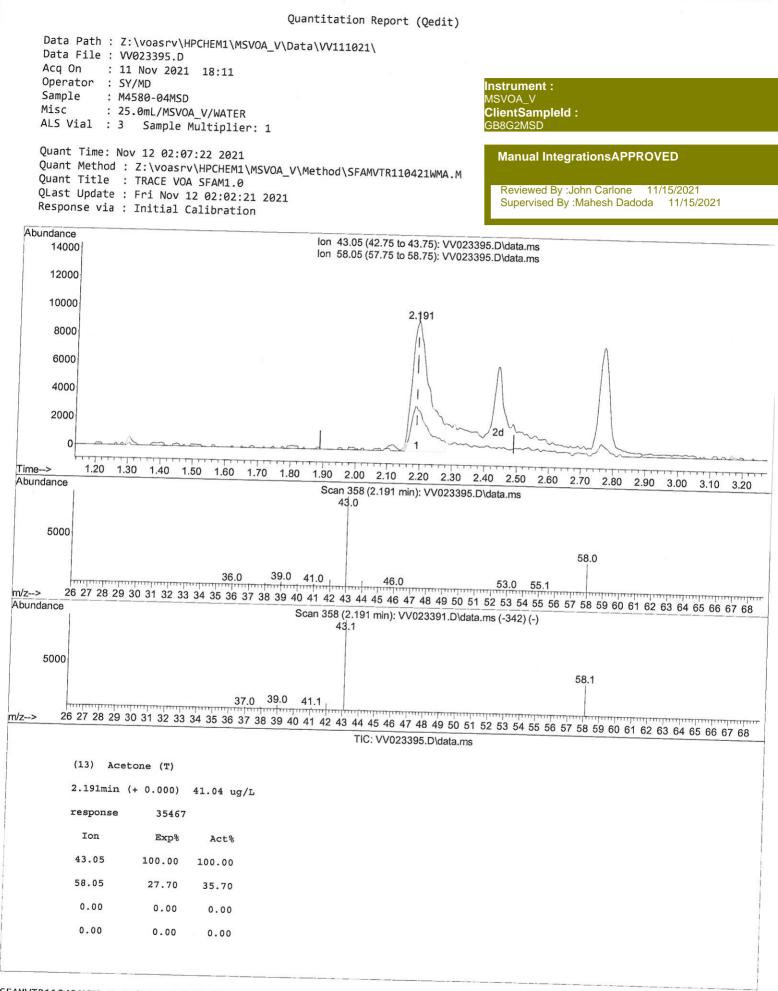
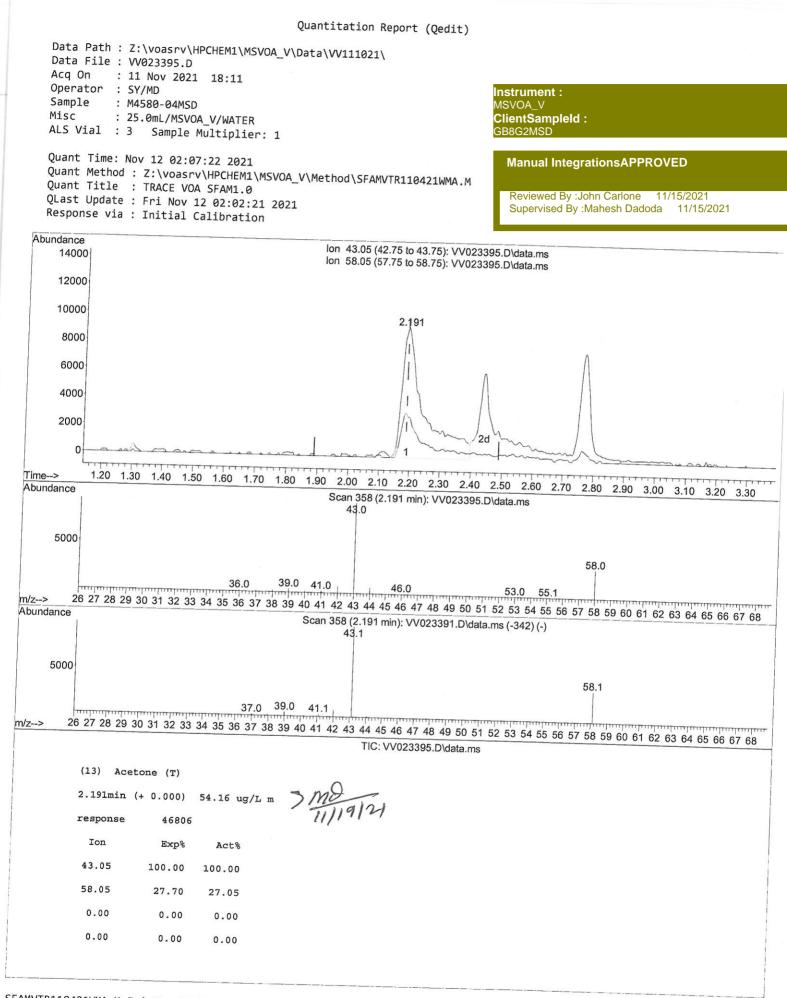


