Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110421\

Data File : VV023200.D

Acq On : 04 Nov 2021 13:32

Operator : SY/MD Sample : VSTD01050

Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Nov 08 12:56:42 2021

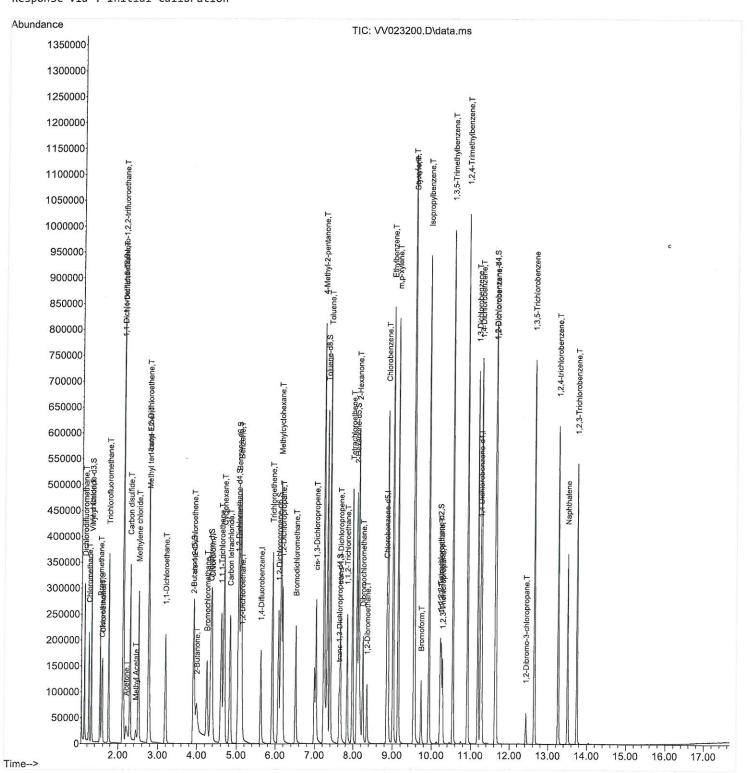
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0

QLast Update : Mon Nov 08 12:49:02 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110421\

Data File: VV023200.D

Acq On : 04 Nov 2021 13:32

Operator : SY/MD Sample : VSTD01050

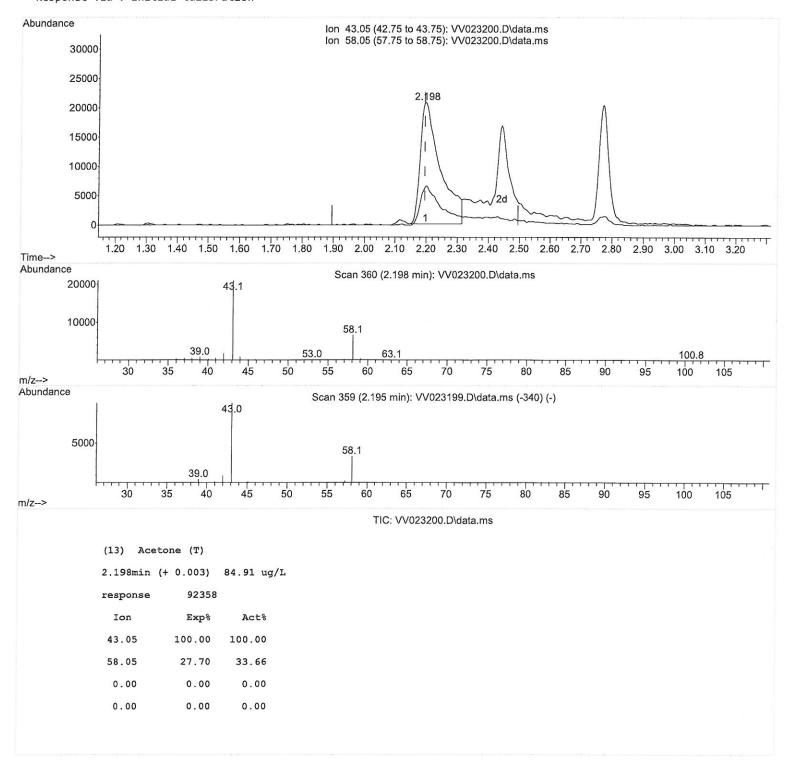
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Nov 08 12:56:42 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Mon Nov 08 12:49:02 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleId : VSTD010250

