Quantitation Report (Qedit)

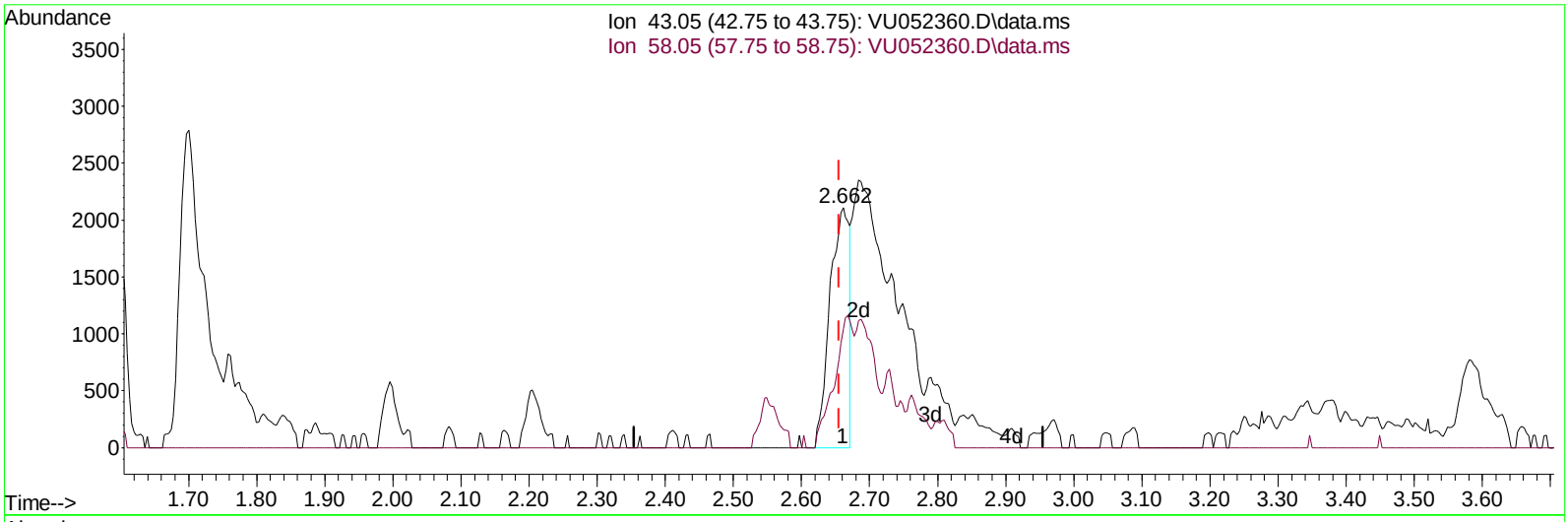
Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU121222\
Data File : VU052360.D
Acq On : 12 Dec 2022 18:08
Operator : JC/MD
Sample : N5961-03
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

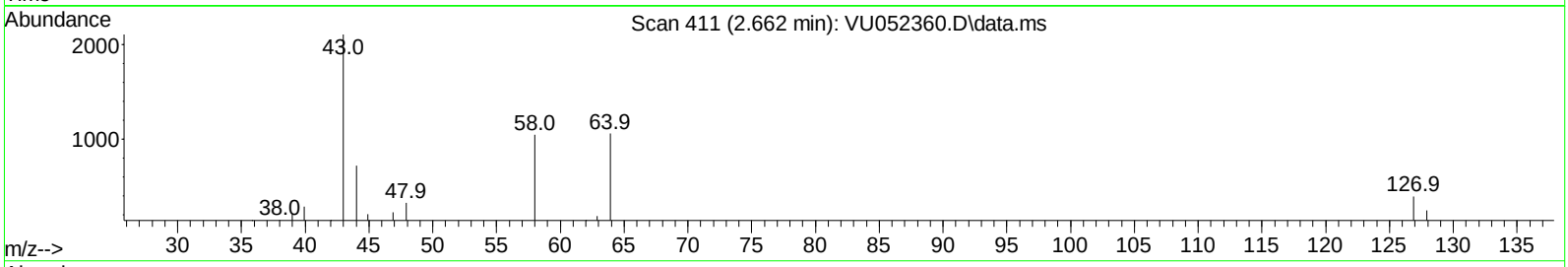
Quant Time: Dec 13 01:06:03 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR111822WMA.M
Quant Title : TRACE VOA SFAM1.0
QLast Update : Tue Dec 13 01:01:45 2022
Response via : Initial Calibration