SFAMVTR110421WMA.M Fri Nov 12 02:24:23 2021



SFAMVTR110421WMA.M Fri Nov 12 02:23:18 2021



	Quant	(QI REVI	ewea
Data Path : Z:\voasrv\HPCHEM:	1\MSVOA V\Data\	W111021\	
Data File : VV023395.D			
Acq On : 11 Nov 2021 18:1	11		
Operator : SY/MD	Instrument : MSVOA_V		
Sample : M4580-04MSD Misc : 25.0mL/MSVOA_V/WA	ClientSampleId :		
ALS Vial : 3 Sample Multip			GB8G2MSD
Quant Time: Nov 12 02:07:22 2 Quant Method : Z:\voasrv\HPCF			Manual IntegrationsAPPROVED
Quant Title : TRACE VOA SFAM	11.0	CHOU (SPANN RI10421WMA.M	
QLast Update : Fri Nov 12 02: Response via : Initial Calibr	02:21 2021		Reviewed By :John Carlone 11/15/2021 Supervised By :Mahesh Dadoda 11/15/2021
Compound	R.T. QIO	n Response Conc Units Dev	(Min)
Internal Standards			
1) 1,4-Difluorobenzene	5.619 114		0.00
28) Chlorobenzene-d5	8.854 11		0.00
58) 1,4-Dichlorobenzene-d4	11.249 152	2 67559 5.000 ug/L	0.00
System Monitoring Compounds			
4) Vinyl Chloride-d3	1.304 65		0.00
Spiked Amount 5.000 7) Chloroethane-d5	Range 40 - 13 1.564 69	· · · · · · · · · · · · · · · · · · ·	
Spiked Amount 5.000	Range 65 - 13		0.00 %
11) 1,1-Dichloroethene-d2	2.108 63		0.00
Spiked Amount 5.000	Range 60 - 12	5 Recovery = 81.400	%
20) 2-Butanone-d5 Spiked Amount 50.000	3.905 46 Range 40 - 13	=	, 0.00
24) Chloroform-d	4.352 84	, , , , , , , , , , , , , , , , , , , ,	° 0.00
Spiked Amount 5.000	Range 70 - 12		
26) 1,2-Dichloroethane-d4	5.034 65	-8, -	0.00
Spiked Amount 5.000 32) Benzene-d6	Range 70 - 13 5.050 84		
	Range 70 - 12		0.00
36) 1,2-Dichloropropane-d6	6.069 67	,	0.00
	Range 60 - 14		
41) Toluene-d8 Spiked Amount 5.000	7.317 98 Range 70 - 13		0.00
43) trans-1,3-Dichloroprop		,	0.00
	Range 55 - 130	0 Recovery = 81.600%	
46) 2-Hexanone-d5	8.092 63	59749 44.645 ug/L	0.00
Spiked Amount 50.000 56) 1,1,2,2-Tetrachloroeth	Range 45 - 130 . 10.217 84	,	
	Range 65 - 120		0.00
66) 1,2-Dichlorobenzene-d4	11.625 152	46706 4.152 ug/L	0.00
Spiked Amount 5.000	Range 80 - 120	Recovery = 83.000%	
Target Compounds		Qva	lue
 2) Dichlorodifluoromethane 2) Chloromethane 	1.127 85	58184 4.553 ug/L	99
3) Chloromethane 5) Vinyl chloride	1.240 50 1.307 62	50455 4.644 ug/L	98
6) Bromomethane	1.307 62 1.519 94	50627 4.666 ug/L 30342 4.375 ug/L	99 100
8) Chloroethane	1.584 64	30163 4.817 ug/L	97
9) Trichlorofluoromethane	1.751 101	76872 4.715 ug/L	99
<pre>10) 1,1,2-Trichloro-1,2,2 12) 1,1-Dichloroethene</pre>		38773 4.724 ug/L	97
13) Acetone	2.118 96 2.191 43	36724 4.699 ug/L 46806m 54.159 ug/L	84 MO
14) Carbon disulfide	2.294 76	133641 4.532 ug/L	99 11/19/21
15) Methyl Acetate	2.442 43	9956 4.070 ug/L	98
16) Methylene chloride 17) Methyl tert-butyl Ether	2.507 84	43437 3.809 ug/L	94
18) trans-1,2-Dichloroethene	2.770 73 2.761 96	87058 5.061 ug/L 46240 4.813 ug/L	96 98
19) 1,1-Dichloroethane	3.188 63	79504 4.902 ug/L	97
21) 2-Butanone	3.986 43	57595 41.220 ug/L	97
22) cis-1,2-Dichloroethene	3.912 96	45662 4.939 ug/L #	88
23) Bromochloromethane	4.249 128	20409 4.787 ug/L	82

