Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111721\

Data File : VV023578.D

Acq On : 17 Nov 2021 18:35

Operator : SY/MD Sample : M4627-01

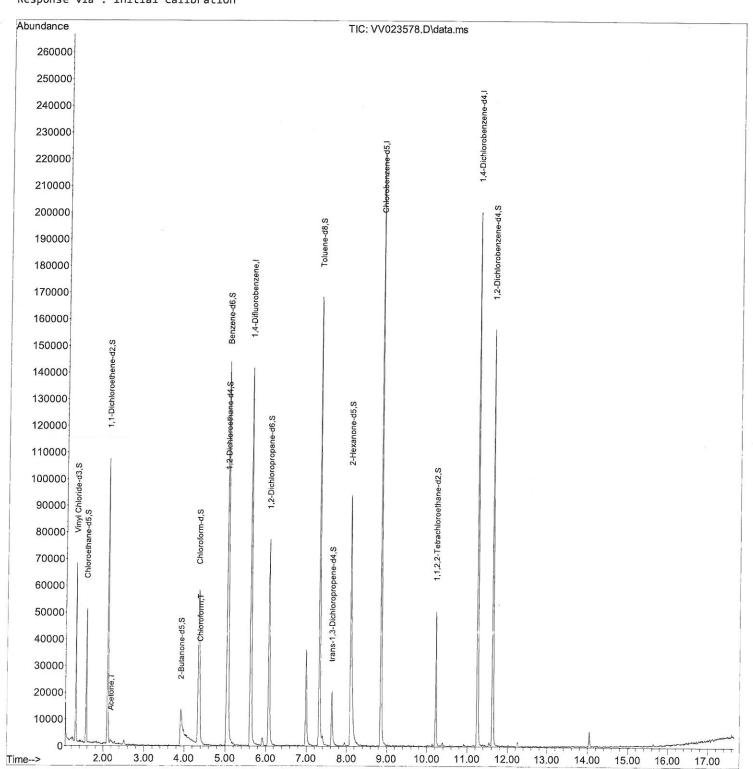
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Quant Time: Nov 18 00:23:40 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Thu Nov 18 00:20:29 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleId : H4630

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111721\

Data File: VV023578.D

Acq On : 17 Nov 2021 18:35

Operator : SY/MD Sample : M4627-01

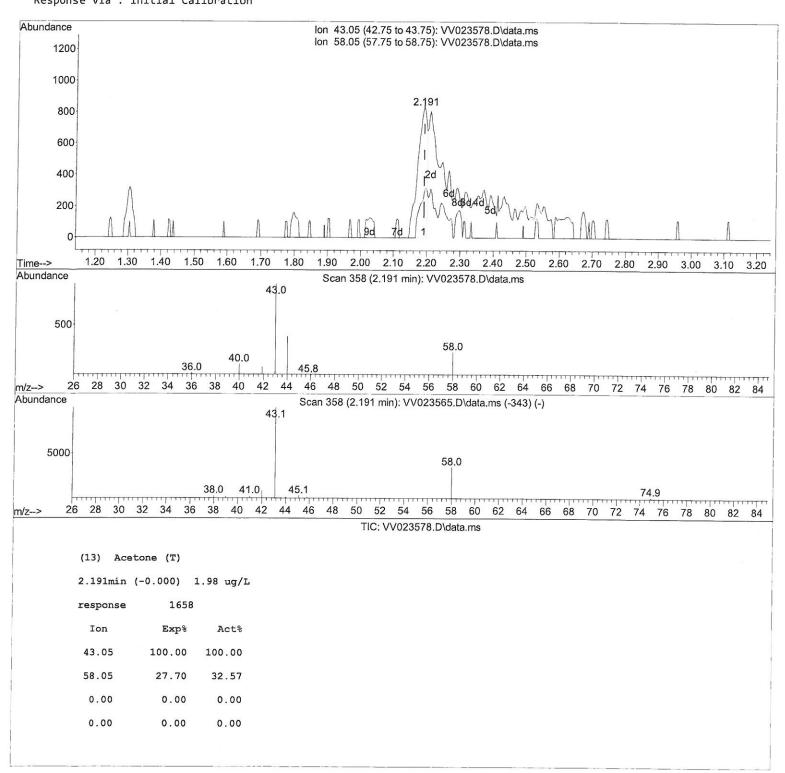
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Quant Time: Nov 18 00:23:40 2021

 $\label{eq:Quant_Method} Quant \ \mbox{Method} : \ Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M$

Quant Title : TRACE VOA SFAM1.0 QLast Update : Thu Nov 18 00:20:29 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleId : H4630

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111721\

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Acq On : 17 Nov 2021 18:35

Operator : SY/MD Sample : M4627-01

Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Quant Time: Nov 18 00:23:40 2021