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110421\

Data File: VV023200.D

Acq On : 04 Nov 2021 13:32

Operator : SY/MD Sample : VSTD01050

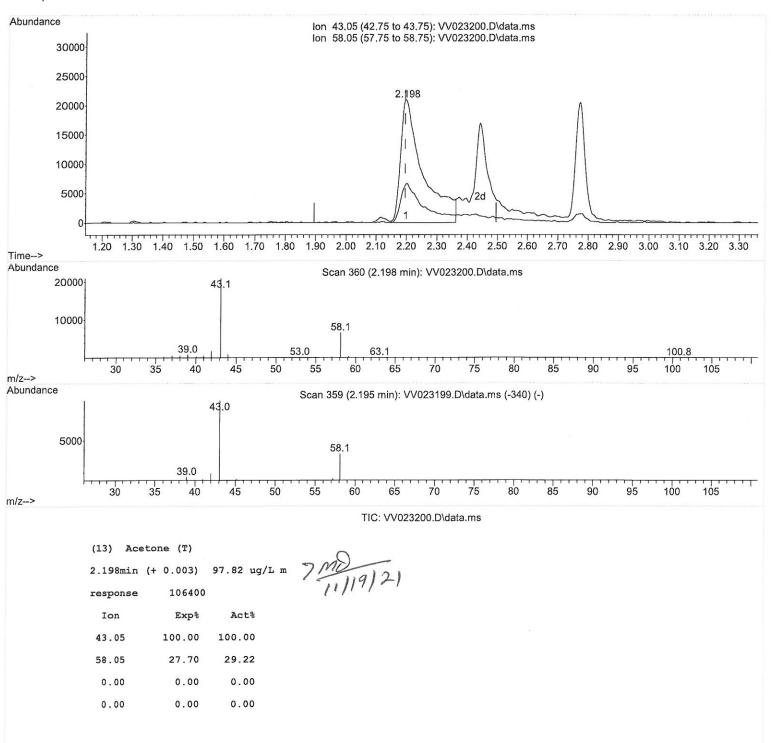
Misc : 25.0mL/MSVOA_V/WATER ALS Vial : 6 Sample Multiplier: 1

Quant Time: Nov 08 12:56:42 2021

 $\label{eq:Quant_Method} {\tt Quant_Method}: {\tt Z:\voasrv\hPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M}$

Quant Title : TRACE VOA SFAM1.0 QLast Update : Mon Nov 08 12:49:02 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleId : VSTD010250

