

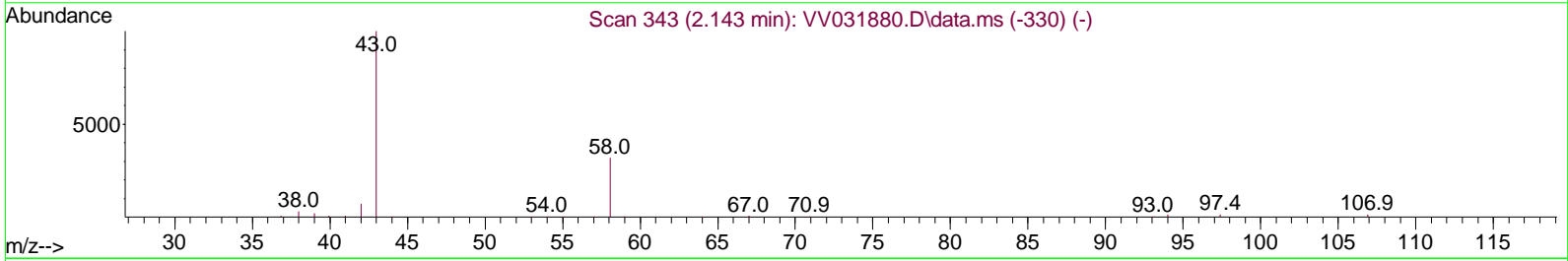
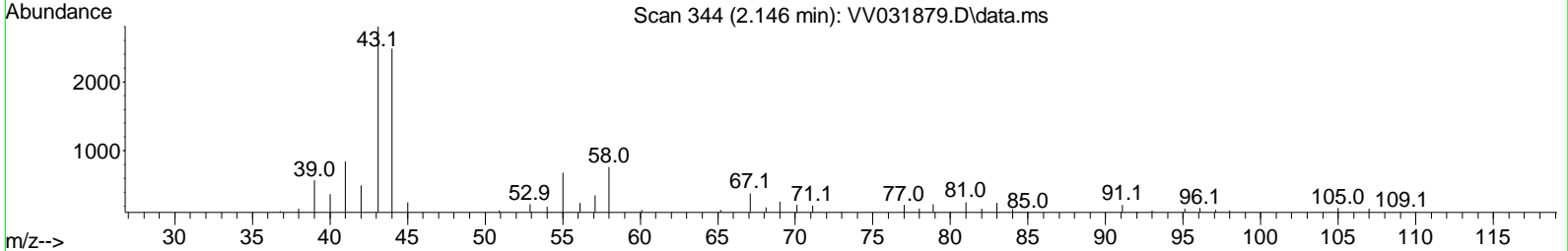
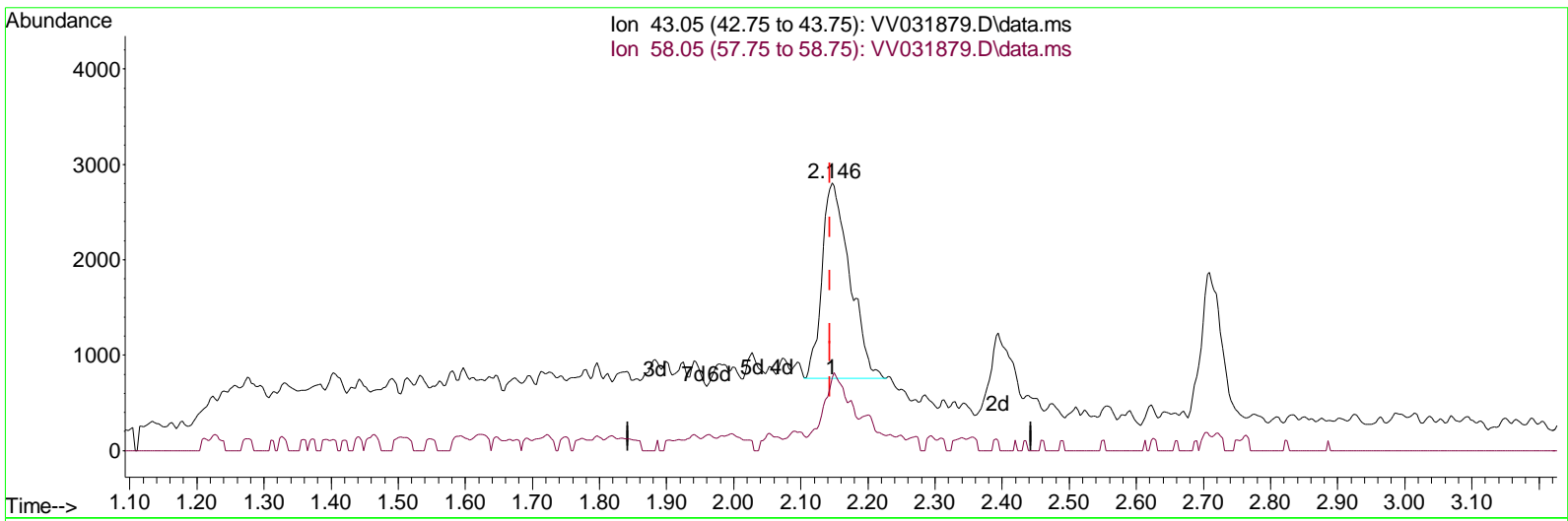
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV082423\
 Data File : VV031879.D
 Acq On : 23 Aug 2023 19:59
 Operator : SY/MD
 Sample : VSTD00173
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_V
ClientSampleId :
 VSTD001273

