Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV111621\

Data File: VV023522.D

Acq On : 16 Nov 2021 10:01

Operator : SY/MD Sample : VSTDCCC005

: 25.0mL/MSVOA\_V/WATER Misc ALS Vial : 2 Sample Multiplier: 1

Quant Time: Nov 17 00:43:27 2021

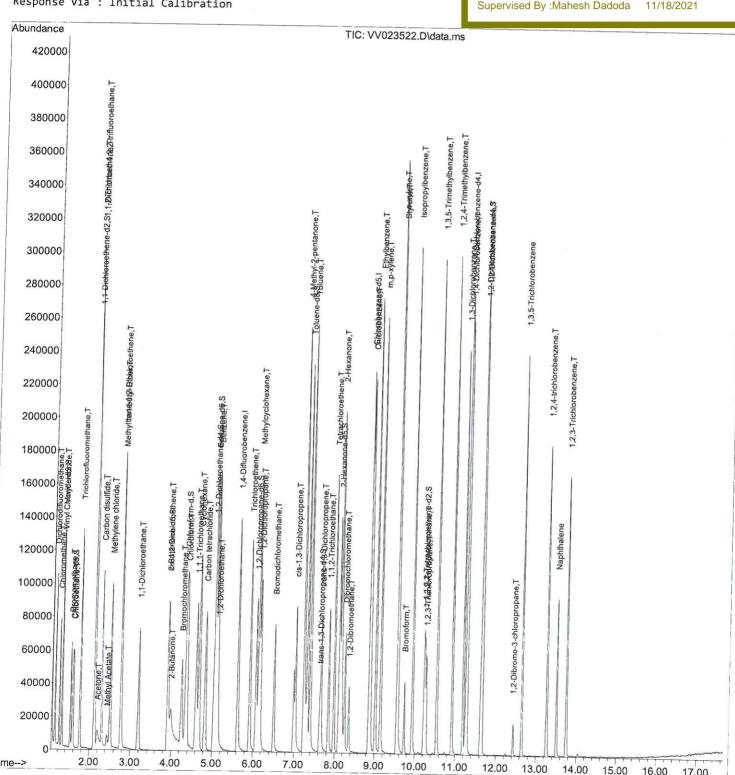
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Tue Nov 16 02:06:43 2021 Response via : Initial Calibration

Instrument: MSVOA\_V **LabSampleld**: STDCCC005

### **Manual IntegrationsAPPROVED**

Reviewed By :John Carlone 11/17/2021 Supervised By :Mahesh Dadoda 11/18/2021



16.00

17.00

15.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV111621\

Data File : VV023522.D

Acq On : 16 Nov 2021 10:01

Operator : SY/MD Sample : VSTDCCC005

Misc : 25.0mL/MSVOA\_V/WATER ALS Vial : 2 Sample Multiplier: 1

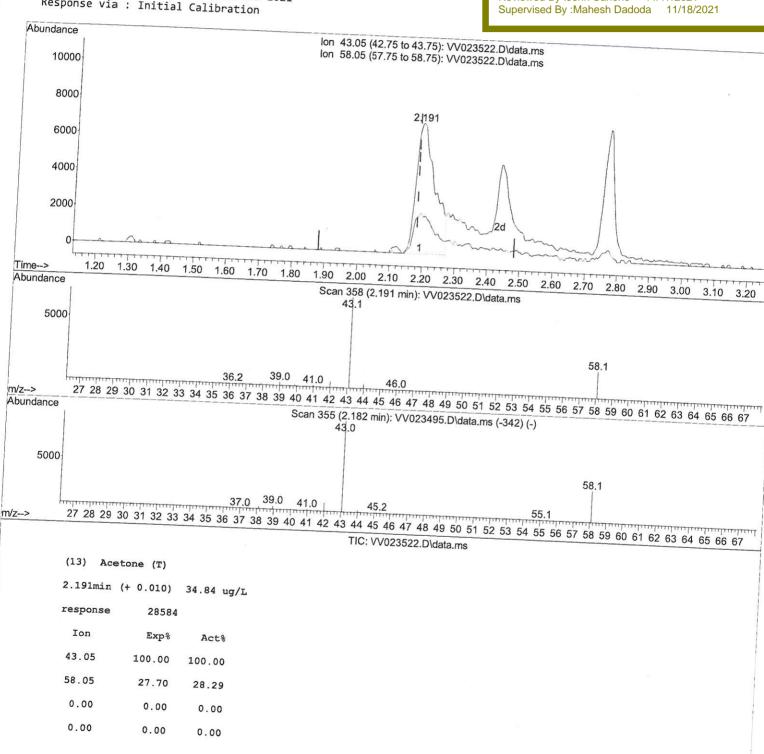
Quant Time: Nov 17 00:43:27 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Tue Nov 16 02:06:43 2021 Response via : Initial Calibration