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\W110421\

Data File: VV023200.D

Acq On : 04 Nov 2021 13:32

Operator : SY/MD Sample : VSTD01050

Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

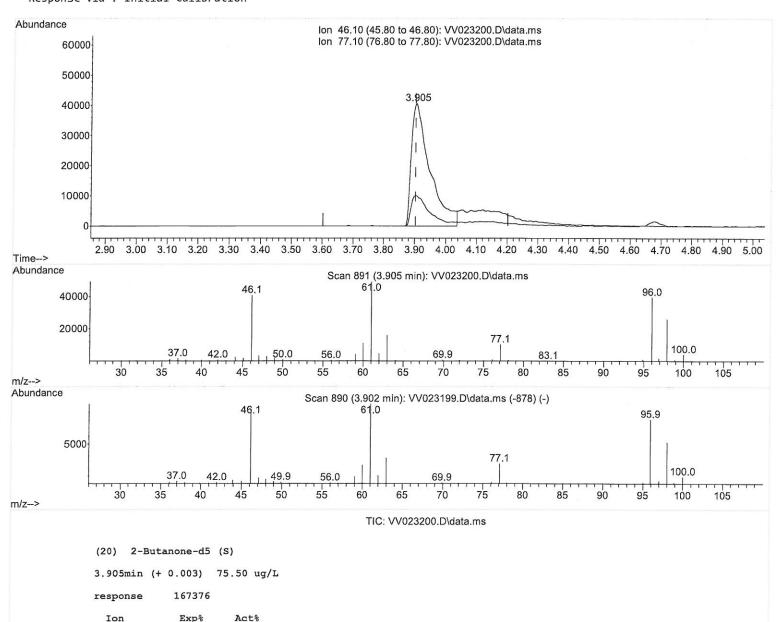
Quant Time: Nov 08 12:56:42 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Mon Nov 08 12:49:02 2021 Response via : Initial Calibration Instrument :
MSVOA_V
ClientSampleId :
VSTD010250

Manual IntegrationsAPPROVED

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



100.00

22.30

0.00

0.00

100.00

23.98

0.00

0.00

46.10

77.10

0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110421\

Data File: VV023200.D

Acq On : 04 Nov 2021 13:32

Operator : SY/MD Sample : VSTD01050

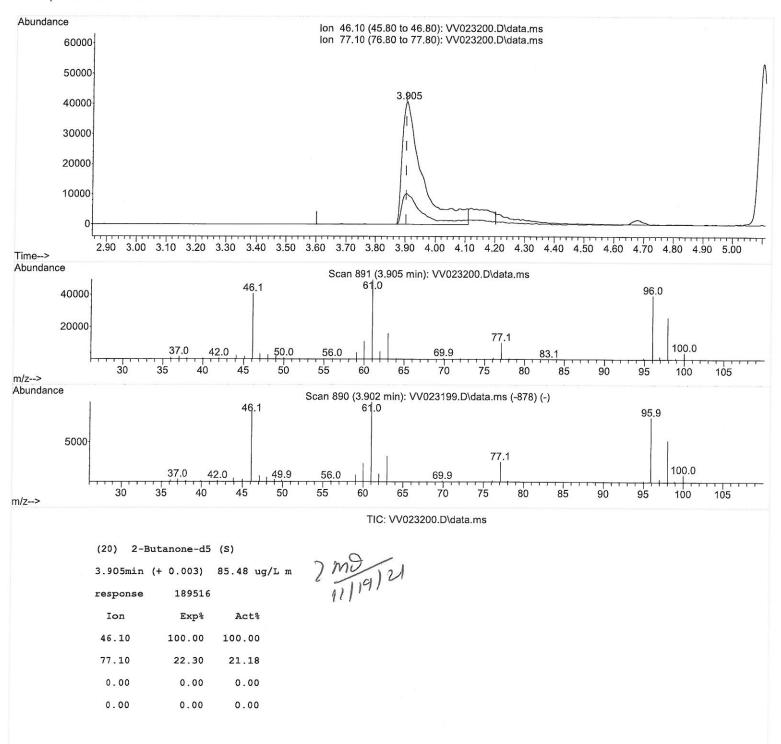
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Nov 08 12:56:42 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Mon Nov 08 12:49:02 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleld : VSTD010250

