Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102221\

Data File : VV022988.D

Acq On : 22 Oct 2021 15:40

Operator : SY/MD Sample : M4265-18

Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 31 Sample Multiplier: 1

Quant Time: Oct 23 01:29:20 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102221WMA.M

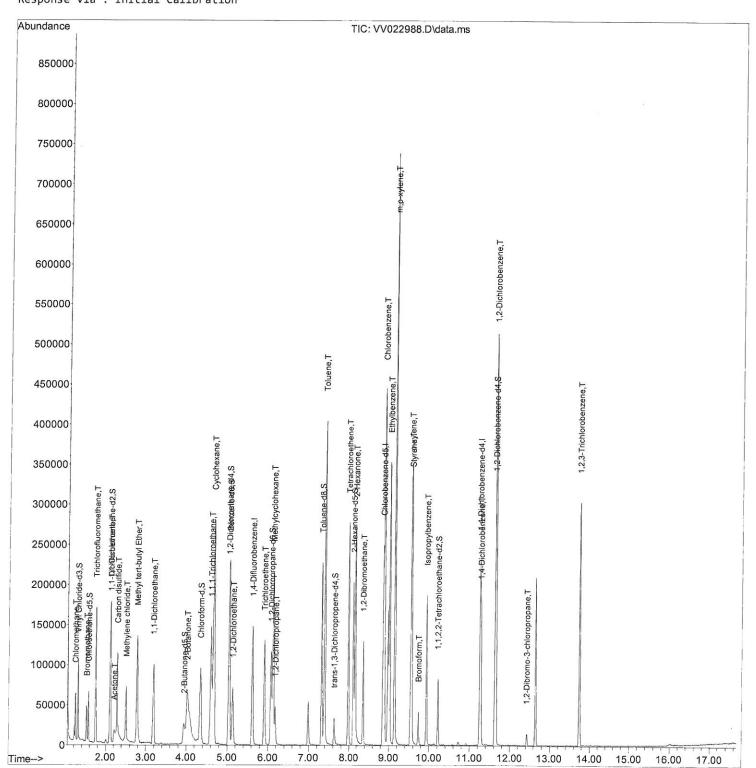
Quant Title : TRACE VOA SFAM1.0

QLast Update : Sat Oct 23 01:14:46 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :John Carlone 10/25/2021 Supervised By :Mahesh Dadoda 10/25/2021



Quantitation Report (Qedit)

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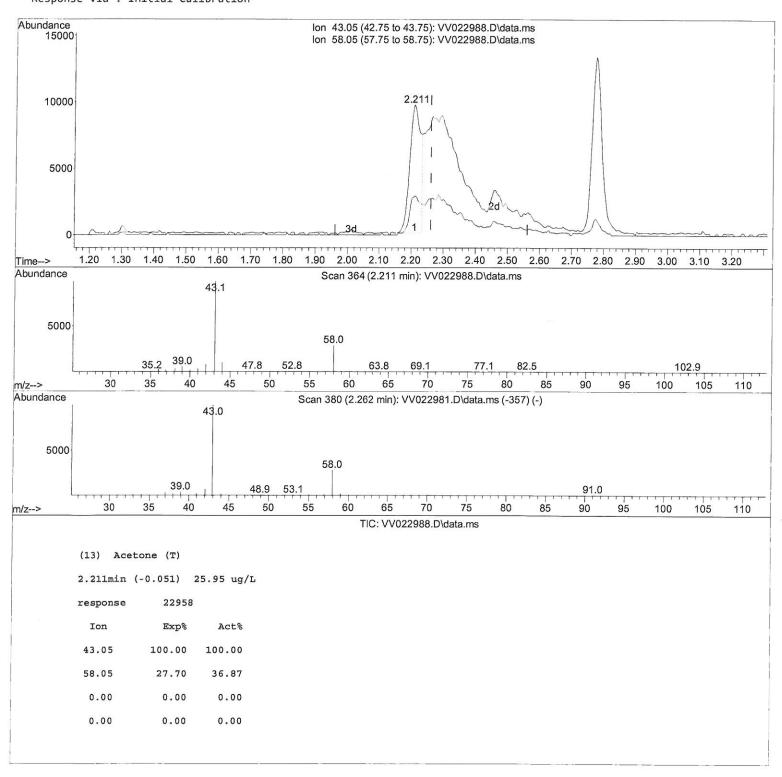
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Quant Title : TRACE VOA SFAM1.0

QLast Update : Sat Oct 23 01:14:46 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleId : GB7K0