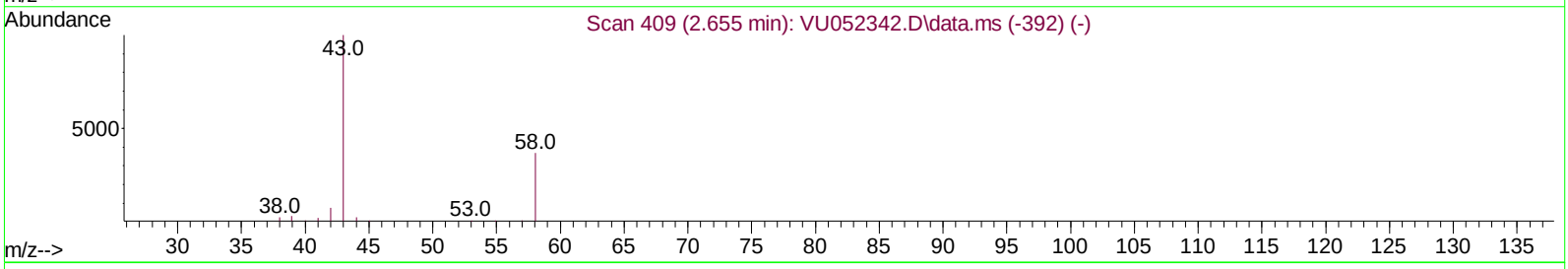
Instrument :
MSVOA_U
ClientSampleId :
F5N34

Manual IntegrationsAPPROVED

Reviewed By :Krupa Patel 12/13/2022
Supervised By :Mahesh Dadoda 12/13/2022







TIC: VU052360.D\data.ms

(13) Acetone (T)

2.662min (+ 0.006) 1.48 ug/L

response 4203

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	38.00	54.65
0.00	0.00	0.00
0.00	0.00	0.00

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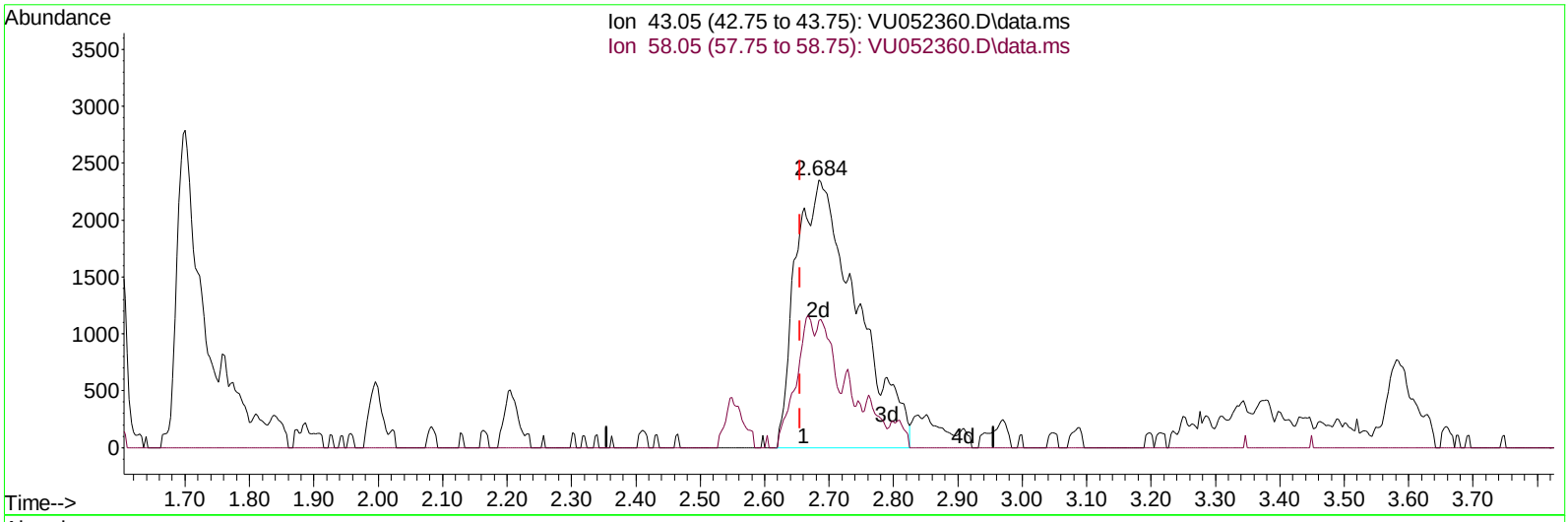