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110421\

Data File : VV023200.D

Acq On : 04 Nov 2021 13:32

Operator : SY/MD Sample : VSTD01050

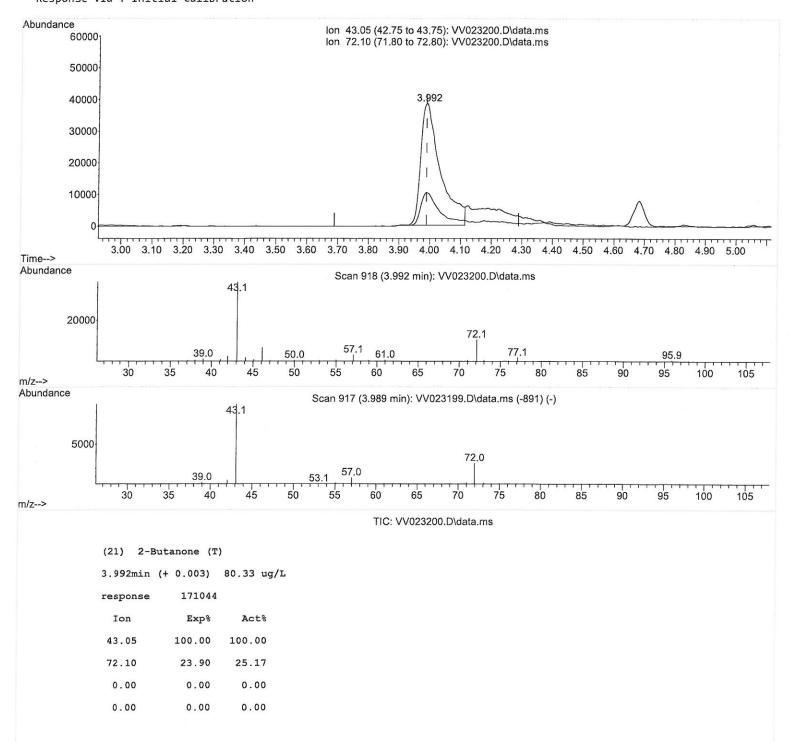
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Nov 08 12:56:42 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Mon Nov 08 12:49:02 2021 Response via : Initial Calibration Instrument:
MSVOA_V
ClientSampleld:
VSTD010250

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110421\

Data File: VV023200.D

Acq On : 04 Nov 2021 13:32

Operator : SY/MD Sample : VSTD01050

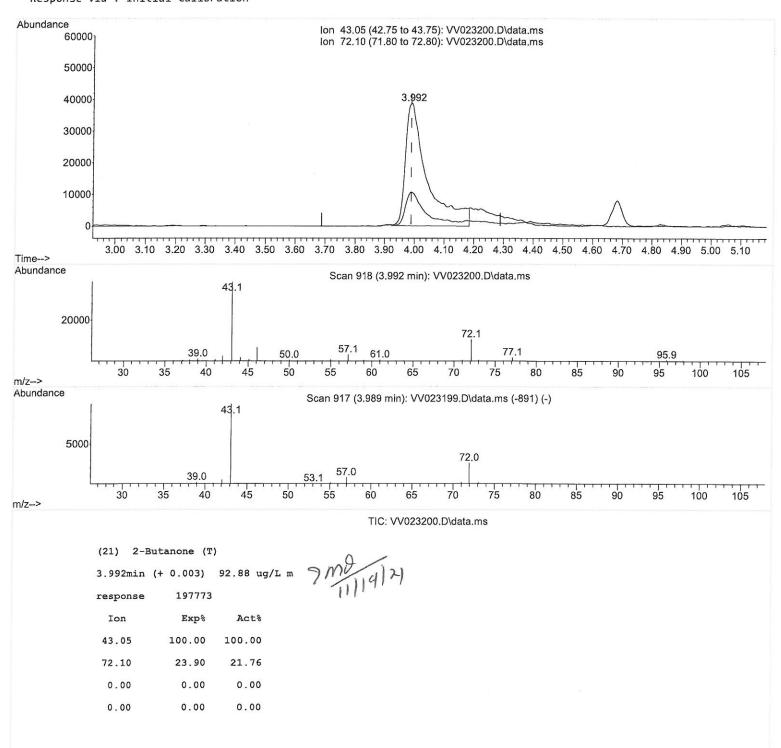
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Nov 08 12:56:42 2021