SFAMVTR110421WMA.M Fri Nov 12 02:24:22 2021

(QT Reviewed)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111021\ Data File : VV023395.D Acq On : 11 Nov 2021 18:11 Operator : SY/MD Sample : M4580-04MSD Misc : 25.0mL/MSVOA_V/WATER ALS Vial : 3 Sample Multiplier: 1

Quant Time: Nov 12 02:07:22 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 : GB8G2MSD

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	4.378	83	85240	4.930 ug/L	 94
27) 1,2-Dichloroethane	5.133	62	45772	4.977 ug/L	99
29) 1,1,1-Trichloroethane	4.609	97	76509	4.960 ug/L	100
30) Cyclohexane	4.677	56	66169	4.787 ug/L	96
Carbon tetrachloride	4.828	117	68390	4.937 ug/L	98
33) Benzene	5.101	78	176297	4.966 ug/L	100
34) Trichloroethene	5.915	95	46179	4.892 ug/L	95
35) Methylcyclohexane	6.130	83	69874	4.689 ug/L	96
37) 1,2-Dichloropropane	6.175	63	40180	4.848 ug/L	98
38) Bromodichloromethane	6.513	83	55587	5.005 ug/L	99
39) cis-1,3-Dichloropropene	7.031	75	58630	4.919 ug/L	96
40) 4-Methyl-2-pentanone	7.230	43	217475	56.581 ug/L	99
42) Toluene	7.387	91	197471	5.201 ug/L	96
44) trans-1,3-Dichloropropene	7.651	75	50442	5.100 ug/L	98
45) 1,1,2-Trichloroethane	7.841	97	29524	4.958 ug/L	98
47) Tetrachloroethene	7.976	164	116017	14.181 ug/L	98
48) 2-Hexanone	8.143	43	156555	58.128 ug/L	99
49) Dibromochloromethane	8.246	129	38395	5.089 ug/L	100
50) 1,2-Dibromoethane	8.355	107	27875	5.051 ug/L	100
51) Chlorobenzene	8.883	112	122554	4.856 ug/L	99
52) Ethylbenzene	9.014	91	200083	4.996 ug/L	99
53) m,p-xylene	9.140	106	79562	5.062 ug/L	96
54) o-xylene	9.545	106	75853	5.145 ug/L	100
55) Styrene	9.561	104	129428	5.124 ug/L	96
57) 1,1,2,2-Tetrachloroethane	10.243	83	32014	4.907 ug/L	98
59) Bromoform	9.735	173	21143	5.240 ug/L #	97
50) Isopropylbenzene	9.931	105	202099	5.213 ug/L	99
51) 1,2,3-Trichloropropane	10.275	75	23564	5.251 ug/L	98
2) 1,3,5-Trimethylbenzene	10.538	105	167725	5.218 ug/L	100
53) 1,2,4-Trimethylbenzene	10.915	105	168345	5.262 ug/L	99
4) 1,3-Dichlorobenzene	11.181	146	99479	5.022 ug/L	98
5) 1,4-Dichlorobenzene	11.275	146	100026	4.944 ug/L	98
57) 1,2-Dichlorobenzene	11.644	146	89815	5.067 ug/L	98
8) 1,2-Dibromo-3-chloropr	12.429	75	4872	5.096 ug/L	85
9) 1,3,5-Trichlorobenzene	12.644	180	74811	4.824 ug/L	98
0) 1,2,4-trichlorobenzene	13.262	180	57936	4.665 ug/L	100
1) Naphthalene	13.503	128	86804	4.740 ug/L	99
 1,2,3-Trichlorobenzene 	13.744	180	53577	4.930 ug/L	99

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