanone	43	6.938	6.932	(0.871)	155488	19.7048	20
	80	Tetrachloroethene	164	6.958	6.952	(0.874)	73486	20.4045	20
	81	Ethyl Methacrylate	69	7.145	7.149	(0.897)	192343	19.7973	20
	82	Dibromochloromethane	129	7.293	7.287	(0.916)	125947	18.3648	18
	83	1,3-Dichloropropane	76	7.361	7.365	(0.925)	199829	20.1860	20
	84	1,2-Dibromoethane	107	7.489	7.483	(0.941)	112228	19.4463	19
	85	n-Butyl Acetate	56	7.971	7.641	(1.001)	77840	15.6484	16
	86	2-Hexanone	43	7.716	7.710	(0.969)	130785	22.2242	22

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
87	1-Chlorohexane	91	7.971	7.965	(1.001)	117690	19.0382	19
88	Chlorobenzene	112	7.971	7.975	(1.001)	310562	20.9451	21
89	1,1,1,2-Tetrachloroethane	131	8.040	8.034	(1.010)	105093	19.5976	20
90	Ethylbenzene	106	8.011	8.005	(1.006)	138653	21.0070	21
91	Xylene (total)mp	106	8.139	8.143	(1.022)	339798	41.1884	41
92	Xylene (total)o	106	8.513	8.516	(1.069)	163897	20.7392	21
93	Styrene	104	8.562	8.566	(1.075)	246063	18.5867	18
94	Bromoform	173	8.581	8.585	(1.078)	67483	17.8346	18
* 95	1,4-Dichlorobenzene-d4	152	10.018	10.012	(1.000)	153272	25.0000	
96	Isopropylbenzene	105	8.798	8.792	(0.878)	393355	20.8108	21
97	Bromobenzene	156	9.123	9.126	(0.911)	97075	20.3398	20
98	1,1,2,2-Tetrachloroethane	83	9.221	9.225	(0.920)	131180	20.1124	20
99	4-Ethyltoluene	105	9.260	9.254	(0.924)	7345	0.39754	0.4
100	1,2,3-Trichloropropane	110	9.329	9.323	(0.931)	40137	19.4231	19
101	trans-1,4-Dichloro-2-Butene	53	9.369	9.363	(0.935)	52429	24.7052	25
102	n-Propylbenzene	91	9.162	9.156	(0.915)	426440	21.0528	21
103	2-Chlorotoluene	91	9.290	9.294	(0.927)	274710	19.6773	20
104	4-Chlorotoluene	91	9.428	9.431	(0.941)	278087	21.1298	21
105	1,3,5-Trimethylbenzene	105	9.329	9.333	(0.931)	314518	20.7746	21
106	tert-Butylbenzene	119	9.605	9.609	(0.959)	311341	21.2623	21
107	1,2,4-Trimethylbenzene	105	9.674	9.668	(0.966)	320742	21.0912	21
108	sec-Butylbenzene	105	9.762	9.766	(0.974)	355512	22.2154	22
109	4-Isopropyltoluene	119	9.890	9.894	(0.987)	364236	21.2599	21
110	1,3-Dichlorobenzene	146	9.949	9.953	(0.993)	166785	19.9790	20
111	1,4-Dichlorobenzene	146	10.028	10.032	(1.001)	171251	20.0994	20
112	1,2-Dichlorobenzene	146	10.392	10.386	(1.037)	166410	20.2557	20
113	Benzyl Chloride	126	10.234	10.238	(1.022)	41876	16.4857	16
114	1,4-Diethylbenzene	119	10.205	10.209	(2.082)	8178	0.85290	0.8
115	n-Butylbenzene	91	10.254	10.258	(1.024)	368449	17.8481	18
118	1,2,4,5-Tetramethylbenzene	119	10.913	10.917	(2.226)	13312	0.91250	0.9
119	1,2-Dibromo-3-chloropropane	75	11.081	11.084	(1.106)	21090	16.1952	16
120	Nitrobenzene	77	11.573	11.566	(1.155)	21885	75.4830	75
121	1,2,4-Trichlorobenzene	180	11.691	11.685	(1.167)	85624	19.1840	19
122	Hexachlorobutadiene	225	11.671	11.675	(1.165)	34520	22.6983	23
123	Naphthalene	128	11.966	11.970	(1.194)	266387	19.6893	20
124	1,2,3-Trichlorobenzene	180	12.133	12.137	(1.211)	75343	18.5007	18
$ 125	Bromofluorobenzene	95	9.044	9.038	(0.903)	145015	25.2196	25
M 126	1,2-Dichloroethene (total)	100				201068	41.2363	41
M 127	Xylene (total)	100				503695	61.9276	62

Data File: L9311.D

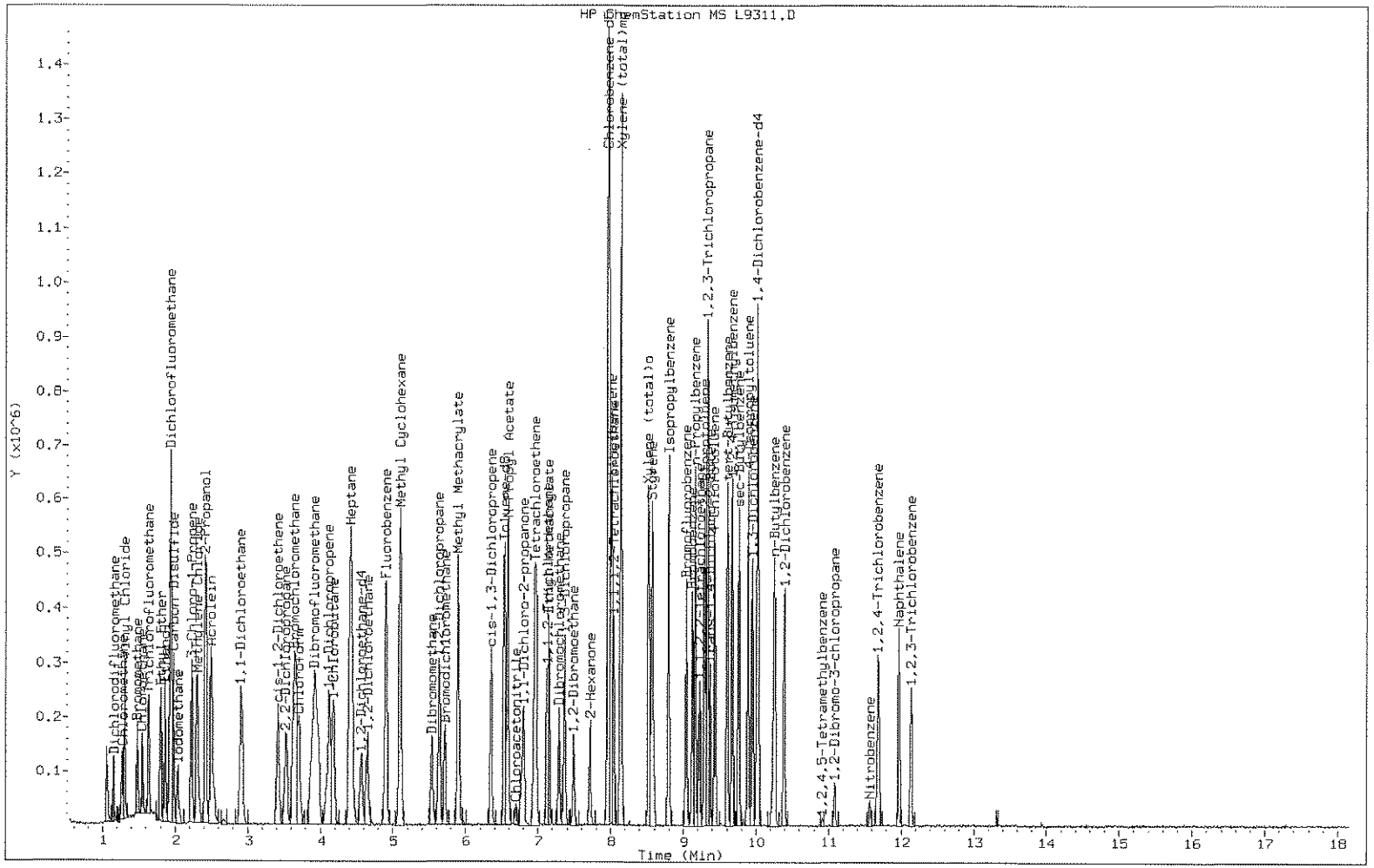
Date: 31-JUL-2007 20:46

Client ID: LCS

Instrument: msl.i

Sample Info: LCS

Operator: D. GAYDA



08/09/2007

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: | Lab Sample ID: LCS 220-8321/2