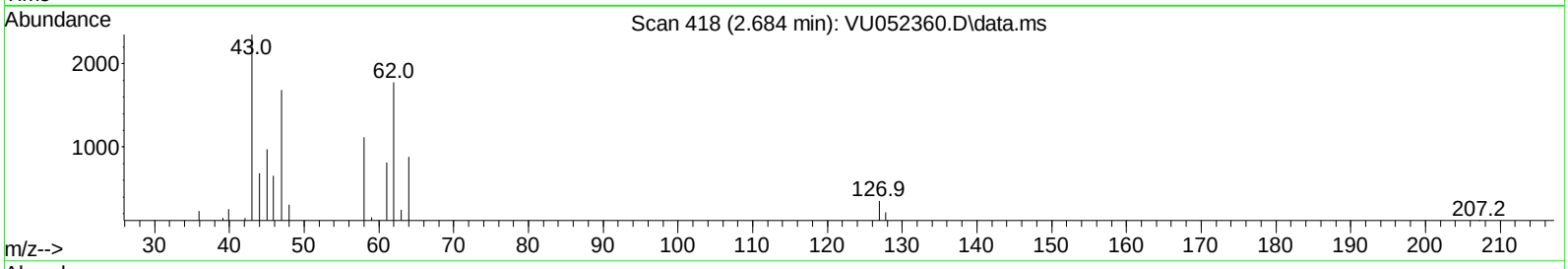

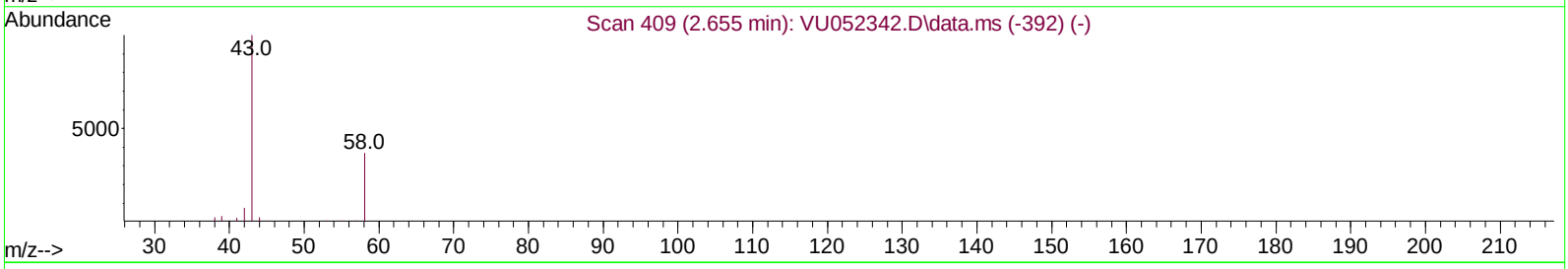

TIC: VU052360.D\data.ms

(13) Acetone (T)

2.684min (+ 0.029) 5.40 ug/L m

response 15361

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	38.00	14.95
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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Manual IntegrationsAPPROVED