Instrument: MSVOA\_V LabSampleId : VSTDCCC005

#### **Manual IntegrationsAPPROVED**



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV111621\

Data File : VV023522.D

: 16 Nov 2021 10:01 Acq On

Operator : SY/MD Sample : VSTDCCC005

Misc : 25.0mL/MSVOA\_V/WATER ALS Vial : 2 Sample Multiplier: 1

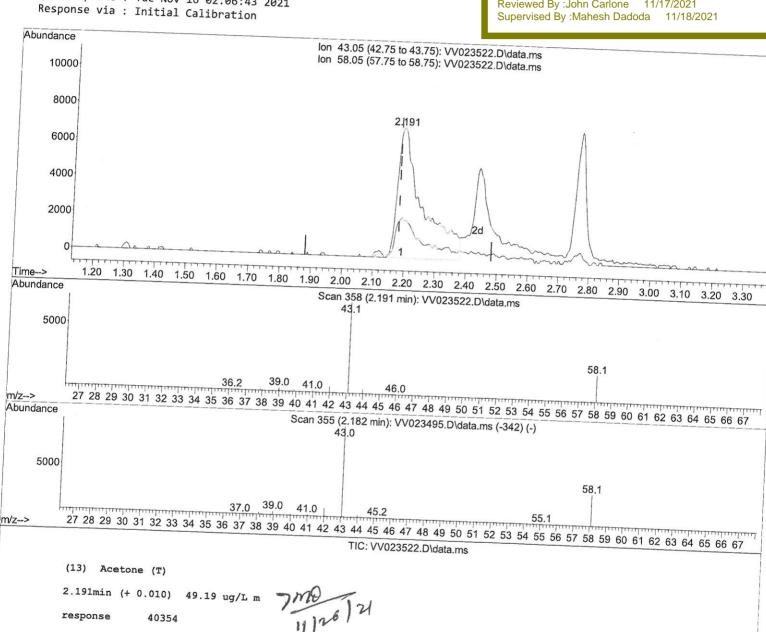
Quant Time: Nov 17 00:43:27 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Tue Nov 16 02:06:43 2021 Instrument: MSVOA\_V LabSampleId : VSTDCCC005

#### **Manual IntegrationsAPPROVED**

Reviewed By :John Carlone 11/17/2021



Ion Ехр% Act% 43.05 100.00 100.00 58.05 27.70 20.04 0.00 0.00 0.00 0.00 0.00 0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV111621\

Data File : VV023522.D

Acq On : 16 Nov 2021 10:01

Operator : SY/MD Sample : VSTDCCC005

Misc : 25.0mL/MSVOA\_V/WATER ALS Vial : 2 Sample Multiplier: 1

Quant Time: Nov 17 00:43:27 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR110421WMA.M

