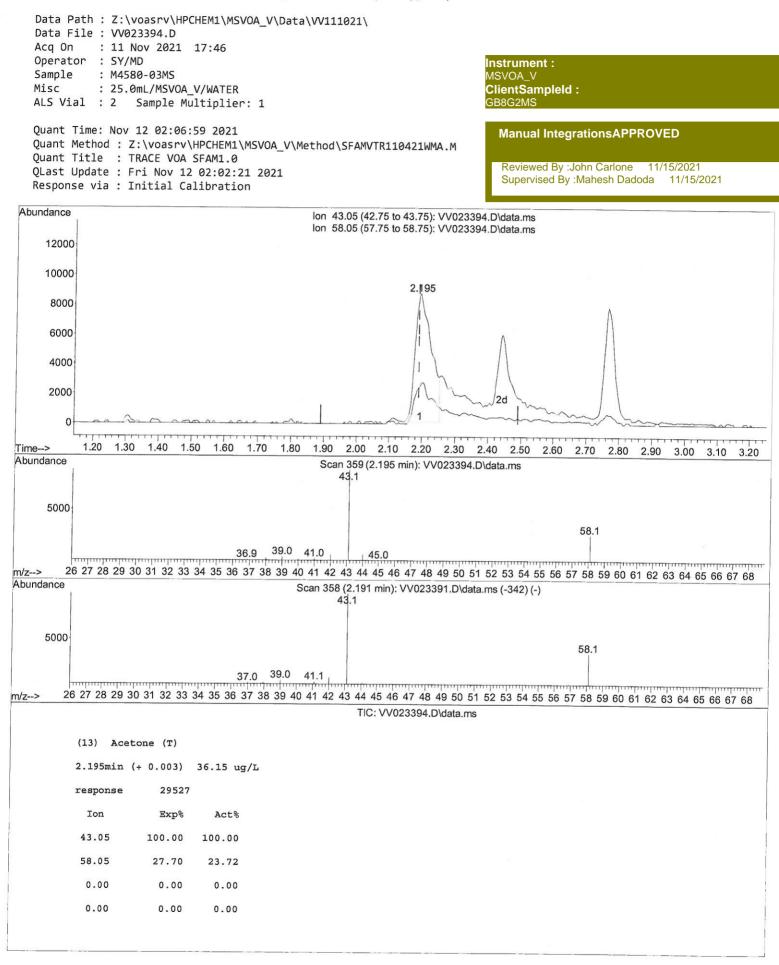
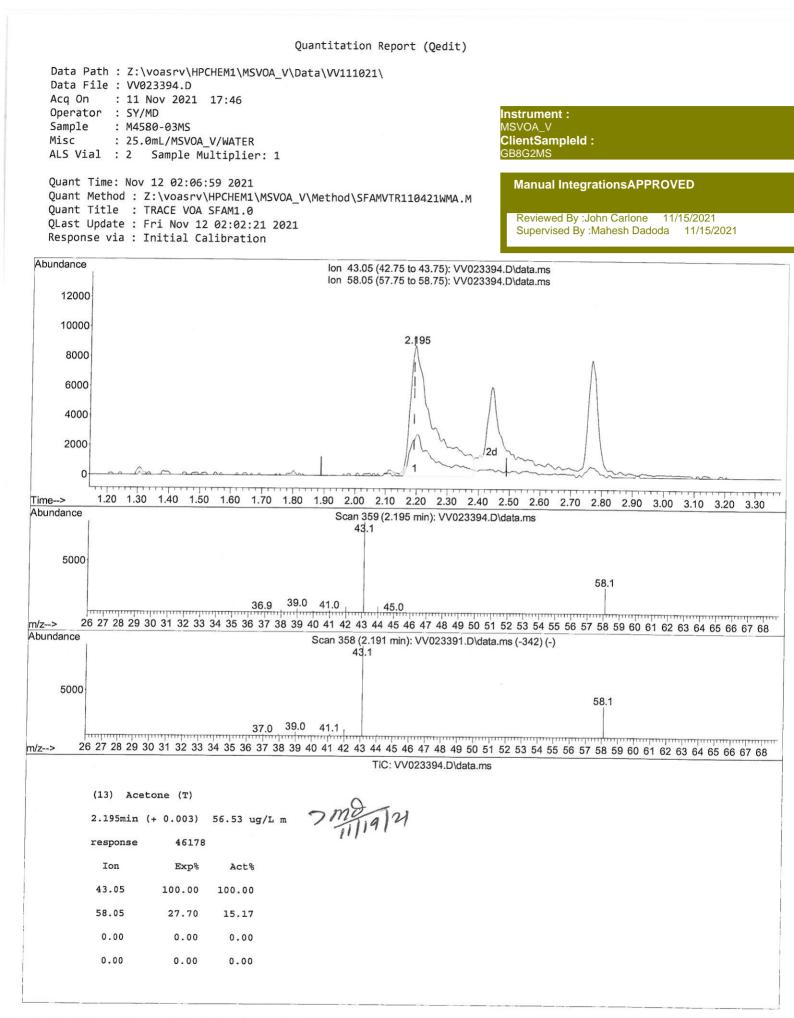