Matrix: Water | Lab File ID: L9342.D

Analysis Method: 8260B | Date Received:

Sample wt/vol: 5 (mL) | Date Analyzed: 08/01/2007 10:34

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8321 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	33.9		10	1.4
71-43-2	Benzene	20.3		5.0	0.40
75-27-4	Bromodichloromethane	19.3		5.0	0.40
75-25-2	Bromoform	18.1		5.0	0.80
74-83-9	Bromomethane	29.0		5.0	1.2
78-93-3	2-Butanone (MEK)	29.7		10	1.2
75-15-0	Carbon disulfide	18.6		5.0	0.90
56-23-5	Carbon tetrachloride	21.5		5.0	1.0
108-90-7	Chlorobenzene	20.2		5.0	0.40
75-00-3	Chloroethane	28.7		5.0	0.80
67-66-3	Chloroform	20.3		5.0	0.70
74-87-3	Chloromethane	24.7		5.0	0.50
124-48-1	Dibromochloromethane	18.9		5.0	0.50
75-34-3	1,1-Dichloroethane	20.7		5.0	0.60
107-06-2	1,2-Dichloroethane	21.1		5.0	0.60
75-35-4	1,1-Dichloroethene	22.9		5.0	0.70
78-87-5	1,2-Dichloropropane	21.6		5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	19.5		5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	19.1		5.0	0.80
100-41-4	Ethylbenzene	20.0		5.0	1.0
591-78-6	2-Hexanone	25.0		10	0.80
75-09-2	Methylene Chloride	19.3		5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	21.8		10	0.70
100-42-5	Styrene	18.7		5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	20.8		5.0	0.40
127-18-4	Tetrachloroethene	21.0		5.0	0.50
108-88-3	Toluene	20.0		5.0	0.30
71-55-6	1,1,1-Trichloroethane	20.7		5.0	0.40
79-00-5	1,1,2-Trichloroethane	20.9		5.0	0.60
79-01-6	Trichloroethene	20.7		5.0	0.70
75-01-4	Vinyl chloride	24.5		5.0	0.80
1330-20-7	Xylenes, Total	60.6		5.0	1.0
156-59-2	cis-1,2-Dichloroethene	20.6		5.0	0.60
156-60-5	trans-1,2-Dichloroethene	20.8		5.0	0.50

STL-CT

Volatile Report SW-846 Method 8260B

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079341.b\L9342.D
Lab Smp Id: LCS Client Smp ID: LCS
Inj Date : 01-AUG-2007 10:34 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : LCS
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079341.b\L8260BNW.m
Meth Date : 02-Aug-2007 09:12 target Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 31 QC Sample: LCS
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1	Fluorobenzene	96	4.906	4.908	(1.000)	450523	25.0000	
2	Dichlorodifluoromethane	85	1.147	1.149	(0.234)	116303	33.6207	34
3	Chloromethane	50	1.265	1.267	(0.258)	149768	24.7160	25
4	Vinyl Chloride	62	1.305	1.307	(0.266)	163742	24.5479	24
5	Bromomethane	94	1.482	1.474	(0.302)	67537	28.9933	29
6	Chloroethane	64	1.551	1.543	(0.316)	101442	28.6944	29
7	Trichlorofluoromethane	101	1.630	1.631	(0.332)	143766	22.2599	22
9	Ethyl Ether	45	1.797	1.799	(0.366)	80196	19.7360	20
11	Freon 141	81	1.866	1.858	(0.380)	155566	19.5231	20
13	Trichlorotrifluoroethane	101	1.944	1.946	(0.396)	92322	20.7010	21
14	1,1-Dichloroethene	96	1.935	1.936	(0.394)	88087	22.9413	23
15	Carbon Disulfide	76	1.974	1.976	(0.402)	349494	18.6471	19
16	Iodomethane	142	2.043	2.035	(0.416)	92726	18.6185	19
19	3-Chloro-1-Propene	41	2.230	2.232	(0.455)	194504	19.6260	20
20	Methylene Chloride	84	2.299	2.300	(0.469)	101764	19.2795	19
21	Acetone	43	2.328	2.330	(0.475)	96938	33.9282	34
22	trans-1,2-Dichloroethene	96	2.426	2.428	(0.495)	100478	20.8362	21
23	Methyl Acetate	43	2.407	2.409	(0.491)	338067	10.0406	10
24	Methyl tert-Butyl Ether	73	2.495	2.497	(0.509)	361553	19.3297	19
25	tert-Butyl alcohol	59	2.535	2.537	(0.517)	92221	88.3791	88
30	Acrylonitrile	53	2.928	2.920	(0.597)	116373	35.2098	35
31	1,1-Dichloroethane	63	2.909	2.901	(0.593)	254983	20.7184	21
33	cis-1,2-Dichloroethene	96	3.420	3.422	(0.697)	106169	20.6037	21

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
34	2,2-Dichloropropane	77	3.528	3.530	(0.719)	181817	21.4179	21
35	Bromochloromethane	128	3.627	3.629	(0.739)	76589	20.7601	21
37	Cyclohexane	84	3.666	3.658	(0.747)	139095	21.1369	21
38	Chloroform	83	3.706	3.707	(0.755)	191133	20.3033	20
39	Ethyl Acetate	43	3.863	3.855	(0.787)	3851	7.64333	8 (M)
40	Methyl Acrylate	55	3.863	3.865	(0.787)	141747	19.6488	20
$ 41	Dibromofluoromethane	111	3.922	3.924	(0.799)	106678	18.3542	18
42	Tetrahydrofuran	42	3.932	3.924	(0.801)	98608	40.3957	40
43	Carbon Tetrachloride	117	3.902	3.904	(0.795)	166337	21.4603	21
44	1,1,1-Trichloroethane	97	3.971	3.973	(0.809)	143179	20.6927	21
45	2-Butanone	43	4.079	4.071	(0.832)	123044	29.6676	30
46	1,1-Dichloropropene	75	4.129	4.131	(0.842)	170824	19.8465	20
49	1-Chlorobutane	56	4.178	4.180	(0.852)	252507	20.4420	20
51	Propionitrile	54	4.404	4.406	(0.898)	241815	211.546	210
52	Benzene	78	4.424	4.426	(0.902)	438278	20.2650	20
53	2-Methyl-2-Propenenitrile	41	4.443	4.436	(0.906)	108589	13.4840	13
$ 55	1,2-Dichloroethane-d4	65	4.571	4.563	(0.932)	128583	18.1457	18
56	1,2-Dichloroethane	62	4.650	4.652	(0.948)	180059	21.0512	21
59	Methyl Cyclohexane	83	5.103	5.095	(1.040)	131063	20.7885	21
60	Trichloroethene	130	5.103	5.105	(1.040)	131905	20.7323	21
63	Dibromomethane	93	5.536	5.538	(1.128)	71114	20.5149	20
64	1,2-Dichloropropane	63	5.634	5.636	(1.148)	154445	21.6421	22
65	Bromodichloromethane	83	5.723	5.724	(1.166)	130026	19.2862	19
66	Methyl Methacrylate	69	5.900	5.902	(1.203)	235011	75.1230	75
70	cis-1,3-Dichloropropene	75	6.352	6.354	(1.295)	199392	19.5356	20
71	Chloroacetonitrile	48	6.706	6.698	(1.367)	16957	58.0785	58
72	2-Nitropropane	41	6.775	6.777	(1.381)	72371	36.1151	36
73	trans-1,3-Dichloropropene	75	6.982	6.984	(1.423)	177398	19.1484	19
74	1,1,2-Trichloroethane	97	7.120	7.122	(1.451)	94216	20.8956	21
* 75	Chlorobenzene-d5	117	7.966	7.968	(1.000)	442311	25.0000	
76	Toluene	91	6.588	6.580	(0.827)	376498	19.9904	20
$ 77	Toluene-d8	98	6.539	6.541	(0.821)	316232	20.1069	20
78	1,1-Dichloro-2-propanone	43	6.805	6.797	(0.854)	427002	92.4514	92
79	4-Methyl-2-Pentanone	43	6.943	6.944	(0.872)	177623	21.8145	22
80	Tetrachloroethene	164	6.962	6.964	(0.874)	78125	21.0224	21
81	Ethyl Methacrylate	69	7.149	7.151	(0.897)	205652	20.5132	20
82	Dibromochloromethane	129	7.297	7.289	(0.916)	133641	18.8846	19
83	1,3-Dichloropropane	76	7.366	7.368	(0.925)	205347	20.1025	20
84	1,2-Dibromoethane	107	7.494	7.495	(0.941)	119837	20.1233	20
86	2-Hexanone	43	7.710	7.712	(0.968)	151709	24.9833	25
87	1-Chlorohexane	91	7.976	7.968	(1.001)	123780	19.4047	19
88	Chlorobenzene	112	7.976	7.978	(1.001)	308471	20.1613	20
89	1,1,1,2-Tetrachloroethane	131	8.045	8.037	(1.010)	110934	20.0477	20
90	Ethylbenzene	106	8.015	8.007	(1.006)	136532	20.0466	20
91	Xylene (total)mp	106	8.143	8.145	(1.022)	345425	40.5769	40
92	Xylene (total)o	106	8.517	8.519	(1.069)	163421	20.0401	20
93	Styrene	104	8.566	8.568	(1.075)	255695	18.7176	19
94	Bromoform	173	8.586	8.588	(1.078)	70691	18.1053	18
* 95	1,4-Dichlorobenzene-d4	152	10.012	10.014	(1.000)	156570	25.0000	
96	Isopropylbenzene	105	8.802	8.804	(0.879)	387699	20.0795	20
97	Bromobenzene	156	9.127	9.129	(0.912)	97685	20.0365	20
98	1,1,2,2-Tetrachloroethane	83	9.225	9.227	(0.921)	138263	20.7519	21
100	1,2,3-Trichloropropane	110	9.333	9.335	(0.932)	44789	21.2177	21
101	trans-1,4-Dichloro-2-Butene	53	9.373	9.375	(0.936)	55493	25.5981	26