Manual Integrations APPROVED

Reviewed By :Romaben Patel 08/24/2023
 Supervised By :Mahesh Dadoda 08/24/2023

Quant Time: Aug 24 04:24:35 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR082323WMA.M
 Quant Title : TRACE VOA SFAM1.0
 QLast Update : Thu Aug 24 04:23:32 2023
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TIC: VV031879.D\data.ms

(13) Acetone (T)

2.146min (+ 0.003) 4.71 ug/L

response	5888
Ion	Exp% Act%
43.05	100.00 100.00
58.05	18.50 28.60
0.00	0.00 0.00
0.00	0.00 0.00

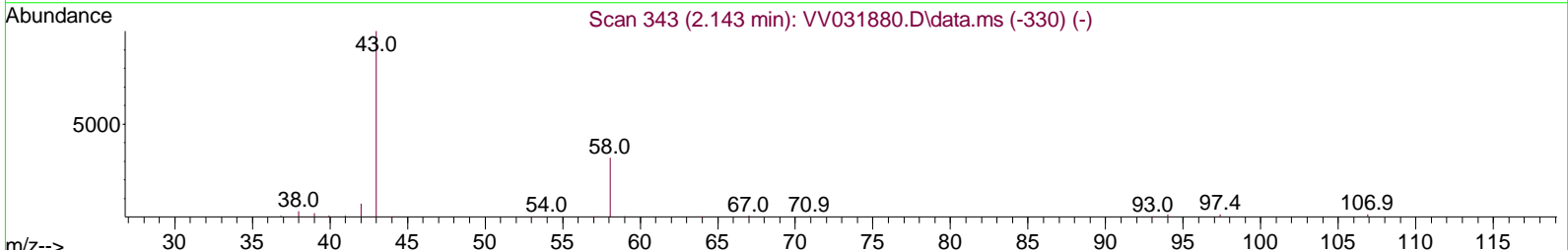
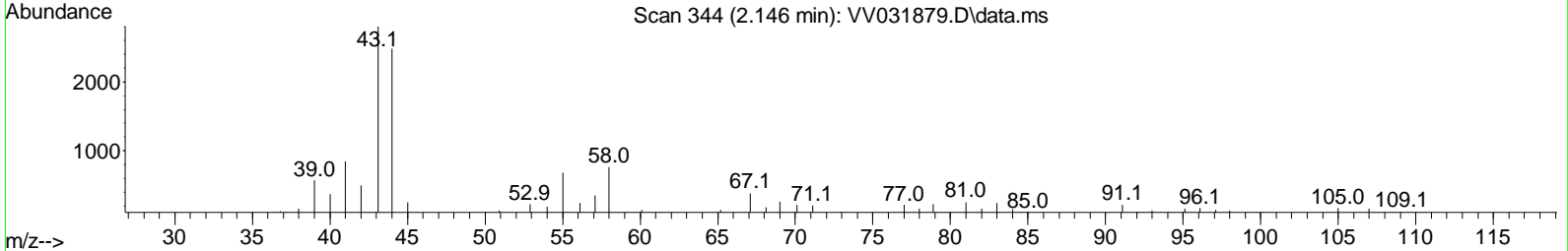
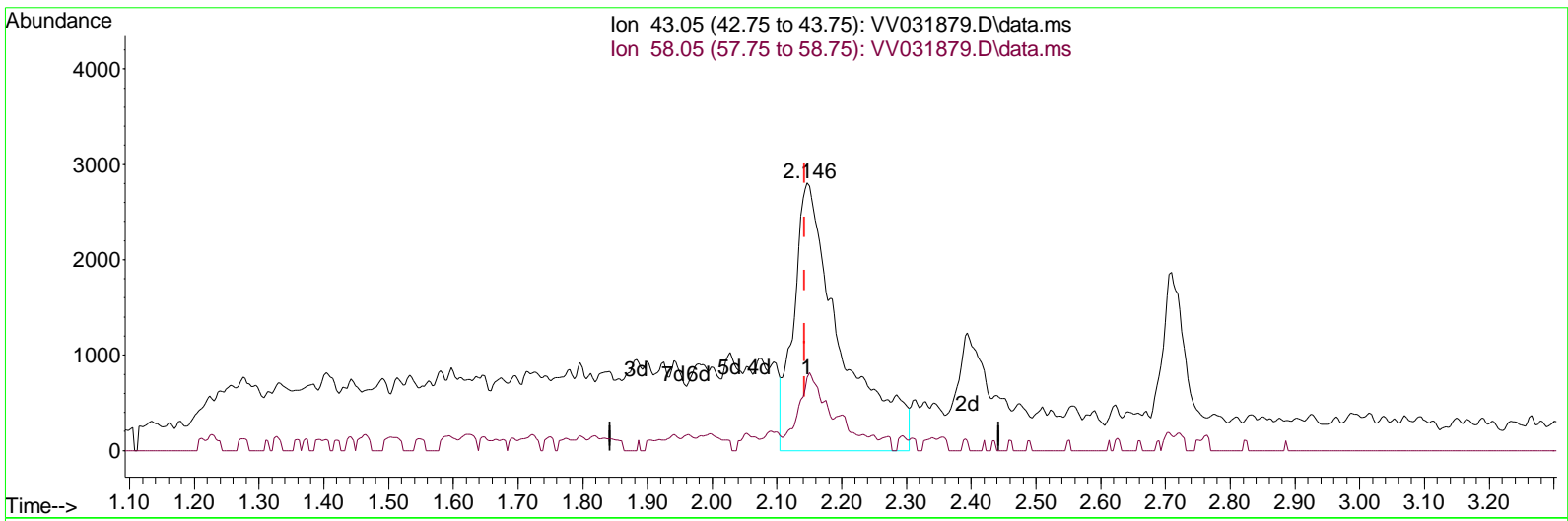
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TIC: VV031879.D\data.ms

(13) Acetone (T)

2.146min (+ 0.003) 11.30 ug/L m

response	14127
Ion	Exp% Act%
43.05	100.00 100.00
58.05	18.50 11.92
0.00	0.00 0.00
0.00	0.00 0.00

