Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111921\

Data File : VV023643.D

Acq On : 19 Nov 2021 15:04

Operator : SY/MD Sample : M4706-18

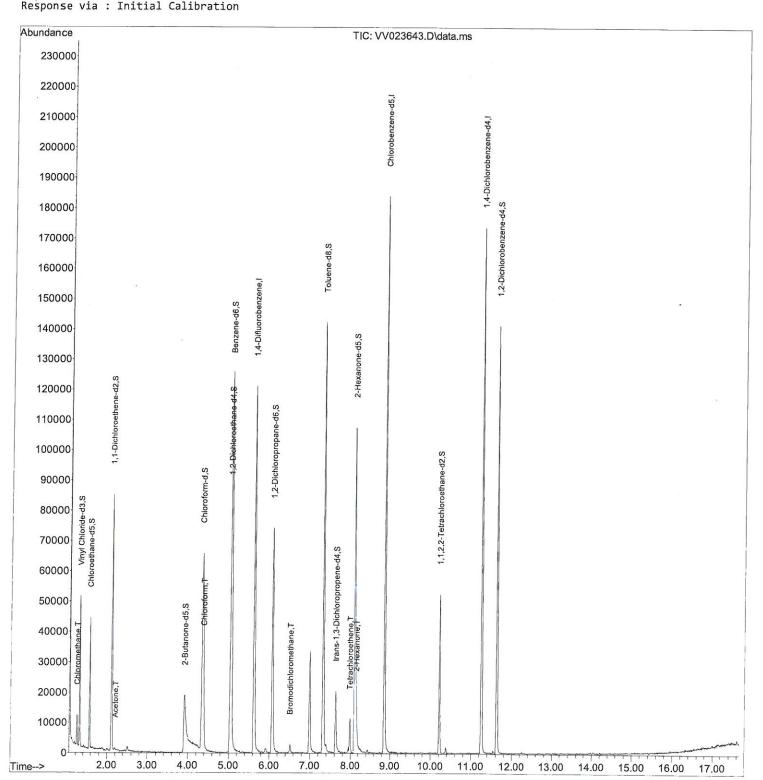
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 22 01:48:16 2021

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

Quant Title : TRACE VOA SFAM1.0 QLast Update : Mon Nov 22 01:44:25 2021 Instrument : MSVOA_V ClientSampleId : B0AB1

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111921\

Data File : VV023643.D

Acq On : 19 Nov 2021 15:04

Operator : SY/MD Sample : M4706-18

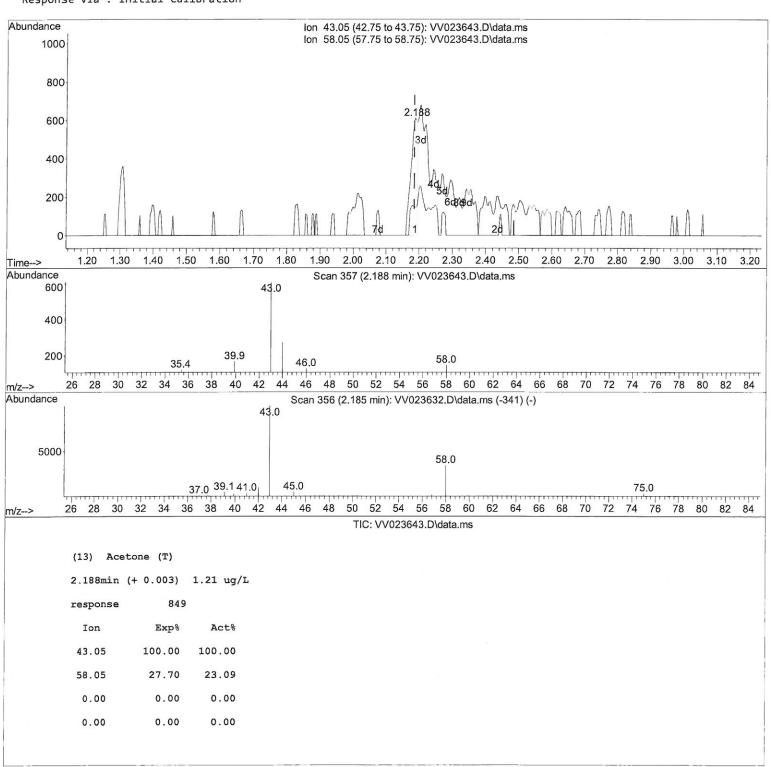
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ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 22 01:48:16 2021

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

Quant Title : TRACE VOA SFAM1.0 QLast Update : Mon Nov 22 01:44:25 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleId : B0AB1