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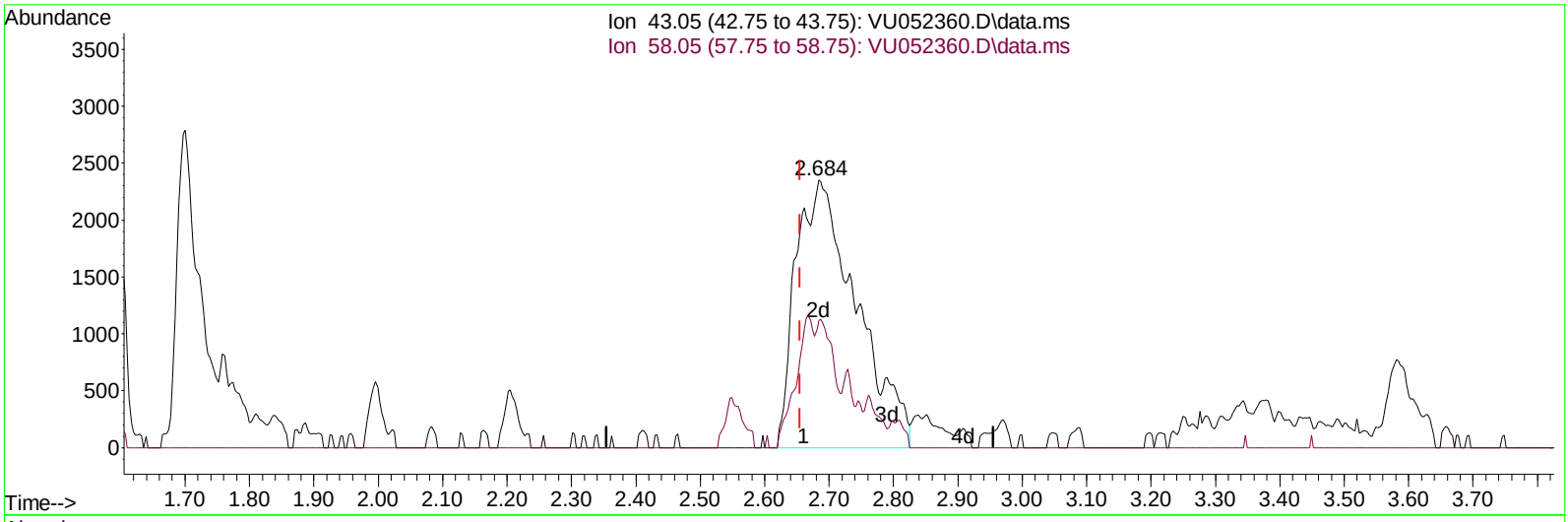