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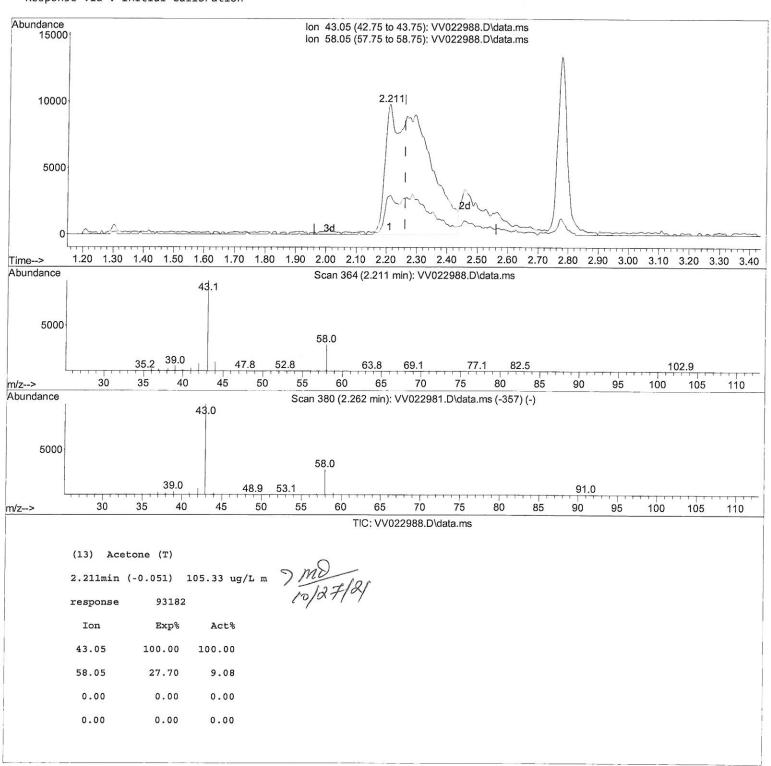
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Compound	R.T.	QIon	Response	Conc Units Dev(Min)
Internal Standards				
1) 1,4-Difluorobenzene	5.616	114	128639	5.000 ug/L 0.00
28) Chlorobenzene-d5	8.854		129680	5.000 ug/L 0.00
58) 1,4-Dichlorobenzene-d4	11.249		58439	5.000 ug/L 0.00
System Monitoring Compounds	4 204		F244F	4.755
4) Vinyl Chloride-d3	1.301	65	53445	4.755 ug/L 0.00
Spiked Amount 5.000	Range 40		Recove	*
7) Chloroethane-d5	1.561		36199	5.206 ug/L -0.02
Spiked Amount 5.000	Range 65		Recove	The state of the s
11) 1,1-Dichloroethene-d2	2.102		78665	4.851 ug/L -0.02
Spiked Amount 5.000	Range 60		Recove	
20) 2-Butanone-d5	3.941	46	87953	48.777 ug/L -0.02
Spiked Amount 50.000	Range 40		Recove	The state of the s
24) Chloroform-d	4.346	84		5.218 ug/L 0.00
Spiked Amount 5.000	Range 70			ery = 104.400%
26) 1,2-Dichloroethane-d4	5.034		44478	5.163 ug/L 0.00
Spiked Amount 5.000	Range 70			ery = 103.200%
32) Benzene-d6	5.044		194030	5.125 ug/L 0.00
Spiked Amount 5.000	Range 70			ery = 102.400%
36) 1,2-Dichloropropane-d6	6.072	67	59935	3.
Spiked Amount 5.000	Range 60			ery = 102.800%
41) Toluene-d8	7.317			4.466 ug/L -0.01
Spiked Amount 5.000 43) trans-1,3-Dichloroprop.	Range 70		Recove	
Spiked Amount 5.000		79 130		4.925 ug/L -0.01
46) 2-Hexanone-d5	Range 55 8.101	63	Recove	
Spiked Amount 50.000	Range 45			
56) 1,1,2,2-Tetrachloroeth.			Recove 39054	
Spiked Amount 5.000	Range 65			4.853 ug/L 0.00 ery = 97.000%
66) 1,2-Dichlorobenzene-d4				5.462 ug/L 0.00
Spiked Amount 5.000	Range 80			ry = 109.200%
Sparked / miledine Steel	nunge oo	120		103.200%
Target Compounds				Qvalue
Chloromethane	1.240	50	38625	4.379 ug/L 95
6) Bromomethane	1.516	94	17461	3.793 ug/L 99
Trichlorofluoromethane	1.748		97062	7.894 ug/L 99
12) 1,1-Dichloroethene	2.111	96	24403	3.705 ug/L # 74
13) Acetone	2.211	43	93182m	
14) Carbon disulfide	2.285	76	107884	6.015 ug/L 98 /0/07/0
16) Methylene chloride	2.503	84	26345	3.642 ug/L 91
17) Methyl tert-butyl Ether	2.777	73	139311	8.860 ug/L 98
19) 1,1-Dichloroethane	3.182	63	94518	7.316 ug/L 99
21) 2-Butanone	4.021	43	209351	120.878 ug/L 82
27) 1,2-Dichloroethane	5.134	62	63230	7.024 ug/L 97
29) 1,1,1-Trichloroethane	4.600	97	118851	8.111 ug/L 99
30) Cyclohexane	4.661	56	142436	11.132 ug/L 98
34) Trichloroethene	5.908	95	45139	5.149 ug/L 98
35) Methylcyclohexane	6.124	83	81869	6.522 ug/L 97
37) 1,2-Dichloropropane	6.175	63	18446	2.097 ug/L 98
42) Toluene	7.387	91	295520	8.356 ug/L 99
47) Tetrachloroethene	7.973	164	62129	8.388 ug/L 99

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48) 2-Hexanone 8.153 43 177753 57.826 ug/L 50) 1,2-Dibromoethane 8.355 107 83330 15.041 ug/L 51) Chlorobenzene 8.883 112 242995 10.460 ug/L 52) Ethylbenzene 9.011 91 235727 6.715 ug/L	(Min)
53) m,p-xylene 9.140 106 193911 13.776 ug/L 54) o-xylene 9.545 106 70306 5.305 ug/L 55) Styrene 9.561 104 123586 5.401 ug/L 59) Bromoform 9.735 173 19369 5.688 ug/L 60) Isopropylbenzene 9.931 105 112940 3.821 ug/L 65) 1,4-Dichlorobenzene 11.275 146 66324 4.198 ug/L 67) 1,2-Dichlorobenzene 11.645 146 186396 12.874 ug/L	(Min) 99 96 99 98 99 99 99 100 98
68) 1,2-Dibromo-3-chloropr 12.429 75 3517 4.333 ug/L # 72) 1,2,3-Trichlorobenzene 13.744 180 92190 11.205 ug/L	84 99

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