Quantitation Report (Qedit)

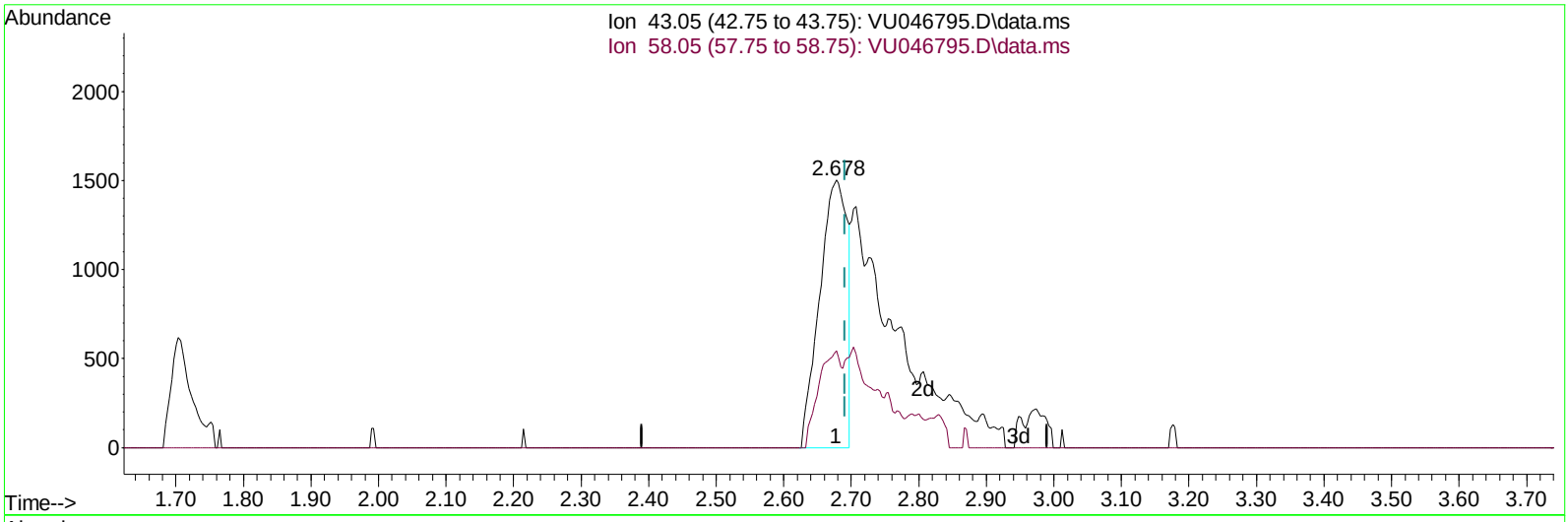
Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU011922\
Data File : VU046795.D
Acq On : 19 Jan 2022 13:21
Operator : SY/MD
Sample : N1179-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

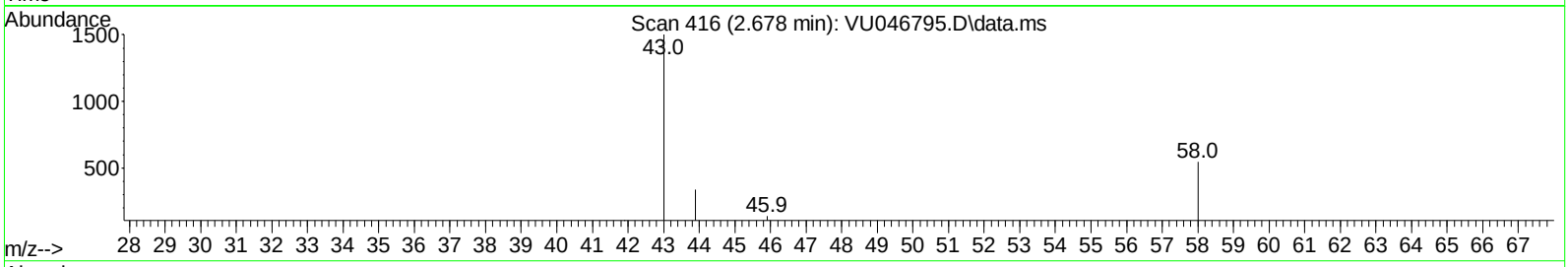
Quant Time: Jan 20 02:01:48 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR011822WMA.M
Quant Title : TRACE VOA SFAM1.0
QLast Update : Thu Jan 20 02:00:32 2022
Response via : Initial Calibration

