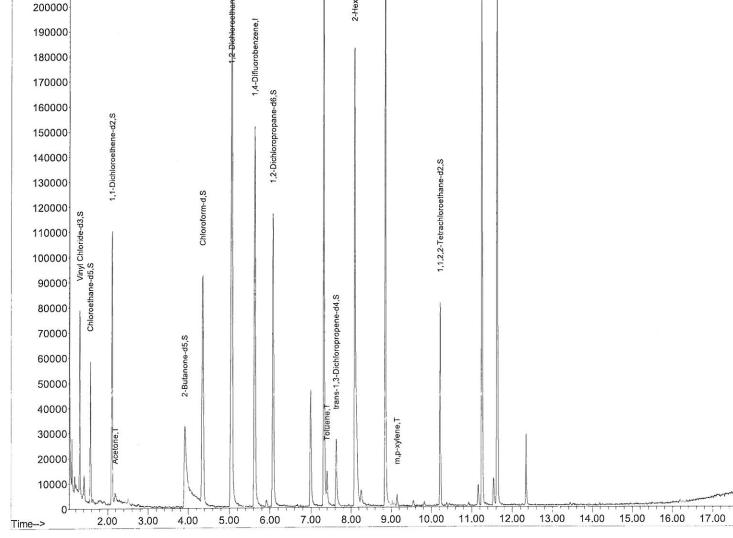
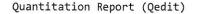
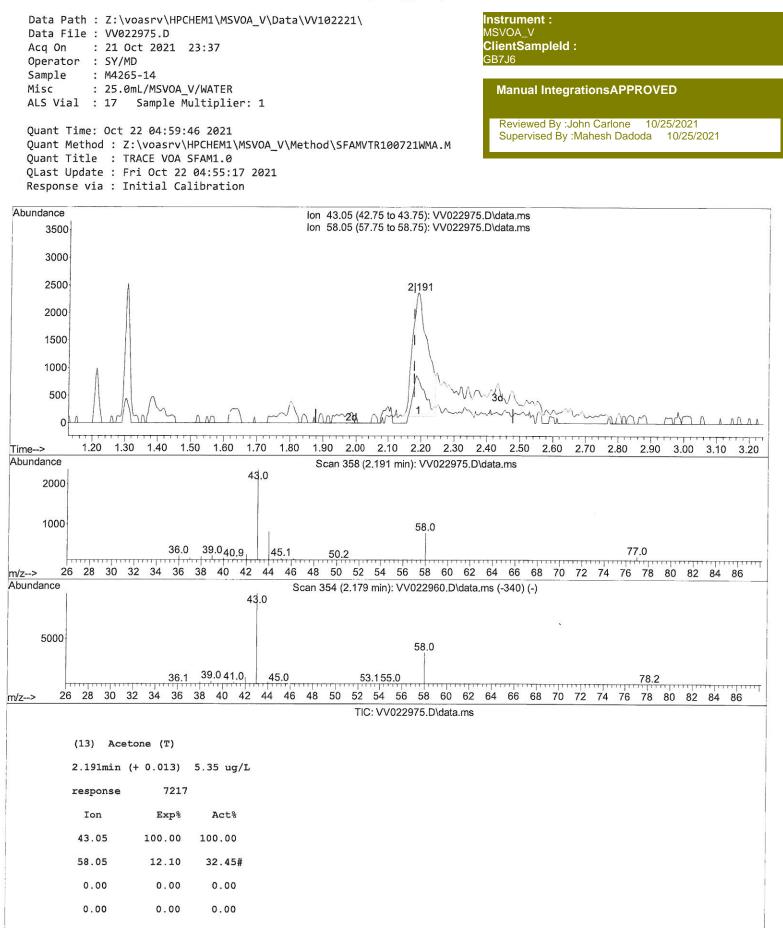
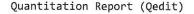
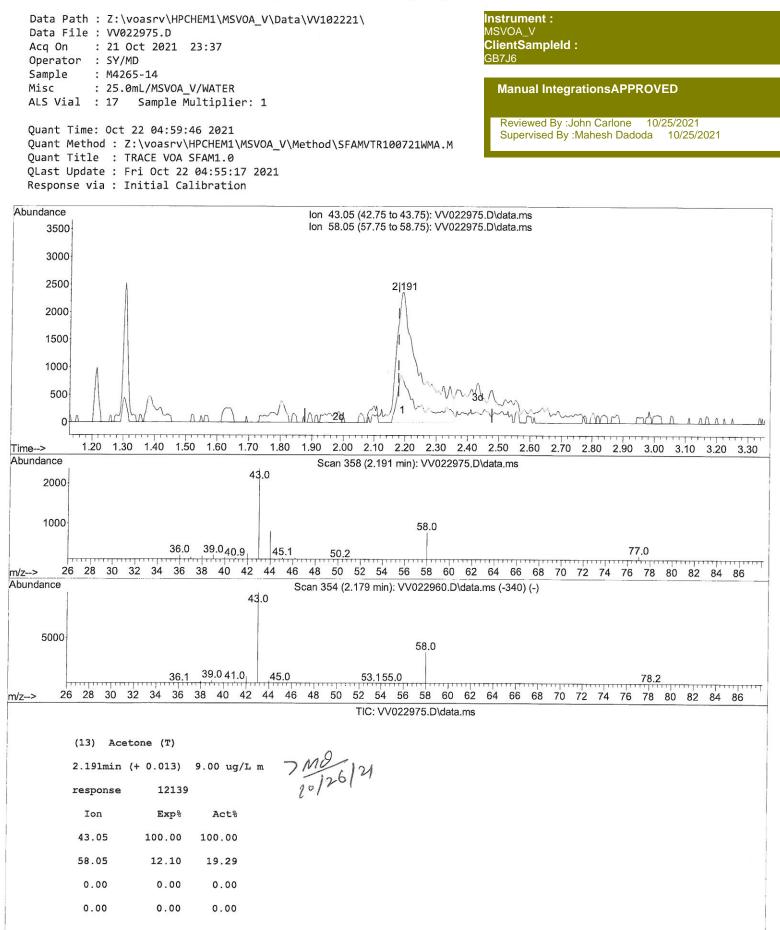
Quantitation Report (QT Reviewed)							
Data File : Acq On : Operator :		Instrument : MSVOA_V ClientSampleId : GB7J6					
Sample : M4265-14 Misc : 25.0mL/MSVOA_V/WATER ALS Vial : 17 Sample Multiplier: 1					Manual IntegrationsAPPROVED		
Quant Time: Oct 22 04:59:46 2021 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100721WMA.M Quant Title : TRACE VOA SFAM1.0					Reviewed By :John Carlone 10/25/2021 Supervised By :Mahesh Dadoda 10/25/2021		
QLast Update : Fri Oct 22 04:55:17 2021 Response via : Initial Calibration							
Abundance 280000			TIC: VV0	22975.D\data.ms			
270000							
260000					- d4, l		
250000					1,4-Dichlorobenzene-d4,1 robenzene-d4,S		
240000	ŵ		-d5.1		Dichlo		
230000	Benzene-d6, S	-oluene-d8,S	d5,S Chlorobenzene-d5.I		1,4-Dichloroben. Dichlorobenzene-d4,S		
220000	Benz	oluen	15,S				
210000	64 Y	T	none-c				
200000	e thane -d4,S		2-Hexanone-d5,S				











Data Path : Z:\voasrv\HPCHEM Data File : VV022975.D Acq On : 21 Oct 2021 23: Openator : SY(MD		Instrument : MSVOA_V ClientSampleId : GB7J6							
Operator : SY/MD Sample : M4265-14		00/00							
Sample : M4265-14 Misc : 25.0mL/MSVOA_V/k ALS Vial : 17 Sample Mult		Manual IntegrationsAPPROVED							
Quant Time: Oct 22 04:59:46 Quant Method : Z:\voasrv\HPC Quant Title : TRACE VOA SFA	HEM1\MSVOA_V\Meth M1.0	Reviewed By :John Carlone 10/25/2021 Supervised By :Mahesh Dadoda 10/25/2021							
QLast Update : Fri Oct 22 04:55:17 2021 Response via : Initial Calibration									
Compound	R.T. QIon	Response Conc Units Dev(Min)						
Internal Standards									
1) 1,4-Difluorobenzene	5.619 114	134073 5.000 ug/L	0.00						
28) Chlorobenzene-d5	8.854 117	134602 5.000 ug/L	0.00						