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	CONCENTRATIONS FINAL (ug/L)
102	n-Propylbenzene	91	9.166	9.158	(0.915)	430176	20.7899	21
103	2-Chlorotoluene	91	9.294	9.296	(0.928)	274812	19.2700	19
104	4-Chlorotoluene	91	9.432	9.434	(0.942)	279345	20.7783	21
105	1,3,5-Trimethylbenzene	105	9.333	9.335	(0.932)	319228	20.6416	21
106	tert-Butylbenzene	119	9.609	9.611	(0.960)	310462	20.7557	21
107	1,2,4-Trimethylbenzene	105	9.668	9.670	(0.966)	325118	20.9286	21
108	sec-Butylbenzene	105	9.766	9.768	(0.975)	363909	22.2612	22
109	4-Isopropyltoluene	119	9.894	9.896	(0.988)	359455	20.5389	20
110	1,3-Dichlorobenzene	146	9.953	9.955	(0.994)	174112	20.4174	20
111	1,4-Dichlorobenzene	146	10.032	10.034	(1.002)	176313	20.2576	20
112	1,2-Dichlorobenzene	146	10.396	10.388	(1.038)	168990	20.1365	20
113	Benzyl Chloride	126	10.239	10.241	(1.023)	50398	19.4226	19
115	n-Butylbenzene	91	10.258	10.260	(1.025)	410157	19.4500	19
118	1,2,4,5-Tetramethylbenzene	119	10.918	10.919	(2.225)	19310	1.29336	1
119	1,2-Dibromo-3-chloropropane	75	11.085	11.087	(1.107)	23084	17.3530	17
120	Nitrobenzene	77	11.577	11.579	(1.156)	28160	95.0801	95
121	1,2,4-Trichlorobenzene	180	11.695	11.697	(1.168)	87464	19.1835	19
122	Hexachlorobutadiene	225	11.675	11.677	(1.166)	35191	22.6521	23
123	Naphthalene	128	11.970	11.972	(1.196)	279788	20.2442	20
124	1,2,3-Trichlorobenzene	180	12.138	12.139	(1.212)	82349	19.7951	20
$ 125	Bromofluorobenzene	95	9.038	9.040	(0.903)	137507	23.4101	23
M 126	1,2-Dichloroethene (total)	100				206647	41.4399	41
M 127	Xylene (total)	100				508846	60.6170	61

QC Flag Legend

M - Compound response manually integrated.

Data File: L9342.D

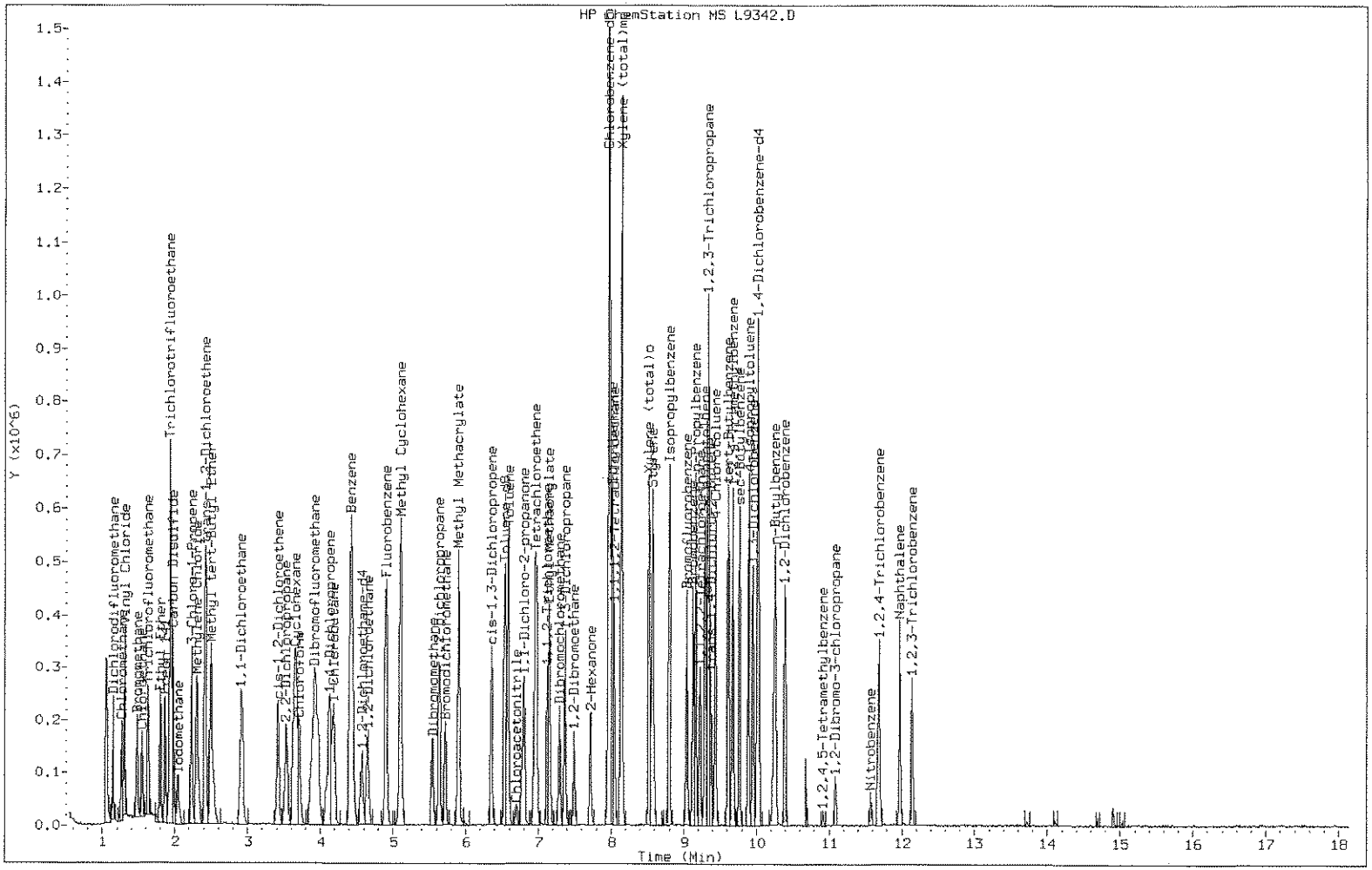
Date: 01-AUG-2007 10:34

Client ID: LCS

Instrument: msl.i

Sample Info: LCS

Operator: D. HUMBERT



08/09/2007

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: Lab Sample ID: LCS 220-8356/2