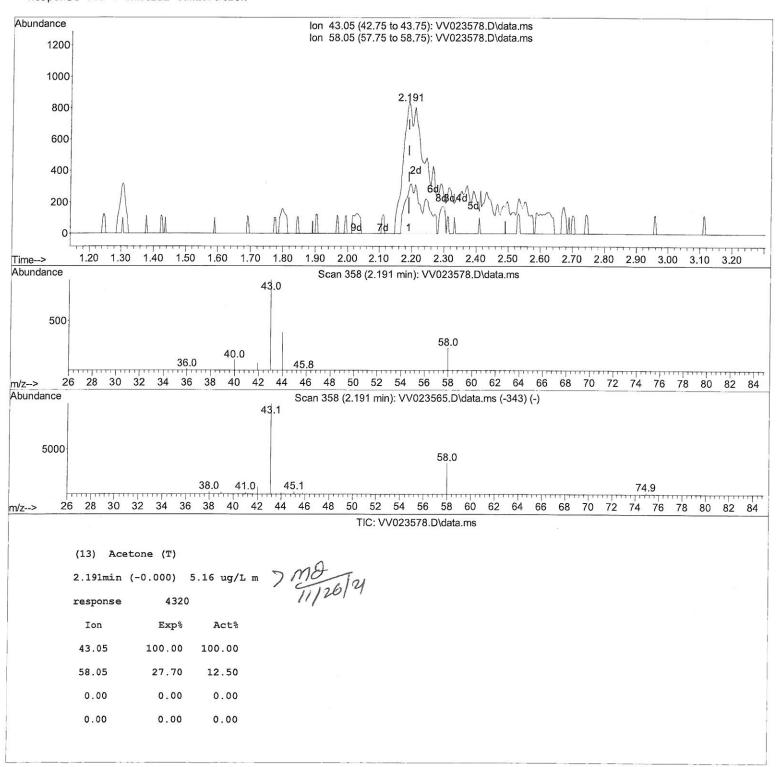
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Quant Title : TRACE VOA SFAM1.0

QLast Update : Thu Nov 18 00:20:29 2021 Response via : Initial Calibration

Instrument : MSVOA_V ClientSampleld : H4630

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111721\

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ALS Vial : 15 Sample Multiplier: 1

Quant Time: Nov 18 00:23:40 2021

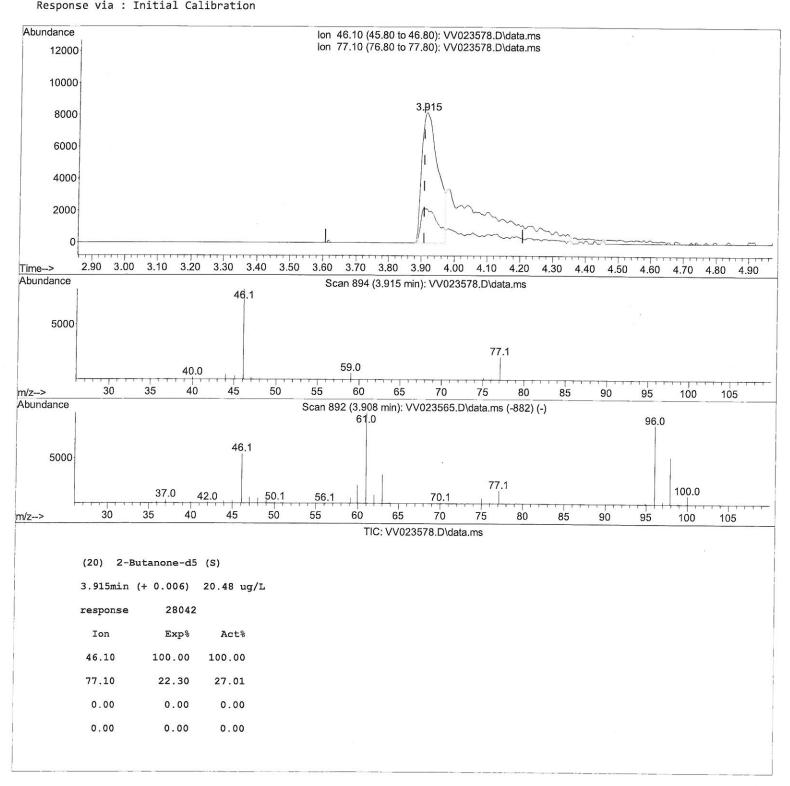
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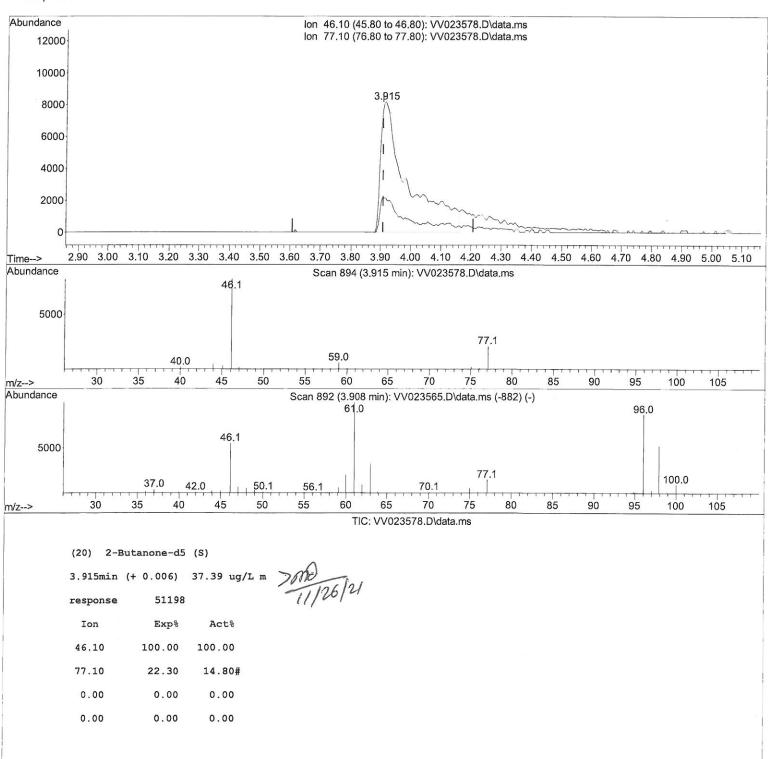
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Quant Time: Nov 18 00:23:40 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