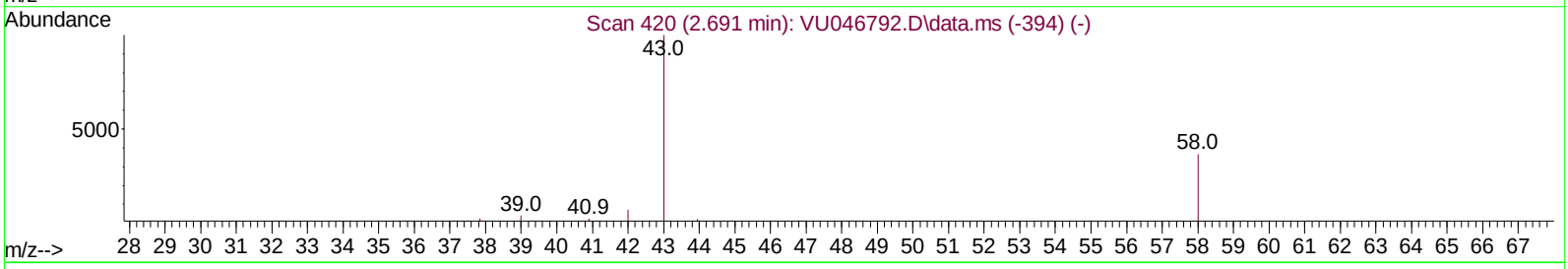
Instrument :
MSVOA_U
ClientSampleId :
C0GG5

Manual IntegrationsAPPROVED

Reviewed By :John Carlone 01/20/2022
Supervised By :Mahesh Dadoda 01/21/2022







TIC: VU046795.D\data.ms

(13) Acetone (T)

2.678min (-0.013) 2.66 ug/L

response 4264

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	35.80	30.28
0.00	0.00	0.00
0.00	0.00	0.00

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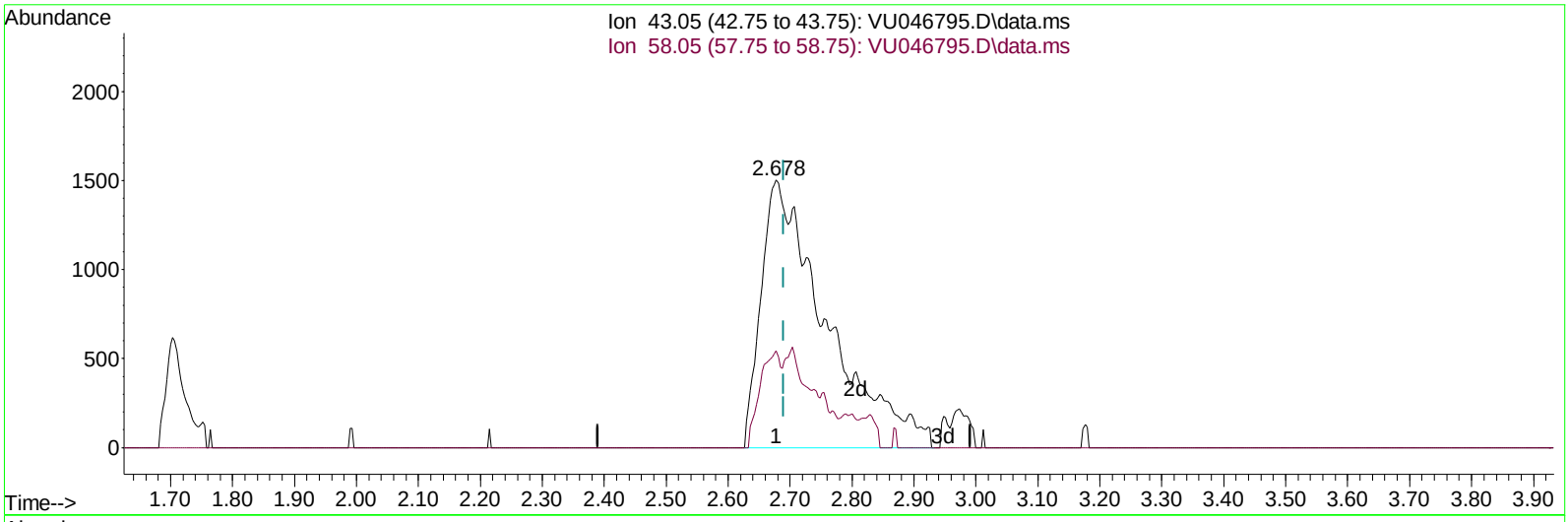