Matrix: Water Lab File ID: L9372.D

Analysis Method: 8260B Date Received:

Sample wt/vol: 5 (mL) Date Analyzed: 08/02/2007 09:59

Level: (low/med) Low Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) Soil Aliquot:

Soil Extract Vol.: % Moisture:

Analy. Batch No.: 8356 Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	33.8		10	1.4
71-43-2	Benzene	20.2		5.0	0.40
75-27-4	Bromodichloromethane	19.1		5.0	0.40
75-25-2	Bromoform	17.6		5.0	0.80
74-83-9	Bromomethane	24.2		5.0	1.2
78-93-3	2-Butanone (MEK)	27.9		10	1.2
75-15-0	Carbon disulfide	19.0		5.0	0.90
56-23-5	Carbon tetrachloride	21.6		5.0	1.0
108-90-7	Chlorobenzene	19.7		5.0	0.40
75-00-3	Chloroethane	26.9		5.0	0.80
67-66-3	Chloroform	20.1		5.0	0.70
74-87-3	Chloromethane	24.3		5.0	0.50
124-48-1	Dibromochloromethane	18.6		5.0	0.50
75-34-3	1,1-Dichloroethane	20.8		5.0	0.60
107-06-2	1,2-Dichloroethane	20.9		5.0	0.60
75-35-4	1,1-Dichloroethene	22.7		5.0	0.70
78-87-5	1,2-Dichloropropane	20.6		5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	19.3		5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	19.8		5.0	0.80
100-41-4	Ethylbenzene	20.4		5.0	1.0
591-78-6	2-Hexanone	24.0		10	0.80
75-09-2	Methylene Chloride	20.2		5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	20.3		10	0.70
100-42-5	Styrene	18.2		5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	21.1		5.0	0.40
127-18-4	Tetrachloroethene	20.2		5.0	0.50
108-88-3	Toluene	19.4		5.0	0.30
71-55-6	1,1,1-Trichloroethane	20.5		5.0	0.40
79-00-5	1,1,2-Trichloroethane	20.1		5.0	0.60
79-01-6	Trichloroethene	20.0		5.0	0.70
75-01-4	Vinyl chloride	23.6		5.0	0.80
1330-20-7	Xylenes, Total	60.7		5.0	1.0
156-59-2	cis-1,2-Dichloroethene	19.9		5.0	0.60
156-60-5	trans-1,2-Dichloroethene	20.2		5.0	0.50

STL-INC

Volatile Report SW-846 Method 8260B

Data file : \\target1_ct\files\chem\VOA\msl.i\L079370.b\L9372.D
Lab Smp Id: LCS Client Smp ID: LCS
Inj Date : 02-AUG-2007 09:59 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : LCS
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 56 QC Sample: LCS
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONMSLNT

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1 Fluorobenzene	96	4.896	4.901	(1.000)	441851	25.0000	
2 Dichlorodifluoromethane	85	1.147	1.143	(0.234)	107678	31.7383	32
3 Chloromethane	50	1.265	1.261	(0.258)	144131	24.2526	24
4 Vinyl Chloride	62	1.304	1.300	(0.267)	154337	23.5921	24
5 Bromomethane	94	1.482	1.477	(0.303)	55251	24.1845	24
6 Chloroethane	64	1.550	1.546	(0.317)	93220	26.8862	27
7 Trichlorofluoromethane	101	1.629	1.625	(0.333)	139081	21.9571	22
9 Ethyl Ether	45	1.796	1.792	(0.367)	78803	19.7738	20
11 Freon 141	81	1.855	1.861	(0.379)	156410	20.0143	20
13 Trichlorotrifluoroethane	101	1.944	1.950	(0.397)	91889	21.0082	21
14 1,1-Dichloroethene	96	1.934	1.930	(0.395)	85620	22.7365	23
15 Carbon Disulfide	76	1.973	1.969	(0.403)	349386	19.0072	19
16 Iodomethane	142	2.032	2.038	(0.415)	69225	14.1725	14
19 3-Chloro-1-Propene	41	2.229	2.225	(0.455)	199039	20.4778	20
20 Methylene Chloride	84	2.298	2.304	(0.469)	104829	20.2500	20
21 Acetone	43	2.328	2.323	(0.476)	94720	33.8026	34
22 trans-1,2-Dichloroethene	96	2.416	2.422	(0.494)	95500	20.1926	20
23 Methyl Acetate	43	2.406	2.402	(0.492)	328148	9.93726	10
24 Methyl tert-Butyl Ether	73	2.495	2.491	(0.510)	364801	19.8861	20
25 tert-Butyl alcohol	59	2.534	2.530	(0.518)	96330	94.1287	94
30 Acrylonitrile	53	2.928	2.924	(0.598)	115109	35.5109	36
31 1,1-Dichloroethane	63	2.898	2.904	(0.592)	250775	20.7764	21

			QUANT SIG					CONCENTRATIONS ON-COLUMN	FINAL
		Compounds	MASS	RT	EXP RT	REL RT	RESPONSE	(ug/L)	(ug/L)
	33	cis-1,2-Dichloroethene	96	3.410	3.416	(0.697)	100361	19.8588	20
	34	2,2-Dichloropropane	77	3.528	3.524	(0.721)	178900	21.4879	21
	35	Bromochloromethane	128	3.626	3.622	(0.741)	73810	20.3995	20
	37	Cyclohexane	84	3.656	3.652	(0.747)	132254	20.4918	20
	38	Chloroform	83	3.705	3.701	(0.757)	185938	20.1392	20
	39	Ethyl Acetate	43	3.853	3.858	(0.787)	4487	9.08043	9(M)
	40	Methyl Acrylate	55	3.863	3.858	(0.789)	135418	19.1399	19
$	41	Dibromofluoromethane	111	3.922	3.927	(0.801)	107778	18.9074	19
	42	Tetrahydrofuran	42	3.922	3.917	(0.801)	92470	38.6247	39
	43	Carbon Tetrachloride	117	3.902	3.898	(0.797)	164267	21.6092	22
	44	1,1,1-Trichloroethane	97	3.961	3.967	(0.809)	139056	20.4913	20
	45	2-Butanone	43	4.069	4.065	(0.831)	113408	27.8809	28
	46	1,1-Dichloropropene	75	4.118	4.124	(0.841)	169710	20.1040	20
	49	1-Chlorobutane	56	4.177	4.173	(0.853)	244532	20.1849	20
	51	Propionitrile	54	4.404	4.399	(0.900)	221216	197.324	200
	52	Benzene	78	4.414	4.419	(0.902)	428384	20.1963	20
	53	2-Methyl-2-Propenenitrile	41	4.443	4.439	(0.908)	110384	13.9759	14
$	55	1,2-Dichloroethane-d4	65	4.561	4.567	(0.932)	129362	18.6139	19
	56	1,2-Dichloroethane	62	4.640	4.645	(0.948)	175541	20.9258	21
	59	Methyl Cyclohexane	83	5.092	5.098	(1.040)	128088	20.7153	21
	60	Trichloroethene	130	5.102	5.098	(1.042)	125094	20.0477	20
	63	Dibromomethane	93	5.535	5.531	(1.131)	67574	19.8762	20
	64	1,2-Dichloropropane	63	5.634	5.629	(1.151)	144138	20.5942	20
	65	Bromodichloromethane	83	5.722	5.718	(1.169)	126019	19.0587	19
	66	Methyl Methacrylate	69	5.899	5.895	(1.205)	229381	74.7624	75
	70	cis-1,3-Dichloropropene	75	6.352	6.348	(1.297)	193582	19.3386	19
	71	Chloroacetonitrile	48	6.696	6.692	(1.368)	20246	70.7045	71
	72	2-Nitropropane	41	6.775	6.771	(1.384)	71135	36.1950	36
	73	trans-1,3-Dichloropropene	75	6.972	6.977	(1.424)	180040	19.8150	20
	74	1,1,2-Trichloroethane	97	7.119	7.125	(1.454)	89053	20.1382	20
*	75	Chlorobenzene-d5	117	7.956	7.961	(1.000)	440917	25.0000	
	76	Toluene	91	6.578	6.584	(0.827)	364045	19.3903	19
$	77	Toluene-d8	98	6.529	6.535	(0.821)	314337	20.0496	20
	78	1,1-Dichloro-2-propanone	43	6.795	6.800	(0.854)	422133	91.6862	92
	79	4-Methyl-2-Pentanone	43	6.942	6.938	(0.873)	164384	20.2524	20
	80	Tetrachloroethene	164	6.962	6.958	(0.875)	74666	20.1551	20
	81	Ethyl Methacrylate	69	7.149	7.145	(0.899)	203152	20.3279	20
	82	Dibromochloromethane	129	7.287	7.292	(0.916)	131274	18.6088	19
	83	1,3-Dichloropropane	76	7.365	7.361	(0.926)	200409	19.6812	20
	84	1,2-Dibromoethane	107	7.493	7.489	(0.942)	119024	20.0499	20
	86	2-Hexanone	43	7.710	7.705	(0.969)	145229	23.9918	24
	87	1-Chlorohexane	91	7.965	7.971	(1.001)	131912	20.7449	21
	88	Chlorobenzene	112	7.975	7.971	(1.002)	300534	19.7047	20
	89	1,1,1,2-Tetrachloroethane	131	8.034	8.040	(1.010)	105829	19.1856	19
	90	Ethylbenzene	106	8.005	8.010	(1.006)	138220	20.3586	20
	91	Xylene (total)mp	106	8.143	8.138	(1.023)	341978	40.2990	40
	92	Xylene (total)o	106	8.516	8.512	(1.070)	165992	20.4197	20
	93	Styrene	104	8.566	8.561	(1.077)	247233	18.1554	18
	94	Bromoform	173	8.585	8.581	(1.079)	68382	17.5693	18
*	95	1,4-Dichlorobenzene-d4	152	10.012	10.018	(1.000)	154291	25.0000	
	96	Isopropylbenzene	105	8.802	8.797	(0.879)	389915	20.4926	20
	97	Bromobenzene	156	9.126	9.122	(0.912)	99568	20.7244	21
	98	1,1,2,2-Tetrachloroethane	83	9.225	9.221	(0.921)	138665	21.1196	21
	100	1,2,3-Trichloropropane	110	9.333	9.329	(0.932)	41681	20.0370	20