SFAMVTR110421WMA.M Fri Nov 12 02:19:17 2021





Data Path : Z:\voasrv\HPCHEM Data File : VV023394.D Acq On : 11 Nov 2021 17: Operator : SY/MD Sample : M4580-03MS Misc : 25.0mL/MSVOA_V/W ALS Vial : 2 Sample Multi	46 ATER	V111021\	Instrument : MSVOA_V ClientSampleId : GB8G2MS
Quant Time: Nov 12 02:06:59 Quant Method : Z:\voasrv\HPC Quant Title : TRACE VOA SFA QLast Update : Fri Nov 12 02 Response via : Initial Calibu	HEM1\MSVOA_V\Met M1.0 :02:21 2021	hod\SFAMVTR110421WMA.M	Manual IntegrationsAPPROVED Reviewed By :John Carlone 11/15/2021 Supervised By :Mahesh Dadoda 11/15/2021
Compound	R.T. QIon	Response Conc Units De	ev(Min)
Internal Standards 1) 1,4-Difluorobenzene 28) Chlorobenzene-d5 58) 1,4-Dichlorobenzene-d4	5.619 114 8.854 117 11.249 152	123860 5.000 ug/L 122448 5.000 ug/L 64352 5.000 ug/L	0.00 0.00 0.00
System Monitoring Compounds 4) Vinyl Chloride-d3 Spiked Amount 5.000	1.304 65 Range 40 - 136	,	0.00 99%
 7) Chloroethane-d5 Spiked Amount 5.000 11) 1,1-Dichloroethene-d2 	1.568 69 Range 65 - 130 2.108 63	64178 4.418 ug/L	0.00 00% 0.00
Spiked Amount 5.000 20) 2-Butanone-d5 Spiked Amount 50.000	Range 60 - 125 3.915 46 Range 40 - 136	48966 36.629 ug/L Recovery = 73.26	0.00
<pre>24) Chloroform-d Spiked Amount 5.000 26) 1,2-Dichloroethane-d4</pre>	4.349 84 Range 70 - 125 5.037 65	32684 4.395 ug/L	0.00 0% 0.00
Spiked Amount 5.000 32) Benzene-d6 Spiked Amount 5.000	Range 70 - 130 5.053 84 Range 70 - 125	133816 4.259 ug/L Recovery = 85.20	0.00 0%
 36) 1,2-Dichloropropane-d6 Spiked Amount 5.000 41) Toluene-d8 Spiked Amount 5.000 	6.069 67 Range 60 - 140 7.317 98	129410 4.396 ug/L	0.00
Spiked Amount 5.000 43) trans-1,3-Dichloroprop. Spiked Amount 5.000 46) 2-Hexanone-d5	Range 55 - 130	15498 4.419 ug/L Recovery = 88.40	0.00 0%
Spiked Amount 50.000 56) 1,1,2,2-Tetrachloroeth. Spiked Amount 5.000	8.092 63 Range 45 - 130 10.217 84 Range 65 - 120	28082 4.222 ug/L	0.00
66) 1,2-Dichlorobenzene-d4 Spiked Amount 5.000	11.625 152 Range 80 - 120	Recovery = 84.400 48054 4.485 ug/L Recovery = 89.600	0.00
Target Compounds			value
 2) Dichlorodifluoromethane 3) Chloromethane 	1.127 85 1.240 50	56902 4.711 ug/L 51860 5.050 ug/L	100 98
5) Vinyl chloride	1.311 62	51423 5.014 ug/L	98
6) Bromomethane8) Chloroethane	1.520 94	29773 4.542 ug/L	97
9) Trichlorofluoromethane	1.584 64 1.751 101	29246 4.942 ug/L 77804 5.049 ug/L	96 99
10) 1,1,2-Trichloro-1,2,2		37577 4.844 ug/L	94
12) 1,1-Dichloroethene	2.118 96	37356 5.058 ug/L	86 M2 121
13) Acetone	2.195 43	46178m 56.534 ug/L	511191
14) Carbon disulfide 15) Methyl Acetate	2.294 76 2.442 43	135405 4.858 ug/L 11406 4.934 ug/L	99 <i>111</i> 97
16) Methylene chloride	2.507 84	43949 4.077 ug/L	97
17) Methyl tert-butyl Ether	2.770 73	84443 5.194 ug/L	95
18) trans-1,2-Dichloroethene		46249 5.094 ug/L	99
19) 1,1-Dichloroethane 21) 2-Butanone	3.188 63 3.995 43	78490 5.120 ug/L 53960 40.860 ug/L	97
22) cis-1,2-Dichloroethene	3.915 96	46119 5.278 ug/L #	97 92
23) Bromochloromethane	4.249 128	20916 5.191 ug/L	83