Manual IntegrationsAPPROVED



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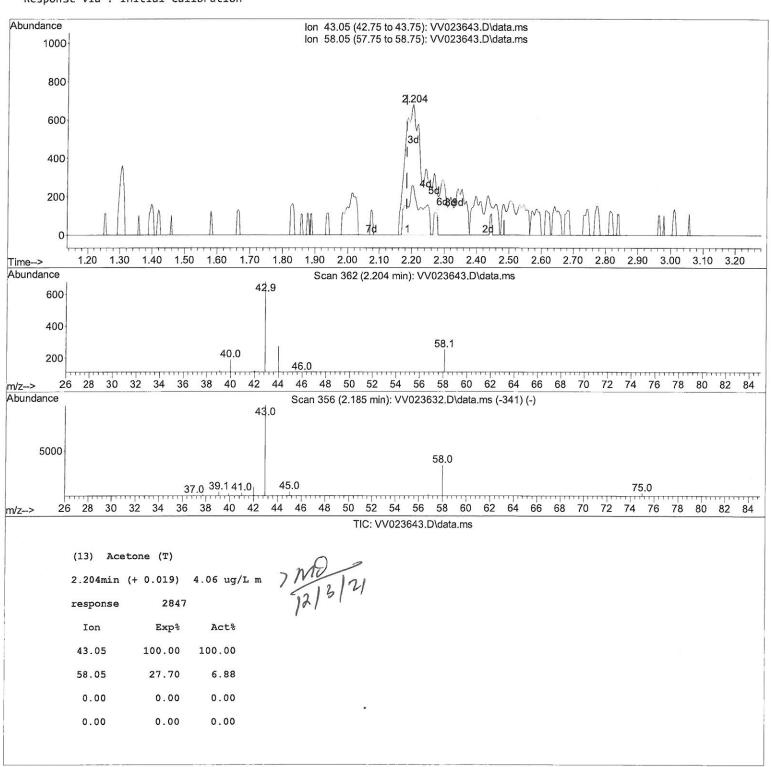
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 22 01:48:16 2021

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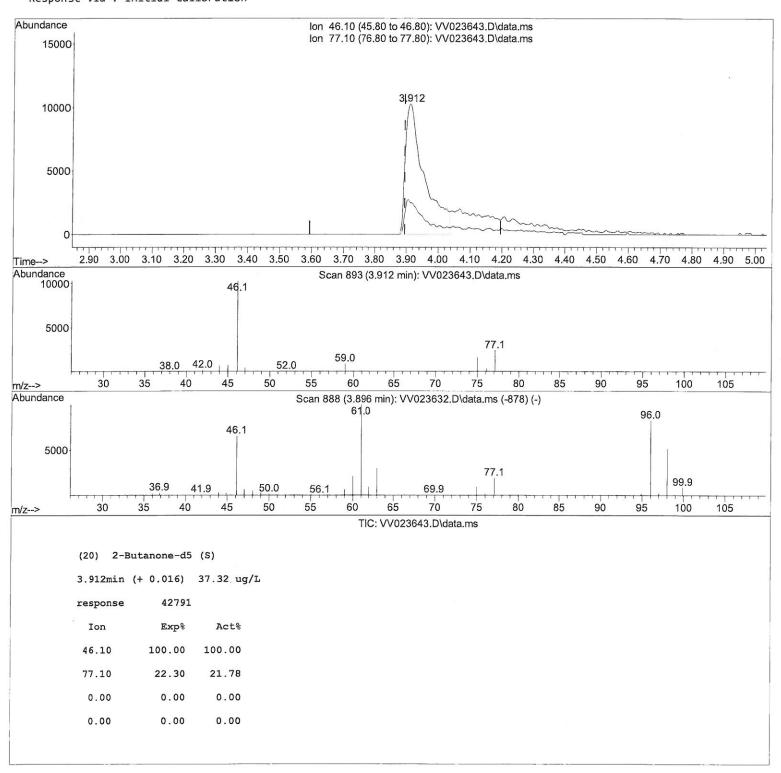
Misc : 25.0mL/MSVOA_V/WATER
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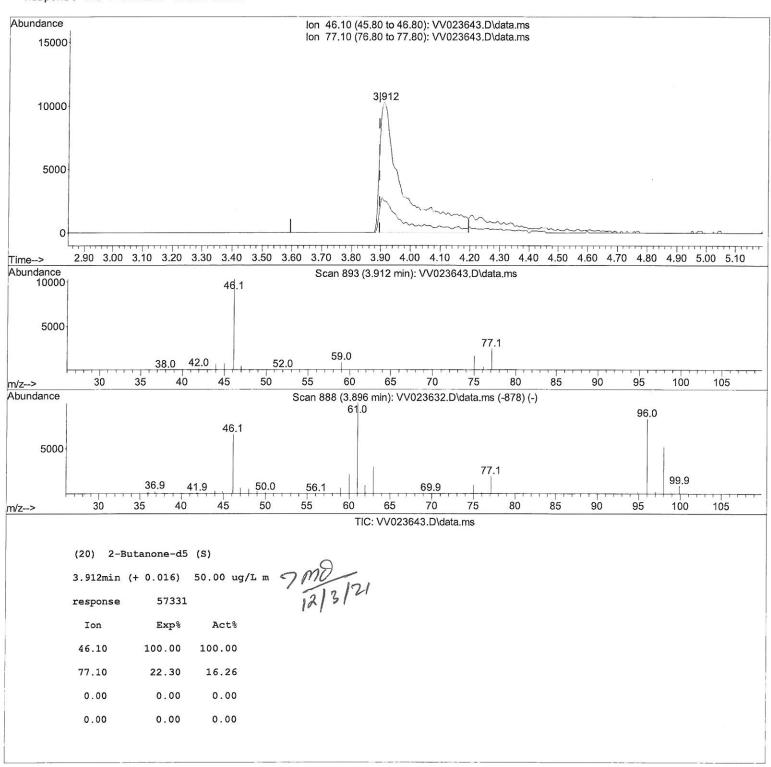
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0

