Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102621\

Data File: VV023048.D

Acq On : 26 Oct 2021 19:14

Operator : SY/MD

Sample : M4277-16DL 10X Misc : 25.0mL/MSVOA_V/WATER ALS Vial : 25 Sample Multiplier: 1

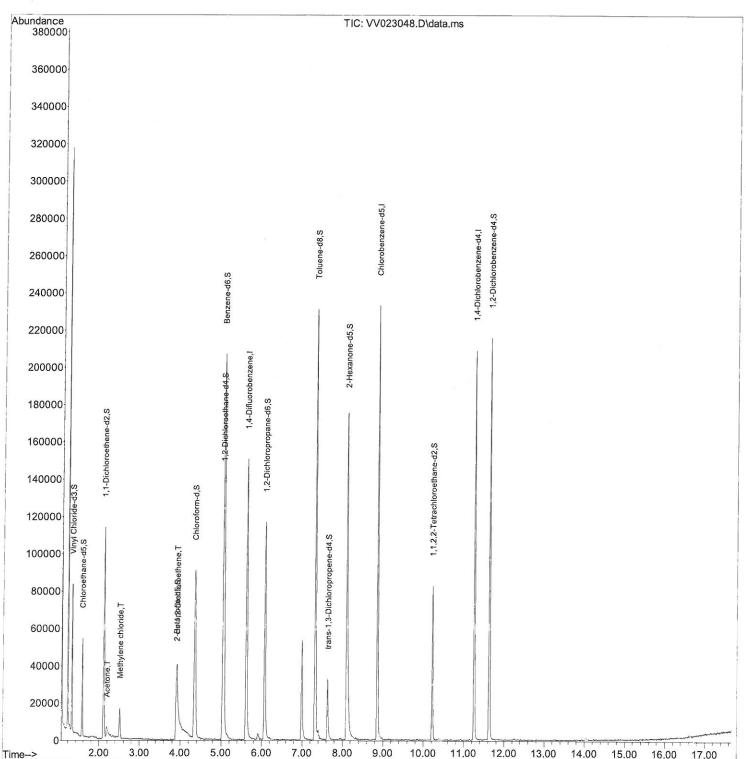
Quant Time: Oct 27 01:35:38 2021

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

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Oct 27 01:29:33 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102621\

Data File : VV023048.D

Acq On : 26 Oct 2021 19:14

Operator : SY/MD

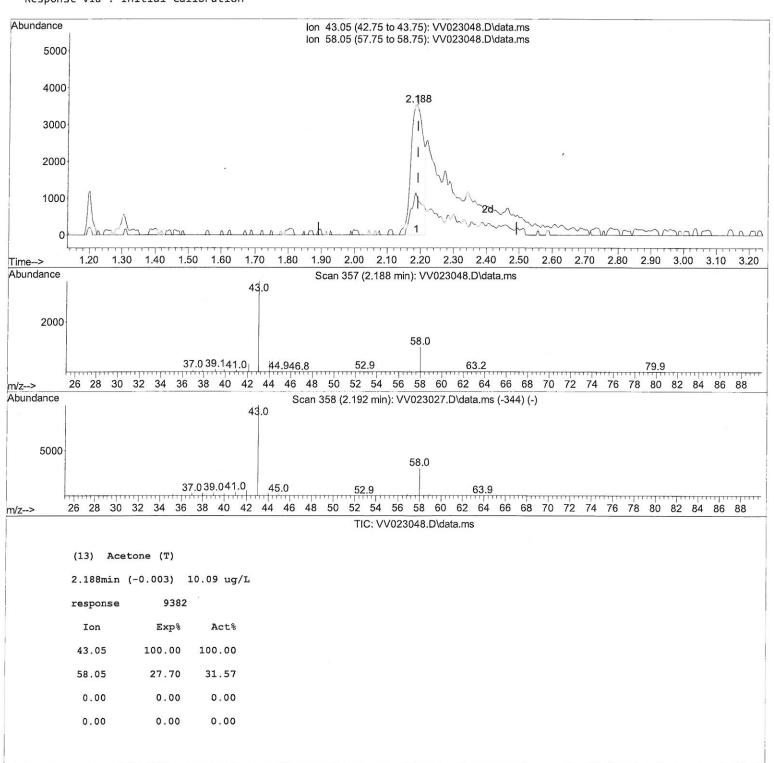
Sample : M4277-16DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Quant Time: Oct 27 01:35:38 2021

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

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Oct 27 01:29:33 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleId : BFGE1DL

Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102621\

Data File: VV023048.D

Acq On : 26 Oct 2021 19:14

Operator : SY/MD

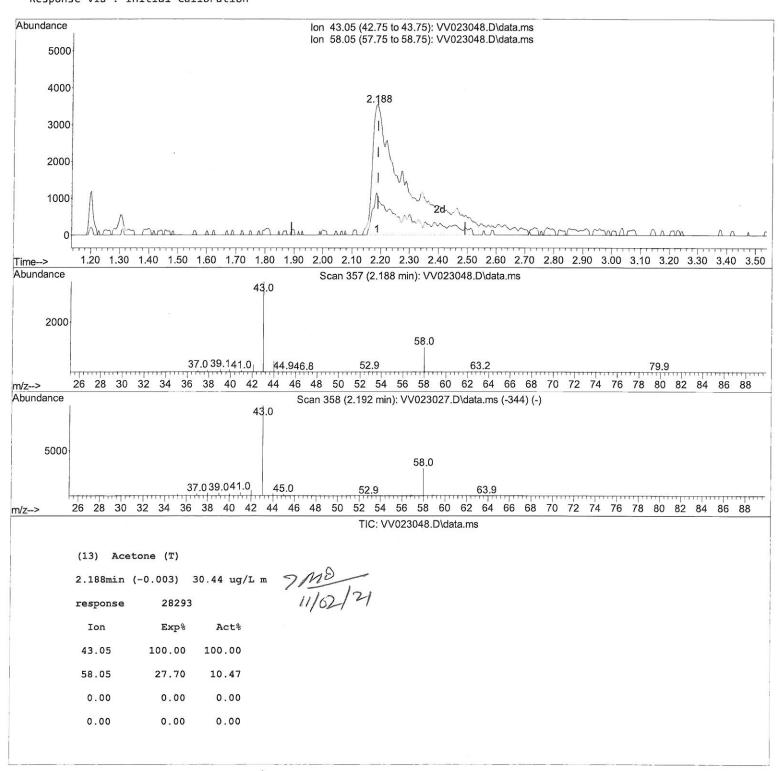
Sample : M4277-16DL 10X Misc : 25.0mL/MSVOA_V/WATER ALS Vial : 25 Sample Multiplier: 1

Quant Time: Oct 27 01:35:38 2021

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

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Oct 27 01:29:33 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleId : BFGE1DL

Manual Integrations APPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102621\

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Manual IntegrationsAPPROVED

Compound	R.T. QIon	Response Conc Units Dev(Min)
Internal Standards		
 1,4-Difluorobenzene 		135154 5.000 ug/L 0.00
28) Chlorobenzene-d5	8.854 117	133072 5.000 ug/L 0.00
58) 1,4-Dichlorobenzene-d4	11.249 152	57531 5.000 ug/L 0.00
System Monitoring Compounds		
4) Vinyl Chloride-d3		49697 4.208 ug/L 0.00
Spiked Amount 5.000		Recovery = 84.200%
7) Chloroethane-d5	1.568 69	32730 4.480 ug/L 0.00
Spiked Amount 5.000	Range 65 - 130	Recovery = 89.600% 58515 3.435 ug/L 0.00 Recovery = 68.600% 81411 42.972 ug/L 0.00
<pre>11) 1,1-Dichloroethene-d2</pre>	2.108 63	58515 3.435 ug/L 0.00
Spiked Amount 5.000	Range 60 - 125	Recovery = 68.600%
20) 2-Butanone-d5	3.902 46	81411 42.972 ug/L 0.00
Spiked Amount 50.000	Range 40 - 130	Recovery = 85.940%
24) Chloroform-d	4.349 84	96234 5.012 ug/L 0.00
Spiked Amount 5.000	Range 70 - 125	Recovery = 100.200%
26) 1,2-Dichloroethane-d4	5.034 65	44703 4.939 ug/L 0.00
Spiked Amount 5.000	Range 70 - 130	Recovery = 98.800%
32) Benzene-d6	5.050 84	189961 4.890 ug/L 0.00
Spiked Amount 5.000	Range 70 - 125	Recovery = 97.800%
36) 1,2-Dichloropropane-d6	6.072 67	59393 4.966 ug/L 0.00
Spiked Amount 5.000	Range 60 - 140	Recovery = 99.400%
41) Toluene-d8	7.317 98	158631 4.546 ug/L 0.00
Spiked Amount 5.000	Range 70 - 130	Recovery = 91.000%
Spiked Amount 5.000 43) trans-1,3-Dichloroprop.	7.629 79	19766 4.717 ug/L 0.00
Spiked Amount 5.000	Range 55 - 130	Recovery = 94.400%
46) 2-Hexanone-d5		73295 47.244 ug/L 0.00
Spiked Amount 50.000		
56) 1,1,2,2-Tetrachloroeth.		
Spiked Amount 5.000		
66) 1,2-Dichlorobenzene-d4	11.625 152	57747 5.626 ug/L 0.00
66) 1,2-Dichlorobenzene-d4 Spiked Amount 5.000	Range 80 - 120	Recovery = 112.600%
Target Compounds		Qvalue
13) Acetone	2.188 43	28293m 30.440 ug/L 7MC
16) Methylene chloride		6607 0.869 ug/L 97 11/02/0
22) cis-1,2-Dichloroethene		

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