 $\label{lem:quant_method} {\tt Quant_Method} : {\tt Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M}$

Quant Title : TRACE VOA SFAM1.0 QLast Update : Mon Nov 08 12:49:02 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleId : VSTD010250

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110421\

Data File : VV023200.D

Acq On : 04 Nov 2021 13:32

Operator : SY/MD Sample : VSTD01050

Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Nov 08 12:56:42 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Mon Nov 08 12:49:02 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleId : VSTD010250

Manual IntegrationsAPPROVED

Compound	R.T.	QIon	Response	Conc Un:	its [Dev(Min)	
Internal Standards							
1) 1,4-Difluorobenzene	5.622	114	158159	5.00	ug/L	0.00	
28) Chlorobenzene-d5	8.853		154688	5.00 (0.00	
58) 1,4-Dichlorobenzene-d4	11.249		87875	5.00 t		0.00	
System Monitoring Compounds	1 207	65	100031	7 20	71	0.00	
4) Vinyl Chloride-d3	1.307		100831	7.30 t	-	0.00	
7) Chloroethane-d5	1.568	69	81384	9.52 (_	0.00	Ω
11) 1,1-Dichloroethene-d2	2.111	63	189965	9.53 t 85.48 t		0.00	Mo a121
20) 2-Butanone-d5	3.905	46	189516m		•	0.00	11/11/11
24) Chloroform-d	4.352	84	217371 99207	9.67 ເ 9.37 ເ		0.00	• •
26) 1,2-Dichloroethane-d4 32) Benzene-d6	5.037 5.053	65 84	425187	9.41 u	0	0.00 0.00	
	6.072	67	121846	8.76 (-	0.00	
<pre>36) 1,2-Dichloropropane-d6 41) Toluene-d8</pre>	7.317	98	412690	10.17 t	-	0.00	
43) trans-1,3-Dichloroprop	7.625	79	48883	10.04		0.00	
46) 2-Hexanone-d5	8.091	63	203223	112.69		0.00	
	10.217	84	88814	9.25 t	100000000000000000000000000000000000000	0.00	
56) 1,1,2,2-Tetrachloroeth 66) 1,2-Dichlorobenzene-d4	11.625	152	152178	9.71 u	-	0.00	
ob) 1,2-bichiolobenzene-u4	11.025	132	132170	5.71 0	45/ L	0.00	
Target Compounds						Qvalue	
Dichlorodifluoromethane	1.127	85	164264	16.37 u		98	
Chloromethane	1.240	50	137279	12.66 u	0.000	97	
Vinyl chloride	1.310	62	137512	12.28 L		100	
6) Bromomethane	1.523	94	89326	15.78 u	٠.	98	
8) Chloroethane	1.587	64	80421	13.73 u	•	99	
Trichlorofluoromethane	1.754	101	207734	13.74 u		99	
10) 1,1,2-Trichloro-1,2,2	2.117	101	105273	12.21 u	3 TO 10 10 10 10 10 10 10 10 10 10 10 10 10	98	^
12) 1,1-Dichloroethene	2.121	96	100432	12.40 u	_	88	MOTIL
13) Acetone	2.198	43	106400m	97.82 u		/	11/19/4
14) Carbon disulfide	2.294	76	382946	17.37 u	1770	100	1. 1
15) Methyl Acetate	2.442	43	39705	9.60 u			
16) Methylene chloride	2.510	84	122148	13.73 u	3753677	98	
17) Methyl tert-butyl Ether	2.773	73	235124	12.16 u		96	
18) trans-1,2-Dichloroethene	2.764	96	125612	14.53 u	-	97	Ω
19) 1,1-Dichloroethane	3.195	63	212399	13.37 u	-	97	M0 12-1
21) 2-Butanone	3.992	43	197773m	92.88 u		# 02	11/19/0
22) cis-1,2-Dichloroethene	3.918	96	125185	12.77 u			1. [
23) Bromochloromethane	4.252	128	56559	12.71 u			
25) Chloroform	4.381	83	222834	10.62 u	_	97	
27) 1,2-Dichloroethane	5.137	62	120643	10.90 u		100	
29) 1,1,1-Trichloroethane	4.612 4.683	97 56	206339 196202	11.80 u 12.86 u		99 97	
30) Cyclohexane31) Carbon tetrachloride	4.831	117	188851	12.56 u		96	
33) Benzene34) Trichloroethene	5.104 5.918	78 95	489263 128093	11.75 u 12.25 u		100 98	
35) Methylcyclohexane	6.137	83	212089	14.16 u		95	
37) 1,2-Dichloropropane	6.175	63	114193	10.88 u	_	98	
38) Bromodichloromethane	6.513	83	151622	10.38 u		99	
39) cis-1,3-Dichloropropene	7.030	75	169567	12.98 u		97	
40) 4-Methyl-2-pentanone	7.230	43	594553	12.38 u	-	97	
42) Toluene	7.391	91	546010	12.94 u		97	
44) trans-1,3-Dichloropropene	7.654	75	144490	13.37 u	_	97	
, 2,5 5261201 091 096110				, u	۰, -		