58) 1,4-Dichlorobenzene-d4	11.252 152	59655 5.000 ug/L	0.00						
System Monitoring Compounds 4) Vinyl Chloride-d3	1 204 65	42224							
Spiked Amount 5.000	1.304 65	43224 4.686 ug/L	0.00						
7) Chloroethane-d5	Range 40 - 130 1.568 69	Recovery = 93.800%							
Spiked Amount 5.000	Range 65 - 130	33416 4.031 ug/L Recovery = 80.600%	0.00						
11) 1,1-Dichloroethene-d2	2.105 63	Recovery = 80.600% 55420 3.063 ug/L	0.00						
Spiked Amount 5.000	Range 60 - 125	Recovery = 61.200%							
20) 2-Butanone-d5	3.905 46	75057 31.032 ug/L	0.02						
Spiked Amount 50.000	Range 40 - 130	Recovery = 62.060%	0.02						
24) Chloroform-d	4.349 84	96382 5.070 ug/L	0.00						
Spiked Amount 5.000	Range 70 - 125	Recovery = 101.400%							
26) 1,2-Dichloroethane-d4	5.037 65	43738 4.741 ug/L	0.00						
Spiked Amount 5.000	Range 70 - 130	Recovery = 94.800%							
32) Benzene-d6	5.050 84	187738 4.685 ug/L	0.00						
Spiked Amount 5.000	Range 70 - 125	Recovery = 93.800%							
36) 1,2-Dichloropropane-d6	6.069 67	58885 4.962 ug/L	0.00						
Spiked Amount 5.000	Range 60 - 140	Recovery = 99.200%							
41) Toluene-d8	7.317 98	156447 4.516 ug/L	0.00						
Spiked Amount 5.000	Range 70 - 130	Recovery = 90.400%							
<pre>43) trans-1,3-Dichloroprop. Spiked Amount 5.000</pre>	7.625 79 Range 55 - 130	17426 4.427 ug/L	0.00						
46) 2-Hexanone-d5	8.091 63	Recovery = 88.600% 77060 48.576 ug/L	0.00						
Spiked Amount 50.000	Range 45 - 130	77060 48.576 ug/L Recovery = 97.160%	0.00						
56) 1,1,2,2-Tetrachloroeth.		40107 4.989 ug/L	0.00						
Spiked Amount 5.000	Range 65 - 120	Recovery = 99.800%							
66) 1,2-Dichlorobenzene-d4	11.625 152		0.00						
Spiked Amount 5.000	Range 80 - 120	Recovery = 109.000%							
Target Compounds		Qval	ue a Q						
13) Acetone	2.191 43	12139m 9.000 ug/L	7 MO						
42) Toluene	7.397 91	9951 0.270 ug/L	94 10/20/21						
53) m,p-xylene	9.146 106	1630 0.111 ug/L	68						

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