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
101	trans-1,4-Dichloro-2-Butene	53	9.372	9.368	(0.936)	67253	31.4811	31
102	n-Propylbenzene	91	9.156	9.162	(0.915)	421763	20.6844	21
103	2-Chlorotoluene	91	9.294	9.289	(0.928)	276433	19.6700	20
104	4-Chlorotoluene	91	9.431	9.437	(0.942)	281614	21.2565	21
105	1,3,5-Trimethylbenzene	105	9.333	9.339	(0.932)	319427	20.9596	21
106	tert-Butylbenzene	119	9.609	9.604	(0.960)	309412	20.9910	21
107	1,2,4-Trimethylbenzene	105	9.668	9.673	(0.966)	328031	21.4280	21
108	sec-Butylbenzene	105	9.766	9.762	(0.975)	358038	22.2255	22
109	4-Isopropyltoluene	119	9.894	9.890	(0.988)	363257	21.0627	21
110	1,3-Dichlorobenzene	146	9.953	9.949	(0.994)	172165	20.4873	20
111	1,4-Dichlorobenzene	146	10.032	10.027	(1.002)	178883	20.8565	21
112	1,2-Dichlorobenzene	146	10.386	10.391	(1.037)	169335	20.4756	20
113	Benzyl Chloride	126	10.238	10.234	(1.023)	54461	21.2985	21
115	n-Butylbenzene	91	10.258	10.254	(1.025)	416229	20.0294	20
118	1,2,4,5-Tetramethylbenzene	119	10.917	10.913	(2.230)	15742	1.07507	1
119	1,2-Dibromo-3-chloropropane	75	11.084	11.080	(1.107)	23462	17.8977	18
120	Nitrobenzene	77	11.566	11.572	(1.155)	28922	99.0953	99
121	1,2,4-Trichlorobenzene	180	11.694	11.690	(1.168)	83685	18.6258	19
122	Hexachlorobutadiene	225	11.675	11.670	(1.166)	34617	22.6118	23
123	Naphthalene	128	11.970	11.966	(1.196)	232573	17.0765	17
124	1,2,3-Trichlorobenzene	180	12.137	12.133	(1.212)	77807	18.9796	19
$ 125	Bromofluorobenzene	95	9.038	9.043	(0.903)	140346	24.2464	24
M 126	1,2-Dichloroethene (total)	100				195861	40.0514	40
M 127	Xylene (total)	100				507970	60.7187	61

QC Flag Legend

M - Compound response manually integrated.

Data File: L9372.D

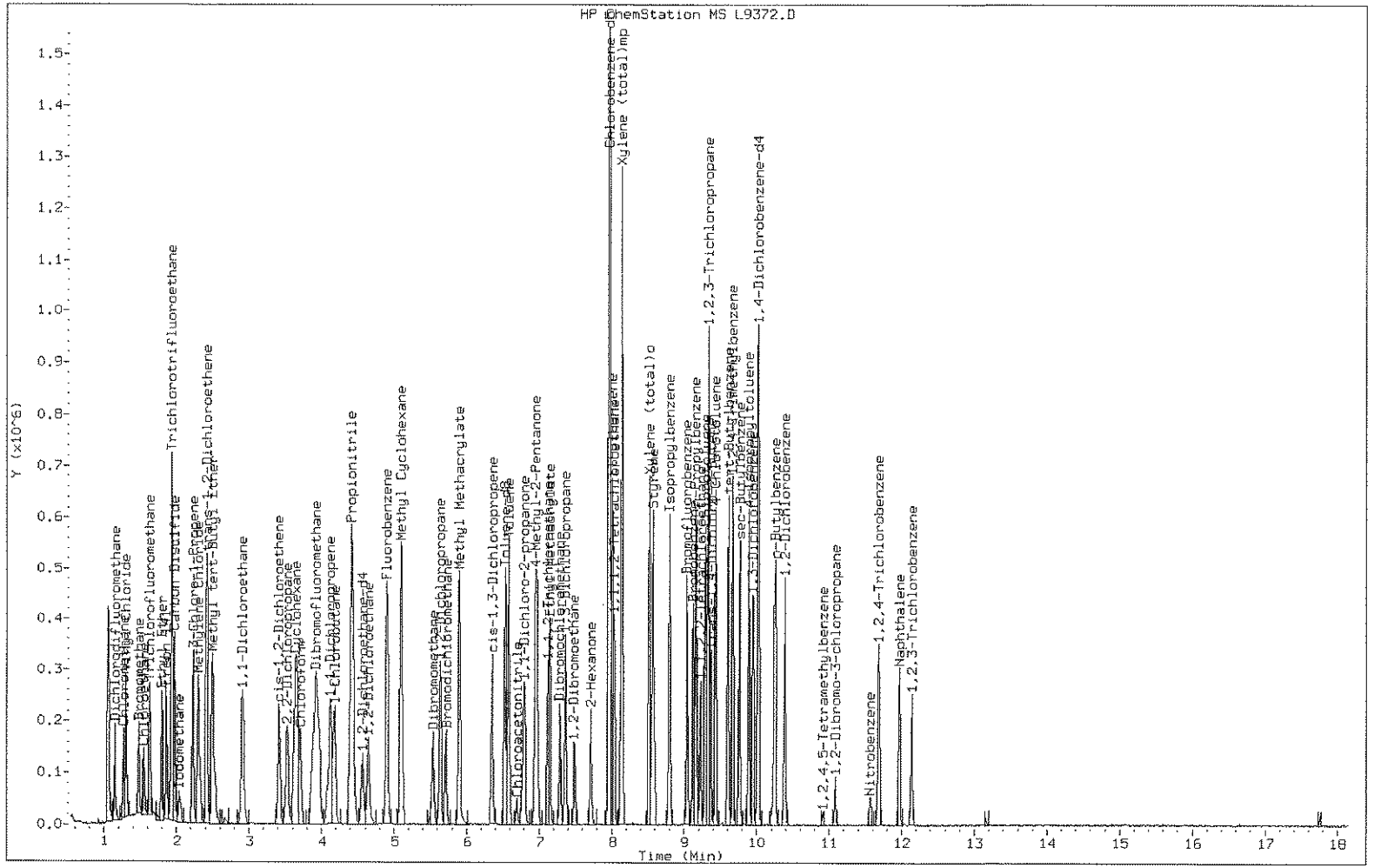
Date: 02-AUG-2007 09:59

Client ID: LCS

Instrument: msl.i

Sample Info: LCS

Operator: D. HUMBERT



08/09/2007

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Connecticut | Job No.: 220-2277-1