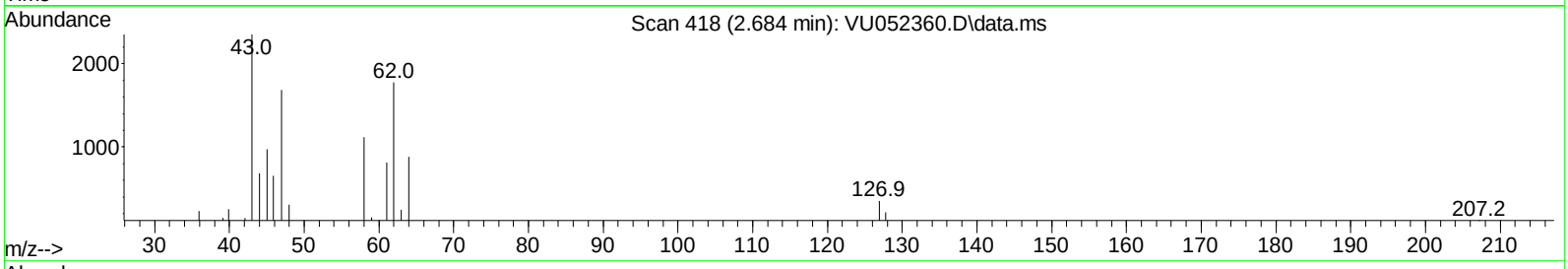

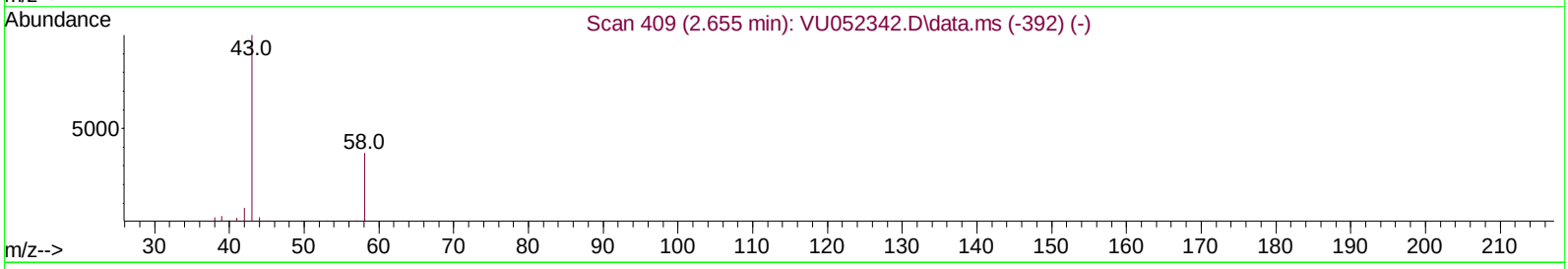

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(13) Acetone (T)

2.684min (+ 0.029) 5.40 ug/L m

response 15361

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	38.00	14.95
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (QT/LSC Reviewed)

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QLast Update : Tue Dec 13 01:01:45 2022
Response via : Initial Calibration

Instrument :
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ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Krupa Patel 12/13/2022
Supervised By :Mahesh Dadoda 12/13/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Difluorobenzene	6.247	114	237406	5.000	ug/L	0.00
28) Chlorobenzene-d5	9.414	117	247138	5.000	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.809	152	111357	5.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.600	65	90444	3.917	ug/L	0.00
Spiked Amount 5.000	Range 40 - 130		Recovery = 78.400%			
7) Chloroethane-d5	1.916	69	79457	4.430	ug/L	0.00
Spiked Amount 5.000	Range 65 - 130		Recovery = 88.600%			
11) 1,1-Dichloroethene-d2	2.568	65	36057	4.356	ug/L	0.00
Spiked Amount 5.000	Range 60 - 125		Recovery = 87.200%			
20) 2-Butanone-d5	4.649	46	281131	67.163	ug/L	0.00
Spiked Amount 50.000	Range 40 - 130		Recovery = 134.320%#			
24) Chloroform-d	5.060	84	171381	5.051	ug/L	0.00
Spiked Amount 5.000	Range 70 - 125		Recovery = 101.000%			
26) 1,2-Dichloroethane-d4	5.700	65	89006	5.193	ug/L	0.00
Spiked Amount 5.000	Range 70 - 130		Recovery = 103.800%			
32) Benzene-d6	5.726	84	350753	4.781	ug/L	0.00
Spiked Amount 5.000	Range 70 - 125		Recovery = 95.600%			
36) 1,2-Dichloropropane-d6	6.690	67	116657	5.138	ug/L	0.00
Spiked Amount 5.000	Range 60 - 140		Recovery = 102.800%			
41) Toluene-d8	7.896	98	280998	4.233	ug/L	0.00
Spiked Amount 5.000	Range 70 - 130		Recovery = 84.600%			
43) trans-1,3-Dichloroprop...	8.179	79	35707	4.315	ug/L	0.00
Spiked Amount 5.000	Range 55 - 130		Recovery = 86.400%			
46) 2-Hexanone-d5	8.632	63	217474	62.018	ug/L	0.00
Spiked Amount 50.000	Range 45 - 130		Recovery = 124.040%			
56) 1,1,2,2-Tetrachloroeth...	10.754	84	107778	5.475	ug/L	0.00
Spiked Amount 5.000	Range 65 - 120		Recovery = 109.400%			
66) 1,2-Dichlorobenzene-d4	12.188	152	116327	5.456	ug/L	0.00
Spiked Amount 5.000	Range 80 - 120		Recovery = 109.200%			
Target Compounds						Qvalue
13) Acetone	2.684	43	15361m	5.404	ug/L	
14) Carbon disulfide	2.793	76	20263	0.340	ug/L	97
27) 1,2-Dichloroethane	5.796	62	25107	1.216	ug/L	96
42) Toluene	7.967	91	19788	0.239	ug/L	97

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

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