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110421\

Data File : VV023200.D

Acq On : 04 Nov 2021 13:32 Operator : SY/MD

: VSTD01050 Sample

: 25.0mL/MSVOA_V/WATER Misc ALS Vial : 6 Sample Multiplier: 1

Quant Time: Nov 08 12:56:42 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Mon Nov 08 12:49:02 2021 Response via : Initial Calibration

Instrument : MSVOA_V ClientSampleId : VSTD010250

Manual IntegrationsAPPROVED

	Compound	R.T.	QIon	Response	Conc Units Dev	/(Min)
45)	1,1,2-Trichloroethane	7.841	97	79628	10.85 ug/L	98
47)	Tetrachloroethene	7.979	164	111767	12.65 ug/L	99
48)	2-Hexanone	8.143	43	423283	115.44 ug/L	97
49)	Dibromochloromethane	8.249	129	105807	12.11 ug/L	99
50)	1,2-Dibromoethane	8.355	107	76710	11.61 ug/L	98
51)	Chlorobenzene	8.882	112	342626	12.36 ug/L	99
52)	Ethylbenzene	9.014	91	582458	13.91 ug/L	98
53)	m,p-xylene	9.140	106	231875	13.81 ug/L	99
54)	o-xylene	9.545	106	219013	13.85 ug/L	99
55)	Styrene	9.561	104	383639	14.05 ug/L	98
57)	1,1,2,2-Tetrachloroethane	10.243	83	90109	10.98 ug/L	99
59)	Bromoform	9.731	173	58414	11.41 ug/L	99
60)	Isopropylbenzene	9.934	105	595966	13.41 ug/L	99
61)	1,2,3-Trichloropropane	10.275	75	64000	10.35 ug/L	97
62)	1,3,5-Trimethylbenzene	10.538	105	506730	14.19 ug/L	99
63)	1,2,4-Trimethylbenzene	10.915	105	518614	14.56 ug/L	99
64)	1,3-Dichlorobenzene	11.181	146	292021	12.49 ug/L	99
65)	1,4-Dichlorobenzene	11.275	146	289442	12.18 ug/L	99
67)	<pre>1,2-Dichlorobenzene</pre>	11.644	146	258233	11.86 ug/L	100
68)	1,2-Dibromo-3-chloropr	12.429	75	14571	11.94 ug/L	93
69)	1,3,5-Trichlorobenzene	12.644	180	226497	12.72 ug/L	98
70)	1,2,4-trichlorobenzene	13.262	180	183625	13.78 ug/L	99
71)	Naphthalene	13.503	128	284595	14.05 ug/L	99
72)	1,2,3-Trichlorobenzene	13.744	180	164548	13.30 ug/L	98

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