SDG No.: 220-2277

Client Sample ID: | Lab Sample ID: MSB 220-8356/24

Matrix: Water | Lab File ID: L9395.D

Analysis Method: 8260B | Date Received:

Sample wt/vol: 5 (mL) | Date Analyzed: 08/02/2007 20:32

Level: (low/med) Low | Dilution Factor: 1

GC Column/ID: RTX-VMS 0.18 (mm) | Soil Aliquot:

Soil Extract Vol.: | % Moisture:

Analy. Batch No.: 8356 | Units: ug/L

CAS No.	Compound Name	Result	Q	RL	MDL
67-64-1	Acetone	48.9		10	1.4
71-43-2	Benzene	53.2		5.0	0.40
75-27-4	Bromodichloromethane	52.0		5.0	0.40
75-25-2	Bromoform	47.0		5.0	0.80
74-83-9	Bromomethane	53.7		5.0	1.2
78-93-3	2-Butanone (MEK)	54.0		10	1.2
75-15-0	Carbon disulfide	50.3		5.0	0.90
56-23-5	Carbon tetrachloride	57.7		5.0	1.0
108-90-7	Chlorobenzene	53.5		5.0	0.40
75-00-3	Chloroethane	64.7		5.0	0.80
67-66-3	Chloroform	54.8		5.0	0.70
74-87-3	Chloromethane	50.7		5.0	0.50
124-48-1	Dibromochloromethane	48.1		5.0	0.50
75-34-3	1,1-Dichloroethane	55.0		5.0	0.60
107-06-2	1,2-Dichloroethane	54.2		5.0	0.60
75-35-4	1,1-Dichloroethene	56.8		5.0	0.70
78-87-5	1,2-Dichloropropane	55.5		5.0	0.90
10061-01-5	cis-1,3-Dichloropropene	49.5		5.0	0.50
10061-02-6	trans-1,3-Dichloropropene	49.7		5.0	0.80
100-41-4	Ethylbenzene	52.8		5.0	1.0
591-78-6	2-Hexanone	55.3		10	0.80
75-09-2	Methylene Chloride	53.9		5.0	0.40
108-10-1	4-Methyl-2-pentanone (MIBK)	56.3		10	0.70
100-42-5	Styrene	54.2		5.0	0.50
79-34-5	1,1,2,2-Tetrachloroethane	52.2		5.0	0.40
127-18-4	Tetrachloroethene	54.6		5.0	0.50
108-88-3	Toluene	52.8		5.0	0.30
71-55-6	1,1,1-Trichloroethane	53.9		5.0	0.40
79-00-5	1,1,2-Trichloroethane	55.0		5.0	0.60
79-01-6	Trichloroethene	53.4		5.0	0.70
75-01-4	Vinyl chloride	53.7		5.0	0.80
1330-20-7	Xylenes, Total	162		5.0	1.0
156-59-2	cis-1,2-Dichloroethene	55.1		5.0	0.60
156-60-5	trans-1,2-Dichloroethene	56.0		5.0	0.50

STL-CT

Volatile Report SW-846 Method 8260B

Data file : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\L9395.D
Lab Smp Id: MSB Client Smp ID: MSB
Inj Date : 02-AUG-2007 20:32 MS Autotune Date: 26-APR-2004 14:21
Operator : D. HUMBERT Inst ID: msl.i
Smp Info : MSB
Misc Info : : ;;; ; 8260 ; 1 ; LLW
Comment :
Method : \\TARGET1_CT\FILES\chem\VOA\msl.i\L079370.b\L8260BNW.m
Meth Date : 02-Aug-2007 12:01 dave Quant Type: ISTD
Cal Date : 26-JUL-2007 15:11 Cal File: L9156.D
Als bottle: 77 QC Sample: BS
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14
Processing Host: CONMSLNT

Concentration Formula: Amt * DF * Uf * 1/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	5.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
* 1	Fluorobenzene	96	4.887	4.901	(1.000)	444303	25.0000	
2	Dichlorodifluoromethane	85	1.149	1.143	(0.235)	165959	48.6468	49
3	Chloromethane	50	1.267	1.261	(0.259)	303271	50.7490	51
4	Vinyl Chloride	62	1.306	1.300	(0.267)	353066	53.6721	54
5	Bromomethane	94	1.473	1.477	(0.301)	123336	53.6887	54
6	Chloroethane	64	1.532	1.546	(0.314)	225484	64.6745	65
7	Trichlorofluoromethane	101	1.621	1.625	(0.332)	345500	54.2441	54
8	Dichlorofluoromethane	67	1.641	1.645	(0.336)	648195	55.7663	56
9	Ethyl Ether	45	1.798	1.792	(0.368)	218391	54.4977	54
10	Ethanol	45	1.857	1.861	(0.380)	184756	585.039	580
11	Freon 141	81	1.857	1.861	(0.380)	423514	53.8940	54
12	Freon 123a	67	1.641	1.645	(0.336)	648195	55.7663	56
13	Trichlorotrifluoroethane	101	1.946	1.950	(0.398)	244765	55.6509	56
14	1,1-Dichloroethene	96	1.926	1.930	(0.394)	215251	56.8446	57
15	Carbon Disulfide	76	1.965	1.969	(0.402)	929917	50.3098	50
16	Iodomethane	142	2.034	2.038	(0.416)	284354	57.8950	58
17	Acrolein	56	2.123	2.127	(0.434)	417380	404.683	400
18	2-Propanol	45	2.211	2.215	(0.452)	63428	62.2106	62
19	3-Chloro-1-Propene	41	2.221	2.225	(0.454)	504324	51.6002	52
20	Methylene Chloride	84	2.300	2.304	(0.471)	280385	53.8635	54
21	Acetone	43	2.319	2.323	(0.475)	137886	48.9356	49
22	trans-1,2-Dichloroethene	96	2.418	2.422	(0.495)	266160	55.9665	56