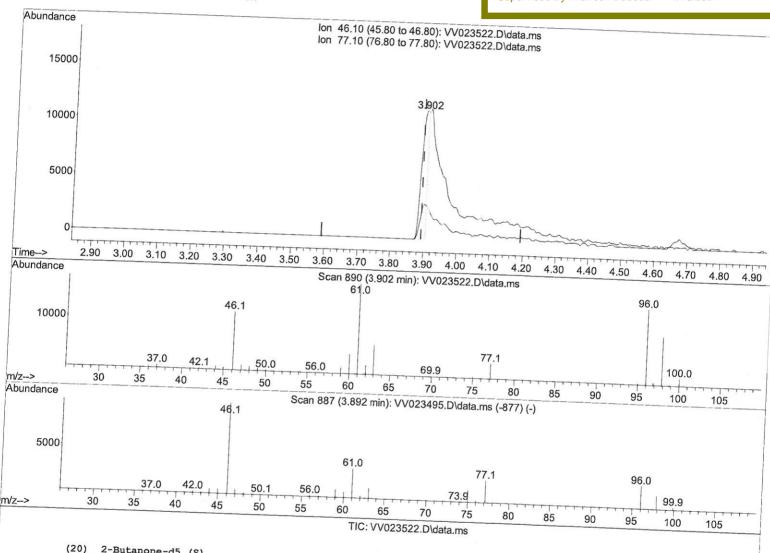
Quant Title : TRACE VOA SFAM1.0

QLast Update : Tue Nov 16 02:06:43 2021 Response via : Initial Calibration

Instrument: MSVOA\_V **LabSampleld**: VSTDCCC005

### **Manual IntegrationsAPPROVED**

Reviewed By :John Carlone 11/17/2021 Supervised By :Mahesh Dadoda 11/18/2021



### (20) 2-Butanone-d5 (S)

3.902min (+ 0.010) 10.64 ug/L

response 14287 Ion Exp% Act% 46.10 100.00 100.00 77.10 22.30 50.88# 0.00 0.00 0.00 0.00 0.00 0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_v\Data\VV111621\

Data File: VV023522.D

Acq On : 16 Nov 2021 10:01

Operator | : SY/MD Sample : VSTDCCC005

Misc : 25.0mL/MSVOA\_V/WATER ALS Vial : 2 Sample Multiplier: 1

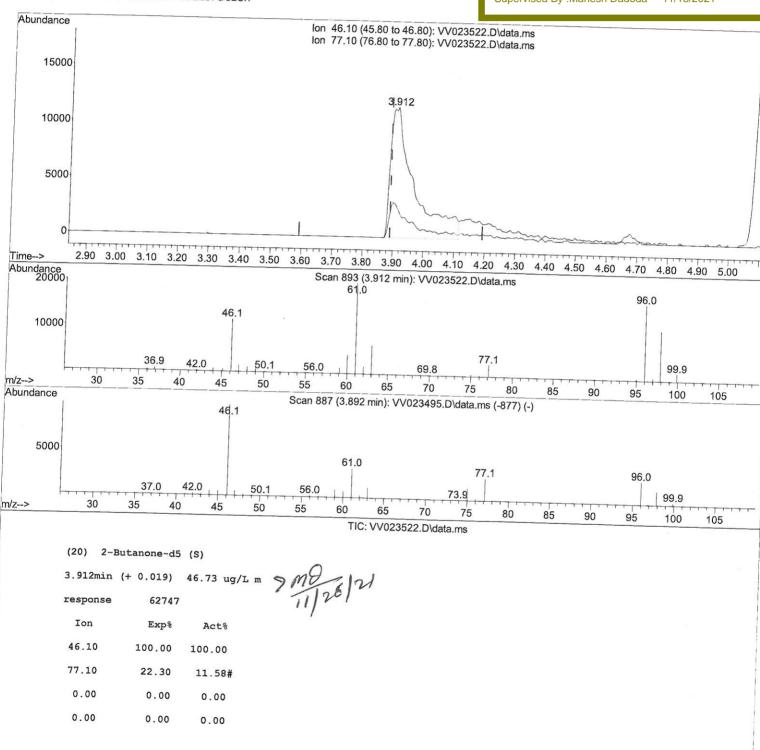
Quant Time: Nov 17 00:43:27 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Tue Nov 16 02:06:43 2021 Response via : Initial Calibration