QLast Update : Mon Nov 22 01:44:25 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleId : B0AB1

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111921\

Data File : VV023643.D

Acq On : 19 Nov 2021 15:04

Operator : SY/MD Sample : M4706-18

: 25.0mL/MSVOA_V/WATER Misc ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 22 01:48:16 2021

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

Quant Title : TRACE VOA SFAM1.0 QLast Update : Mon Nov 22 01:44:25 2021 Response via: Initial Calibration

Instrument : MSVOA_V ClientSampleId : B0AB1

Manual IntegrationsAPPROVED

Response via : initial Calib	racion		
Compound	R.T. QIon	Response Conc Un	its Dev(Min)
Internal Standards			
1) 1,4-Difluorobenzene	5.619 114	106239 5.000	ug/L 0.00
28) Chlorobenzene-d5	8.853 117		ug/L 0.00
58) 1,4-Dichlorobenzene-d4			ug/L 0.00
System Monitoring Compounds			
4) Vinyl Chloride-d3	1.304 65	28513 4.284	ug/L 0.00
Spiked Amount 5.000	Range 40 - 130	Recovery =	85.600%
7) Chloroethane-d5	1.568 69	25251 4.655	
Spiked Amount 5.000	Range 65 - 130	Recovery =	
11) 1,1-Dichloroethene-d2	2.108 63		
Spiked Amount 5.000	Range 60 - 125		65.800%
20) 2-Butanone-d5	3.912 46	170	2177 -
Spiked Amount 50.000	Range 40 - 130		100.000%
24) Chloroform-d	4.349 84	63005 4.442	10/0/
Spiked Amount 5.000	Range 70 - 125		88.800%
26) 1,2-Dichloroethane-d4	5.034 65	29860 4.682	
Spiked Amount 5.000	Range 70 - 130		93.600%
32) Benzene-d6	5.053 84	112652 4.132	
Spiked Amount 5.000	Range 70 - 125		82.600%
36) 1,2-Dichloropropane-d6	6.072 67	34491 4.298	
Spiked Amount 5.000	Range 60 - 140		86.000%
41) Toluene-d8	7.317 98	95152 3.724	
Spiked Amount 5.000	Range 70 - 130		74.400%
43) trans-1,3-Dichloroprop.			
Spiked Amount 5.000	Range 55 - 130		81.400%
46) 2-Hexanone-d5	8.091 63	45837 40.939	
Spiked Amount 50.000	Range 45 - 130		81.880%
56) 1,1,2,2-Tetrachloroeth.			
Spiked Amount 5.000	Range 65 - 120		85.400%
66) 1,2-Dichlorobenzene-d4			
Spiked Amount 5.000	Range 80 - 120	Recovery =	92.800%
Target Compounds			Ovalue
3) Chloromethane	1.240 50	6017 0.683	
13) Acetone	2.204 43	2847m 4.064	WIX
25) Chloroform	4.378 83	15084 1.076	, , ,
38) Bromodichloromethane	6.516 83		ug/L # 87 12/5/01
47) Tetrachloroethene	7.976 164	2979 0.435	
48) 2-Hexanone	8.149 43		ug/L # 62

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