			QUANT SIG					CONCENTRATIONS ON-COLUMN	FINAL
		Compounds	MASS	RT	EXP RT	REL RT	RESPONSE	(ug/L)	(ug/L)
	23	Methyl Acetate	43	2.398	2.402	(0.491)	1833474	55.2164	55
	24	Methyl tert-Butyl Ether	73	2.487	2.491	(0.509)	995912	53.9897	54
	25	tert-Butyl alcohol	59	2.526	2.530	(0.517)	304976	296.362	300
	26	Acetonitrile	41	2.654	2.658	(0.543)	688739	585.956	580
	27	Isopropyl ether	45	2.782	2.786	(0.569)	1183381	55.1498	55
	28	tert-Butyl ethyl ether	59	3.107	3.111	(0.636)	1172591	54.3860	54
	29	2-Chloro-1,3-Butadiene	88	2.880	2.884	(0.589)	187265	54.4554	54
	30	Acrylonitrile	53	2.920	2.924	(0.597)	375918	115.330	120
	31	1,1-Dichloroethane	63	2.900	2.904	(0.593)	667755	55.0175	55
	32	Vinyl Acetate	43	3.107	3.111	(0.636)	814495	50.2874	50
	33	cis-1,2-Dichloroethene	96	3.402	3.416	(0.696)	280081	55.1148	55
	34	2,2-Dichloropropane	77	3.520	3.524	(0.720)	394439	47.1151	47
	35	Bromochloromethane	128	3.618	3.622	(0.740)	194247	53.3895	53
	36	1-Bromopropane	43	3.608	3.612	(0.738)	511863	55.9238	56
	37	Cyclohexane	84	3.648	3.652	(0.746)	352038	54.2447	54
	38	Chloroform	83	3.697	3.701	(0.756)	508964	54.8222	55
	39	Ethyl Acetate	43	3.844	3.858	(0.787)	60303	121.363	120
	40	Methyl Acrylate	55	3.844	3.858	(0.787)	400061	56.2323	56
$	41	Dibromofluoromethane	111	3.913	3.927	(0.801)	110193	19.2244	19
	42	Tetrahydrofuran	42	3.903	3.917	(0.799)	272450	113.174	110
	43	Carbon Tetrachloride	117	3.884	3.898	(0.795)	440914	57.6819	58
	44	1,1,1-Trichloroethane	97	3.963	3.967	(0.811)	367754	53.8932	54
	45	2-Butanone	43	4.061	4.065	(0.831)	220741	53.9687	54
	46	1,1-Dichloropropene	75	4.110	4.124	(0.841)	460970	54.3056	54
	47	tert-Amyl methyl ether	73	4.563	4.567	(0.934)	986899	52.7652	53
	48	tert-Butyl formate	57	3.107	3.111	(0.636)	339809	54.6460	55
	49	1-Chlorobutane	56	4.169	4.173	(0.853)	667872	54.8253	55
	50	Heptane	43	4.395	4.399	(0.899)	270105	49.1840	49
	51	Propionitrile	54	4.395	4.399	(0.899)	644582	571.792	570
	52	Benzene	78	4.405	4.419	(0.901)	1133718	53.1545	53
	53	2-Methyl-2-Propenenitrile	41	4.425	4.439	(0.905)	438961	55.2709	55
	54	Isobutyl alcohol	42	4.681	4.685	(0.958)	139754	486.588	490
$	55	1,2-Dichloroethane-d4	65	4.553	4.567	(0.932)	130671	18.6985	19
	56	1,2-Dichloroethane	62	4.641	4.645	(0.950)	456939	54.1699	54
	59	Methyl Cyclohexane	83	5.084	5.098	(1.040)	331955	53.3899	53
	60	Trichloroethene	130	5.094	5.098	(1.042)	335106	53.4080	53
	61	Isopropyl Acetate	43	4.671	4.685	(0.956)	242035	101.502	100
	62	N-Butanol	56	5.468	5.482	(1.119)	143753	621.188	620
	63	Dibromomethane	93	5.527	5.531	(1.131)	184389	53.9369	54
	64	1,2-Dichloropropane	63	5.625	5.629	(1.151)	390288	55.4559	55
	65	Bromodichloromethane	83	5.714	5.718	(1.169)	345675	51.9903	52
	66	Methyl Methacrylate	69	5.891	5.895	(1.205)	334874	108.544	110
	67	1,4-Dioxane	58	5.930	5.934	(1.213)	32944	618.378	620
	68	N-Propyl Acetate	43	6.294	6.298	(1.288)	80744	143.669	140
	69	2-Chloroethylvinylether	63	6.294	6.298	(1.288)	117956	68.7912	69
	70	cis-1,3-Dichloropropene	75	6.344	6.348	(1.298)	498006	49.4756	49
	71	Chloroacetonitrile	48	6.688	6.692	(1.368)	155525	540.138	540
	72	2-Nitropropane	41	6.767	6.771	(1.384)	207009	104.749	100
	73	trans-1,3-Dichloropropene	75	6.973	6.977	(1.427)	454103	49.7023	50
	74	1,1,2-Trichloroethane	97	7.111	7.125	(1.455)	244754	55.0425	55
*	75	Chlorobenzene-d5	117	7.957	7.961	(1.000)	439881	25.0000	
	76	Toluene	91	6.580	6.584	(0.827)	989689	52.8384	53
$	77	Toluene-d8	98	6.530	6.535	(0.821)	316267	20.2202	20
	78	1,1-Dichloro-2-propanone	43	6.796	6.800	(0.854)	1211968	263.856	260

		Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (ug/L)	FINAL (ug/L)
	79	4-Methyl-2-Pentanone	43	6.934	6.938	(0.871)	455801	56.2877	56
	80	Tetrachloroethene	164	6.954	6.958	(0.874)	201685	54.5705	54
	81	Ethyl Methacrylate	69	7.141	7.145	(0.897)	532231	53.3819	53
	82	Dibromochloromethane	129	7.288	7.292	(0.916)	338383	48.0806	48
	83	1,3-Dichloropropane	76	7.357	7.361	(0.925)	531219	52.2912	52
	84	1,2-Dibromoethane	107	7.485	7.489	(0.941)	316992	53.5239	54
	85	n-Butyl Acetate	56	7.642	7.646	(0.960)	279010	54.6576	55
	86	2-Hexanone	43	7.701	7.705	(0.968)	334047	55.3145	55
	87	1-Chlorohexane	91	7.967	7.971	(1.001)	316596	49.9062	50
	88	Chlorobenzene	112	7.967	7.971	(1.001)	814642	53.5382	54
	89	1,1,1,2-Tetrachloroethane	131	8.036	8.040	(1.010)	279804	50.8448	51
	90	Ethylbenzene	106	8.006	8.010	(1.006)	357804	52.8254	53
	91	Xylene (total)mp	106	8.134	8.138	(1.022)	910126	107.503	110
	92	Xylene (total)o	106	8.508	8.512	(1.069)	444627	54.8251	55
	93	Styrene	104	8.557	8.561	(1.075)	735768	54.1578	54
	94	Bromoform	173	8.577	8.581	(1.078)	182593	47.0238	47
*	95	1,4-Dichlorobenzene-d4	152	10.013	10.018	(1.000)	165255	25.0000	
	96	Isopropylbenzene	105	8.793	8.797	(0.878)	1062283	52.1258	52
	97	Bromobenzene	156	9.118	9.122	(0.911)	259531	50.4356	50
	98	1,1,2,2-Tetrachloroethane	83	9.217	9.221	(0.920)	367314	52.2328	52
	99	4-Ethyltoluene	105	9.256	9.260	(0.924)	1046477	52.5324	52
	100	1,2,3-Trichloropropane	110	9.325	9.329	(0.931)	118958	53.3919	53
	101	trans-1,4-Dichloro-2-Butene	53	9.364	9.368	(0.935)	119701	52.3145	52
	102	n-Propylbenzene	91	9.157	9.162	(0.915)	1162309	53.2209	53
	103	2-Chlorotoluene	91	9.285	9.289	(0.927)	789469	52.4487	52
	104	4-Chlorotoluene	91	9.433	9.437	(0.942)	746231	52.5893	52
	105	1,3,5-Trimethylbenzene	105	9.335	9.339	(0.932)	875432	53.6314	54
	106	tert-Butylbenzene	119	9.600	9.604	(0.959)	795792	50.4060	50
	107	1,2,4-Trimethylbenzene	105	9.669	9.673	(0.966)	868881	52.9923	53
	108	sec-Butylbenzene	105	9.758	9.762	(0.974)	899807	52.1505	52
	109	4-Isopropyltoluene	119	9.886	9.890	(0.987)	937546	50.7550	51
	110	1,3-Dichlorobenzene	146	9.945	9.949	(0.993)	464715	51.6312	52
	111	1,4-Dichlorobenzene	146	10.023	10.027	(1.001)	475122	51.7206	52
	112	1,2-Dichlorobenzene	146	10.387	10.391	(1.037)	456752	51.5652	52
	113	Benzyl Chloride	126	10.230	10.234	(1.022)	123145	44.9641	45
	114	1,4-Diethylbenzene	119	10.200	10.204	(2.087)	512089	52.9150	53
	115	n-Butylbenzene	91	10.250	10.254	(1.024)	1067065	47.9417	48
	118	1,2,4,5-Tetramethylbenzene	119	10.909	10.913	(2.232)	815849	55.4094	55
	119	1,2-Dibromo-3-chloropropane	75	11.076	11.080	(1.106)	73259	52.1770	52
	120	Nitrobenzene	77	11.568	11.572	(1.155)	126827	405.716	400
	121	1,2,4-Trichlorobenzene	180	11.686	11.690	(1.167)	269854	56.0766	56
	122	Hexachlorobutadiene	225	11.666	11.670	(1.165)	81327	49.5982	50
	123	Naphthalene	128	11.962	11.966	(1.195)	822675	56.3968	56
	124	1,2,3-Trichlorobenzene	180	12.129	12.133	(1.211)	239551	54.5571	54
$	125	Bromofluorobenzene	95	9.030	9.043	(0.902)	146483	23.6276	24
M	126	1,2-Dichloroethene (total)	100				546241	111.081	110
M	127	Xylene (total)	100				1354753	162.328	160