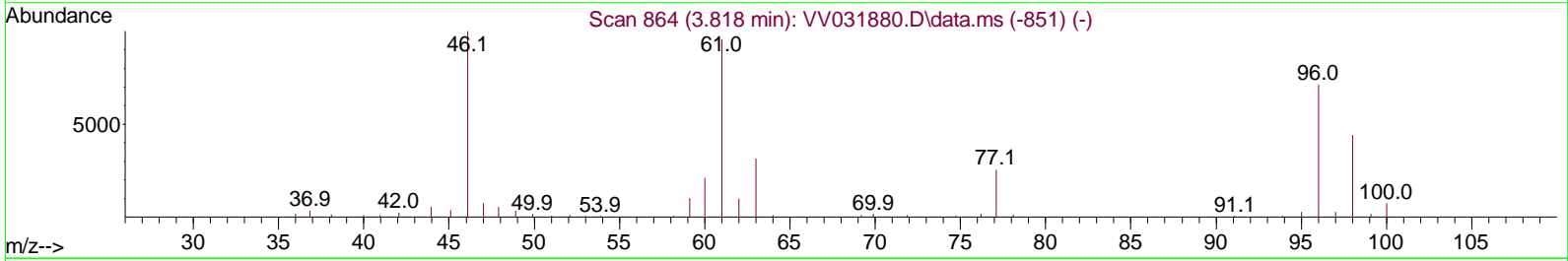
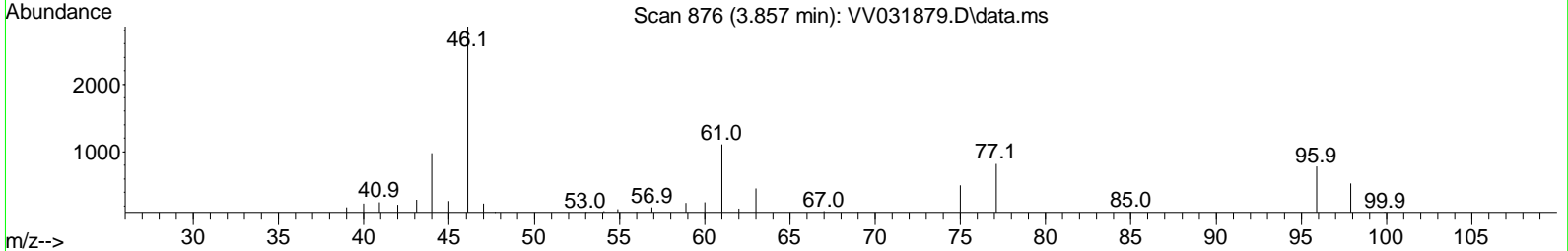
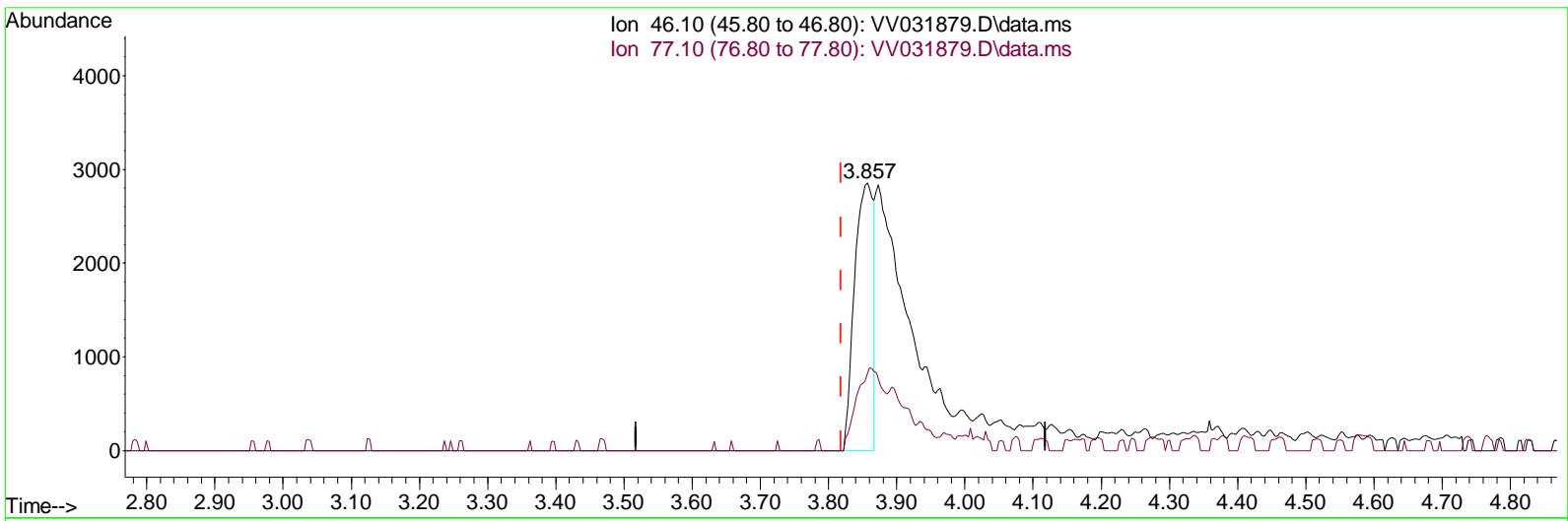
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TIC: VV031879.D\data.ms