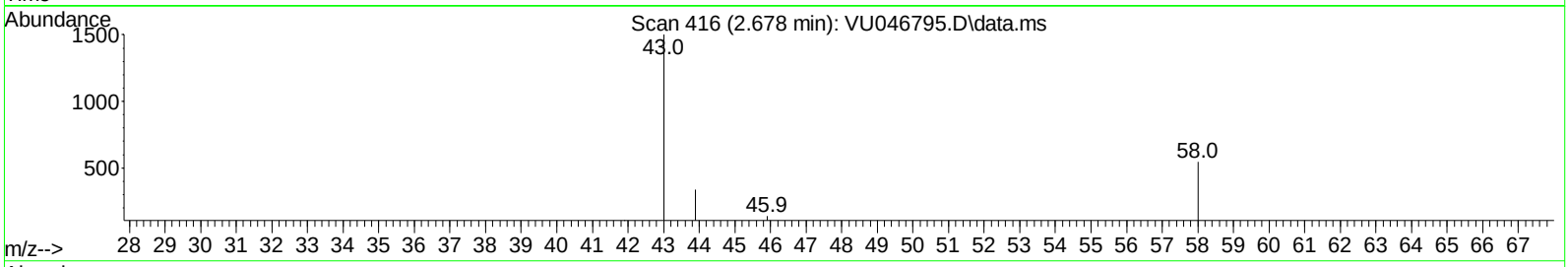

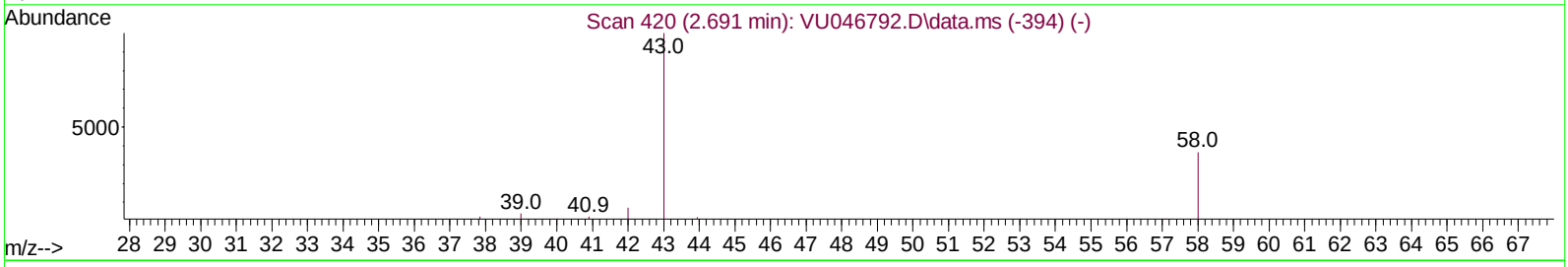

TIC: VU046795.D\data.ms

(13) Acetone (T)

2.678min (-0.013) 6.83 ug/L m

response 10947

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	35.80	11.79
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (QT/LSC Reviewed)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU011922\
Data File : VU046795.D
Acq On : 19 Jan 2022 13:21
Operator : SY/MD
Sample : N1179-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Jan 20 02:01:48 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR011822WMA.M
Quant Title : TRACE VOA SFAM1.0
QLast Update : Thu Jan 20 02:00:32 2022
Response via : Initial Calibration

Instrument :
MSVOA_U
ClientSampleId :
C0GG5

Manual IntegrationsAPPROVED

Reviewed By :John Carlone 01/20/2022
Supervised By :Mahesh Dadoda 01/21/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Difluorobenzene	6.253	114	117368	5.000	ug/L	0.00
28) Chlorobenzene-d5	9.417	117	123132	5.000	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.812	152	61026	5.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.601	65	45958	4.621	ug/L	0.00
Spiked Amount 5.000	Range 40 - 130		Recovery = 92.400%			
7) Chloroethane-d5	1.919	69	38227	4.878	ug/L	0.00
Spiked Amount 5.000	Range 65 - 130		Recovery = 97.600%			
11) 1,1-Dichloroethene-d2	2.572	65	19545	4.941	ug/L	0.00
Spiked Amount 5.000	Range 60 - 125		Recovery = 98.800%			
20) 2-Butanone-d5	4.655	46	163849	55.612	ug/L	0.00
Spiked Amount 50.000	Range 40 - 130		Recovery = 111.220%			
24) Chloroform-d	5.067	84	85640	4.905	ug/L	0.00
Spiked Amount 5.000	Range 70 - 125		Recovery = 98.200%			
26) 1,2-Dichloroethane-d4	5.703	65	48470	5.098	ug/L	0.00
Spiked Amount 5.000	Range 70 - 130		Recovery = 102.000%			
32) Benzene-d6	5.729	84	176411	4.875	ug/L	0.00
Spiked Amount 5.000	Range 70 - 125		Recovery = 97.400%			
36) 1,2-Dichloropropane-d6	6.694	67	54864	5.049	ug/L	0.00
Spiked Amount 5.000	Range 60 - 140		Recovery = 101.000%			
41) Toluene-d8	7.899	98	149854	4.524	ug/L	0.00
Spiked Amount 5.000	Range 70 - 130		Recovery = 90.400%			
43) trans-1,3-Dichloroprop...	8.182	79	21588	4.524	ug/L	0.00
Spiked Amount 5.000	Range 55 - 130		Recovery = 90.400%			
46) 2-Hexanone-d5	8.639	63	144575	51.169	ug/L	0.00
Spiked Amount 50.000	Range 45 - 130		Recovery = 102.340%			
56) 1,1,2,2-Tetrachloroeth...	10.758	84	53997	4.950	ug/L	0.00
Spiked Amount 5.000	Range 65 - 120		Recovery = 99.000%			
66) 1,2-Dichlorobenzene-d4	12.195	152	60080	5.250	ug/L	0.00
Spiked Amount 5.000	Range 80 - 120		Recovery = 105.000%			
Target Compounds						Qvalue
3) Chloromethane	1.520	50	1588	0.175	ug/L	95
13) Acetone	2.678	43	10947m	6.831	ug/L	

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

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