Data File: L9395.D

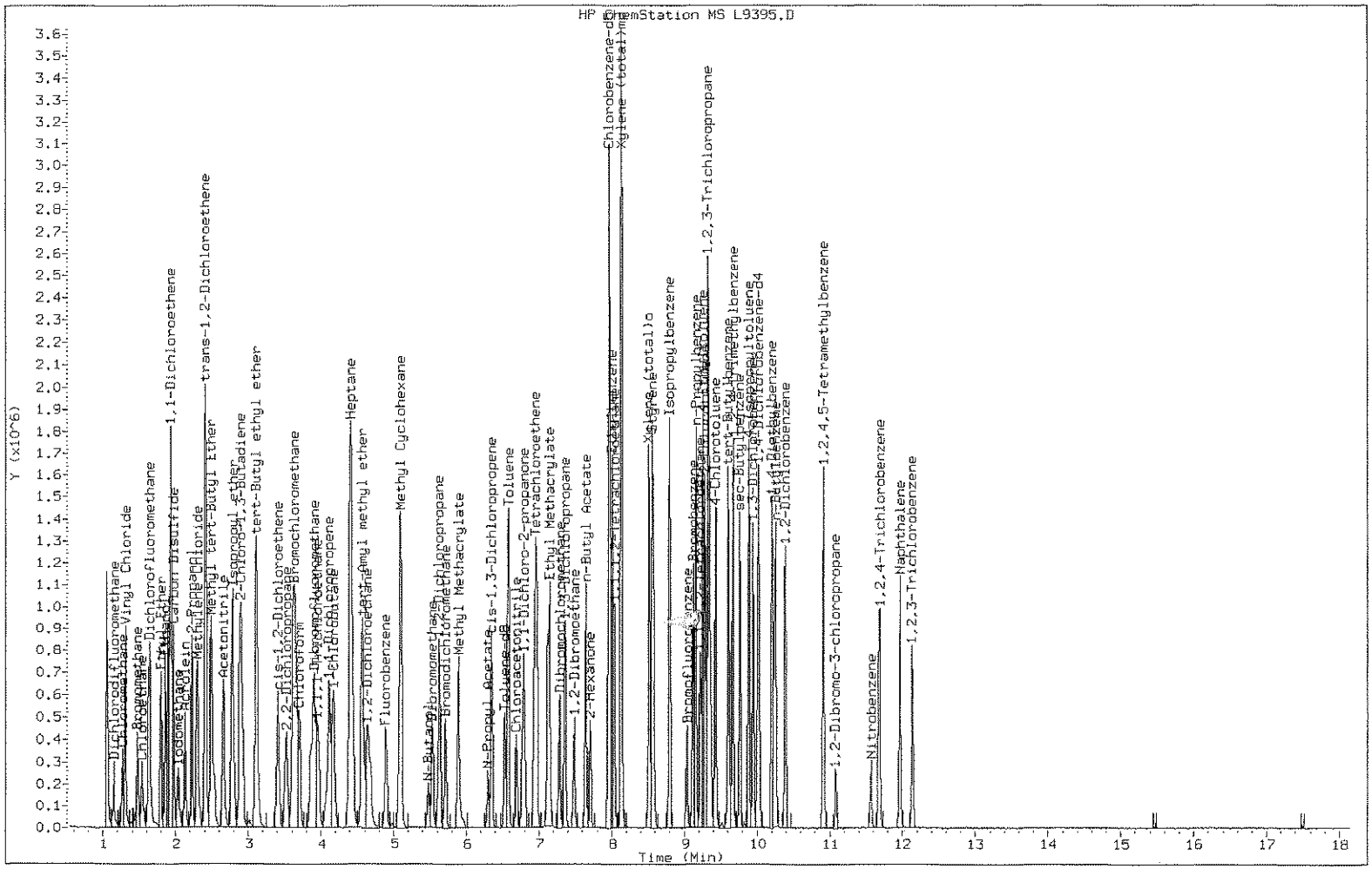
Date: 02-AUG-2007 20:32

Client ID: MSB

Instrument: msl.i

Sample Info: MSB

Operator: D. HUMBERT



08/09/2007

DATA USABILITY SUMMARY REPORT
FOR JULY 2007 GROUNDWATER SAMPLING
LAB REPORT #220-2277
90-30 METROPOLITAN AVENUE SITE
REGO PARK, QUEENS, NEW YORK

This DUSR was prepared using the entire original laboratory report, including the sample data summary report and the extended data package. The sampling event included 24 groundwater samples collected from the groundwater monitoring well network.

Sample Collection Procedures

The samples were collected in laboratory-provided glassware utilizing dedicated sampling equipment. Samples for Quality Assurance/Quality Control (QA/QC) were also obtained to evaluate field sampling methods and laboratory procedures. All sample collection was conducted under Chain of Custody (COC) procedures and in accordance with the QA/QC procedures presented in the approved Remedial Action Work Plan (November 2005).

Sample Analyses

The samples were transmitted via overnight courier and analyzed by TestAmerica Laboratories, Inc. (formerly Severn-Trent Laboratories, Inc.) at their Shelton, Connecticut facility, which is NYSDOH-certified for the analyses performed. The samples were analyzed for Target Compound List (TCL) volatile organic compounds (VOCs) using Method 5035B/8260B. The analytical and prep methods and analytes are appropriate for the intended use of the data. The sample holding times were met and no problems with sample receipt or handling were reported by the laboratory.

A trip blank sample was used to verify that cross-contamination of samples did not occur in the field or laboratory. The trip blank sample was noted to contain acetone, 2-butanone (MEK) and methylene chloride at low estimated concentrations. These VOCs are likely laboratory contaminants and do not appear to significantly affect the sample data.

Equipment blank samples were collected to evaluate potential contamination from field sampling procedures. No VOCs were detected in the equipment blank samples with the exception of low estimated concentrations of methylene chloride and toluene in one equipment blank sample and low estimated detections of acetone, toluene and methylene chloride in the other equipment blank sample. The acetone and methylene chloride detections are B-qualified, indicating that they were noted in an associated laboratory blank sample. These detections appear to be associated with laboratory-related contamination and do not appear to have significantly affected the primary sample results. The toluene detections are also low and estimated and toluene was not detected in any of the primary samples. Therefore, these low-level detections do not indicate significant levels of laboratory or field-related contamination.

A duplicate sample was collected and utilized to evaluate the precision of the laboratory analysis. The results from the duplicate sample (A-9D) and the associated parent sample (A-9) are nearly identical and, therefore, the laboratory results are likely to be reasonably precise.

A matrix spike/matrix spike duplicate (MS/MSD) sample was prepared to evaluate the effect of the matrix on the reliability of the analytical results. Spiking occurs in the laboratory prior to sample preparation and analysis. One MS/MSD sample was collected during this sampling event. Based

on information provided by the analytical laboratory, the MS/MSD results were within QC limits. Therefore, it appears that no matrix-related effects have significantly affected the analytical results.

Method blank (MB) samples were analyzed by the laboratory to evaluate the potential for cross-contamination associated with the sample preparation and analysis. The MB results show low estimated concentrations of acetone and methylene chloride in one sample. The associated primary sample data are B-qualified, indicating that low-level detections of these compounds. any detections of VOCs above the reporting limits and, therefore, cross-contamination associated with sample preparation or analysis does not appear to significantly affect the sample data.

A laboratory control sample (LCS) was used by the laboratory to verify the accuracy and precision of the analyses. The LCS percent recoveries (%Rs) and relative percent differences (RPDs) were all within established guidelines and, therefore, the sample data do not appear to have been affected.

Conclusions

The groundwater samples were collected in accordance with NYSDEC guidance. No field or laboratory conditions occurred that would result in non-valid analytical data. The data appear to be adequate for their intended purpose.

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