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Manual IntegrationsAPPROVED

Compound	R.T. OIon	Response Conc Un	its Dev(Min)
Internal Standards			
 1,4-Difluorobenzene 	5.619 114	126860 5.000	ug/L 0.00
28) Chlorobenzene-d5	8.853 117	122625 5.000	ug/L 0.00
58) 1,4-Dichlorobenzene-d4	11.249 152	55001 5.000	ug/L 0.00
			0,000 € 0,000 (0.000 (
System Monitoring Compounds	i		
Vinyl Chloride-d3	1.304 65	38830 4.886	ug/L 0.00
Spiked Amount 5.000	Range 40 - 130	Recovery =	97.800%
7) Chloroethane-d5	1.564 69	28995 4.476	ug/L 0.00
Spiked Amount 5.000	Range 65 - 130	Recovery =	89.600%
11) 1,1-Dichloroethene-d2	2.105 63	53679 3.608	ug/L 0.00
Spiked Amount 5.000	Range 60 - 125	Recovery =	72.200%
20) 2-Butanone-d5	3.915 46	51198m 37.393	ug/L 0.00 7/10/12/
Spiked Amount 50.000	Range 40 - 130	Recovery =	74.780%
24) Chloroform-d	4.349 84	56004 3.307	ug/L 0.00
Spiked Amount 5.000	Range 70 - 125	Recovery =	66.200%#
26) 1,2-Dichloroethane-d4	5.034 65	30070 3.948	ug/L 0.00
Spiked Amount 5.000	Range 70 - 130	Recovery =	79.000%
32) Benzene-d6	5.053 84	130513 4.148	ug/L 0.00
Spiked Amount 5.000	Range 70 - 125	Recovery =	83.000%
36) 1,2-Dichloropropane-d6	6.072 67	37189 4.015	ug/L 0.00
Spiked Amount 5.000	Range 60 - 140	Recovery =	80.400%
41) Toluene-d8	7.317 98	111530 3.783	ug/L 0.00
Spiked Amount 5.000	Range 70 - 130	Recovery =	75.600%
43) trans-1,3-Dichloroprop	7.625 79	13166 3.749	ug/L 0.00
Spiked Amount 5.000	Range 55 - 130	Recovery =	75.000%
46) 2-Hexanone-d5	8.095 63	41245 31.920	ug/L 0.00
Spiked Amount 50.000	Range 45 - 130	Recovery =	63.840%
56) 1,1,2,2-Tetrachloroeth		23115 3.470	ug/L 0.00
Spiked Amount 5.000	Range 65 - 120		69.400%
66) 1,2-Dichlorobenzene-d4			ug/L 0.00
Spiked Amount 5.000	Range 80 - 120	Recovery =	89.400%
Tanget Compounds			
Target Compounds	2 101 42	4220 5 454	Qvalue 7 M
13) Acetone 25) Chloroform	2.191 43	4320m 5.164	
25) CHIOLOLOLUM	4.375 83	9867 0.590	ug/L 100 ///20/5/

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