Instrument: MSVOA\_V **LabSampleld**: VSTDCCC005

### **Manual IntegrationsAPPROVED**



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV111621\

Data File : VV023522.D

Acq On : 16 Nov 2021 10:01

Operator : SY/MD Sample : VSTDCCC005

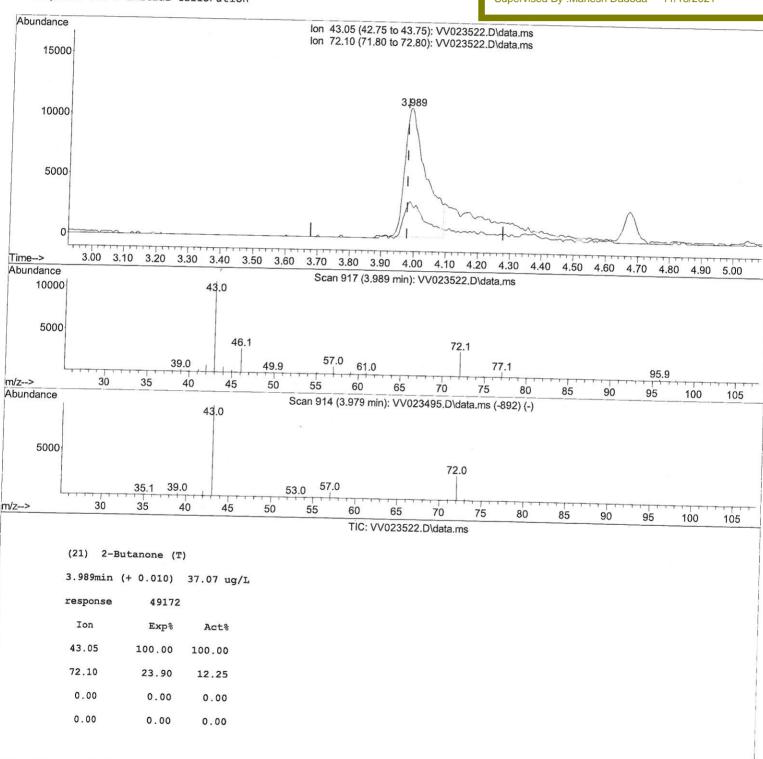
Misc : 25.0mL/MSVOA\_V/WATER
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Nov 17 00:43:27 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Tue Nov 16 02:06:43 2021 Response via : Initial Calibration Instrument: MSVOA\_V LabSampleId: VSTDCCC005

### **Manual IntegrationsAPPROVED**



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV111621\

Data File: VV023522.D

Acq On : 16 Nov 2021 10:01 Operator : SY/MD

Sample : VSTDCCC005

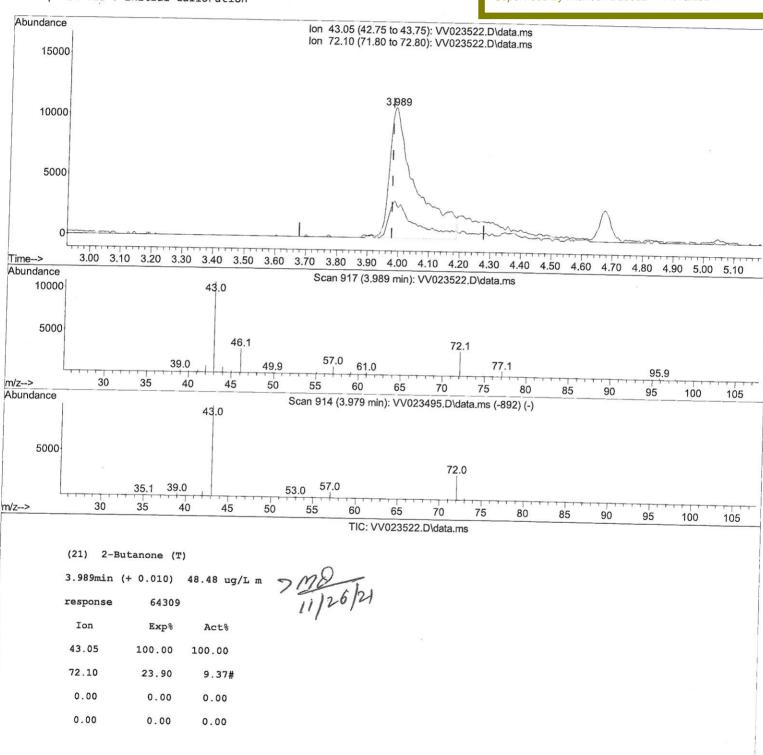
Misc : 25.0mL/MSVOA\_V/WATER
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Nov 17 00:43:27 2021