SFAMVTR110421WMA.M Fri Nov 12 02:19:15 2021

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(QT Reviewed)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111021\ Data File : VV023394.D Acq On : 11 Nov 2021 17:46 Operator : SY/MD Sample : M4580-03MS Misc : 25.0mL/MSVOA_V/WATER ALS Vial : 2 Sample Multiplier: 1

Quant Time: Nov 12 02:06:59 2021 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M Quant Title : TRACE VOA SFAM1.0 QLast Update : Fri Nov 12 02:02:21 2021 Response via : Initial Calibration

Instrument : MSVOA_V ClientSampleId : GB8G2MS

Manual IntegrationsAPPROVED

Reviewed By : John Carlone 11/15/2021 Supervised By :Mahesh Dadoda 11/15/2021

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
25) Chloroform	4.375	83	85367	5.224 ug/L	 97
27) 1,2-Dichloroethane	5.134		44457	5.115 ug/L	100
29) 1,1,1-Trichloroethane	4.609	97	76184	5.123 ug/L	100
30) Cyclohexane	4.680	56	66829	5.015 ug/L	98
31) Carbon tetrachloride	4.828	117	69084	5.172 ug/L	97
33) Benzene	5.101	78	176408	5.154 ug/L	100
34) Trichloroethene	5.915	95	47174	5.183 ug/L	98
35) Methylcyclohexane	6.133	83	70812	4.929 ug/L	97
37) 1,2-Dichloropropane	6.175	63	40519	5.071 ug/L	98
38) Bromodichloromethane	6.510	83	55379	5.172 ug/L	95
39) cis-1,3-Dichloropropene	7.027	75	58694	5.107 ug/L	100
40) 4-Methyl-2-pentanone	7.230	43	217802	58.776 ug/L	98
42) Toluene	7.387	91	197039	5.383 ug/L	98
44) trans-1,3-Dichloropropene	7.651	75	48863	5.124 ug/L	100
45) 1,1,2-Trichloroethane	7.841	97	29205	5.087 ug/L	96
47) Tetrachloroethene	7.976	164	115678	14.666 ug/L	98
48) 2-Hexanone	8.143	43	157866	60.798 ug/L	99
49) Dibromochloromethane	8.246	129	38978	5.359 ug/L	95
50) 1,2-Dibromoethane	8.355	107	28334	5.326 ug/L #	98
51) Chlorobenzene	8.883	112	123206	5.064 ug/L	99
52) Ethylbenzene	9.011	91	201607	5.222 ug/L	97
53) m,p-xylene	9.140	106	79587	5.253 ug/L	99
54) o-xylene	9.545	106	76145	5.357 ug/L	100
55) Styrene	9.561	104	131129	5.385 ug/L	100
57) 1,1,2,2-Tetrachloroethane	10.243	83	33941	5.396 ug/L	98
59) Bromoform	9.731	173	20950	5.450 ug/L	99
60) Isopropylbenzene	9.931	105	201330	5.452 ug/L	100
61) 1,2,3-Trichloropropane	10.275	75	24429	5.715 ug/L	98
62) 1,3,5-Trimethylbenzene	10.538	105	162939	5.321 ug/L	98
63) 1,2,4-Trimethylbenzene	10.915	105	169513	5.562 ug/L	99
64) 1,3-Dichlorobenzene	11.182	146	100341	5.318 ug/L	98
65) 1,4-Dichlorobenzene		146	98436	5.108 ug/L	98
67) 1,2-Dichlorobenzene	11.645	146	91804	5.437 ug/L	98
68) 1,2-Dibromo-3-chloropr	12.429	75	5011	5.502 ug/L	86
69) 1,3,5-Trichlorobenzene	12.645	180	74139	5.018 ug/L	97
70) 1,2,4-trichlorobenzene	13.262	180	58404	4.937 ug/L	97
71) Naphthalene	13.503	128	84866	4.865 ug/L	99
72) 1,2,3-Trichlorobenzene	13.744	180	52515	5.073 ug/L	99

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