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Quant Title : TRACE VOA SFAM1.0 QLast Update : Tue Nov 16 02:06:43 2021 Response via : Initial Calibration Instrument: MSVOA\_V LabSampleId: VSTDCCC005

#### **Manual IntegrationsAPPROVED**



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV111621\

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Instrument : MSVOA\_V **LabSampleld**: VSTDCCC005

### **Manual IntegrationsAPPROVED**

- Wat						_	
Compound	R.T	. QIon	Response	Conc Ur	nits Dev(	Min)	
Internal Standards							
1) 1,4-Difluorobenze	no 5 61	6 114	124400	F 000	/1	0.00	
28) Chlorobenzene-d5	8.85		124408 123428		ug/L	0.00	
58) 1,4-Dichlorobenze			67780		ug/L ug/L	0.00	
30) 1, 1 510H101 05CH2C	11.24	112	07780	3.000	, ug/L	0.00	
System Monitoring Comp	ounds						
4) Vinyl Chloride-d3		4 65	39443	5.061	ug/L	0.00	
	.000 Range 40		Recover		101.200%		
7) Chloroethane-d5	1.568		32803	-	ug/L	0.00	
Spiked Amount 5	.000 Range 65	5 - 130	Recover		103.200%		
11) 1,1-Dichloroethene	e-d2 2.108	63	75177	5.153	ug/L	0.00	
Spiked Amount 5	.000 Range 60	9 - 125	Recover		103.000%		= who
20) 2-Butanone-d5	3.912	2 46	62747m			0.02	7 Mo /21
Spiked Amount 50.	.000 Range 40	- 130	Recover	ry =	93.460%		7/120101
24) Chloroform-d	4.349	84	79795	4.804	ug/L	0.00	/
	.000 Range 70	9 - 125	Recover	-y =	96.000%		
26) 1,2-Dichloroethane	e-d4 5.034	65	36601	4.900	ug/L	0.00	
	.000 Range 70		Recover	ry =	98.000%		
32) Benzene-d6	5.050		154095	4.866	ug/L	0.00	
	000 Range 70		Recover	-			
36) 1,2-Dichloropropan			44087	4.729	ug/L	0.00	
	000 Range 60		Recover		94.600%		
41) Toluene-d8	7.313		150780	5.081		0.00	
	000 Range 70		Recover		101.600%		
43) trans-1,3-Dichloro			17401	4.923		0.00	
•	000 Range 55		Recover		98.400%		
46) 2-Hexanone-d5	8.092		60535	46.544	A CONTRACTOR OF THE PARTY OF TH	0.00	
	000 Range 45		Recover		93.080%		
56) 1,1,2,2-Tetrachlor Spiked Amount 5.			32467	4.843	98 (57) (C. 10) (C. 10) (C. 10)	0.00	
66) 1,2-Dichlorobenzen		152	Recover		96.800%	0.00	
and the second s	000 Range 80		56076	4.969		0.00	
Spiked Amount 3.	ooo kange oo	- 120	Recover	y =	99.400%		
Target Compounds					Qval	110	
2) Dichlorodifluorome	thane 1.127	85	53335	4.397		97	
3) Chloromethane	1.240		48216	4.675	_	93	
<li>5) Vinyl chloride</li>	1.310		49100	4.767		100	
6) Bromomethane	1.523		25497	3.872		97	
<ol><li>8) Chloroethane</li></ol>	1.584		29451	4.954		97	
<ol><li>Trichlorofluorometh</li></ol>		101	75766	4.895		99	
10) 1,1,2-Trichloro-1,2	2,2 2.118	101	39571	5.079		96	
12) 1,1-Dichloroethene	2.118	96	35926	4.843		97	$\cap$
13) Acetone	2.191	43	40354m	49.186		2	monila
14) Carbon disulfide	2.294	76	118806	4.244		100	11/26/4
15) Methyl Acetate	2.436	43	10723	4.618	ug/L	100	
16) Methylene chloride	2.507	84	41641	3.846	553	98	
17) Methyl tert-butyl E		73	79742	4.883		96	
18) trans-1,2-Dichloroe	ethene 2.761	96	41623	4.564		96	
19) 1,1-Dichloroethane	3.188	63	73145	4.750		98	-0
21) 2-Butanone	3.989	43	64309m	48.482			I'M
22) cis-1,2-Dichloroeth		96	41992	4.784		89	71/26/21
23) Bromochloromethane	4.249	128	19291	4.766	ug/L	83	/ /

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV111621\

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Acq On : 16 Nov 2021 10:01
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Sample : VSTDCCC005
Misc : 25.0mL/MSVOA\_V/WATER ALS Vial : 2 Sample Multiplier: 1

Quant Time: Nov 17 00:43:27 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Tue Nov 16 02:06:43 2021 Response via : Initial Calibration

Instrument: MSVOA\_V **LabSampleld**: VSTDCCC005

### **Manual IntegrationsAPPROVED**

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
25) Chloroform	4.375	 83	78520	4.704 (1	
27) 1,2-Dichloroethane	5.130		43396	4.784 ug/L	99
29) 1,1,1-Trichloroethane	4.606	97	71794	4.971 ug/L	97
30) Cyclohexane	4.674	56	60859	4.789 ug/L	99
31) Carbon tetrachloride	4.825	117	65262	4.531 ug/L	98
33) Benzene	5.098	78	163690	4.847 ug/L	96
34) Trichloroethene	5.915	95	44664	4.745 ug/L	100
35) Methylcyclohexane	6.130	83	66384	4.868 ug/L 4.584 ug/L	96
37) 1,2-Dichloropropane	6.172	63	38834		96
38) Bromodichloromethane	6.510	83	52406	4.822 ug/L 4.856 ug/L	99
39) cis-1,3-Dichloropropene	7.027	75	55397	4.782 ug/L	96
40) 4-Methyl-2-pentanone	7.227	43	204464	54.739 ug/L	99
42) Toluene	7.387	91	186620	5.058 ug/L	98 98
44) trans-1,3-Dichloropropene		75	47101	4.900 ug/L	98 97
45) 1,1,2-Trichloroethane	7.838	97	29001	5.012 ug/L	97
47) Tetrachloroethene	7.976	164	39234	4.935 ug/L	97
48) 2-Hexanone	8.140	43	150551	57.520 ug/L	99
49) Dibromochloromethane	8.246	129	37208	5.075 ug/L	93
50) 1,2-Dibromoethane	8.352	107	26696	4.978 ug/L	97
51) Chlorobenzene	8.879	112	116266	4.740 ug/L	99
52) Ethylbenzene	9.011	91	191141	4.911 ug/L	97
53) m,p-xylene	9.136	106	75190	4.923 ug/L	95
54) o-xylene	9.542	106	71597	4.997 ug/L	97
55) Styrene	9.561	104	127965	5.213 ug/L	98
57) 1,1,2,2-Tetrachloroethane	10.239	83	30743	4.849 ug/L	96
59) Bromoform	9.731	173	19970	4.933 ug/L	99
60) Isopropylbenzene	9.931	105	193421	4.973 ug/L	99
61) 1,2,3-Trichloropropane	10.272	75	23014	5.111 ug/L	98
62) 1,3,5-Trimethylbenzene	10.538	105	159109	4.933 ug/L	99
63) 1,2,4-Trimethylbenzene	10.915	105	161570	5.033 ug/L	98
64) 1,3-Dichlorobenzene	11.181	146	99812	5.022 ug/L	99
65) 1,4-Dichlorobenzene	11.271	146	99231	4.889 ug/L	99
67) 1,2-Dichlorobenzene	11.641	146	90725	5.102 ug/L	99
68) 1,2-Dibromo-3-chloropr	12.429	75	4423	4.611 ug/L	88
69) 1,3,5-Trichlorobenzene	12.644	180	73570	4.728 ug/L	98
70) 1,2,4-trichlorobenzene	13.262	180	56940	4.570 ug/L	99
71) Naphthalene	13.503	128	76534	4.165 ug/L	100
72) 1,2,3-Trichlorobenzene	13.744	180	51111	4.688 ug/L	98

<sup>(#) =</sup> qualifier out of range (m) = manual integration (+) = signals summed