(20) 2-Butanone-d5 (S)

3.857min (+ 0.039) 2.76 ug/L

response	5493	
Ion	Exp%	Act%
46.10	100.00	100.00
77.10	18.70	43.36#
0.00	0.00	0.00
0.00	0.00	0.00

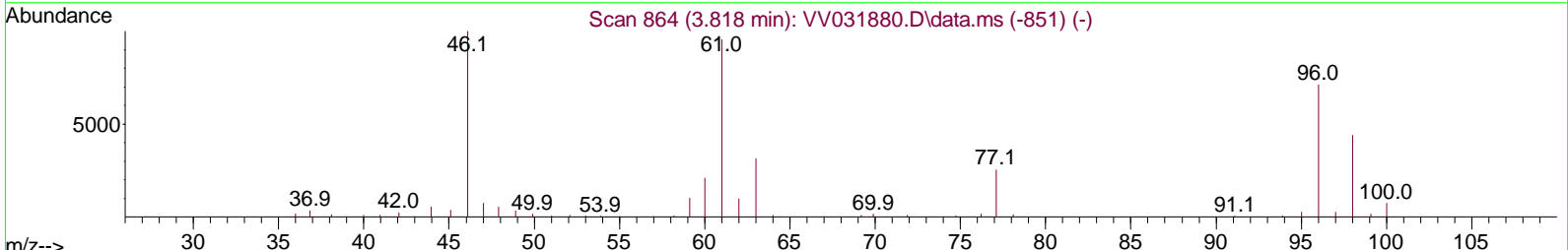
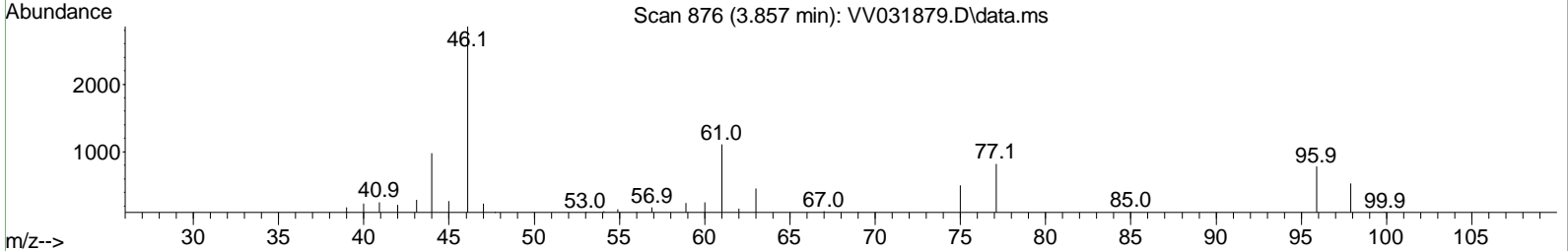
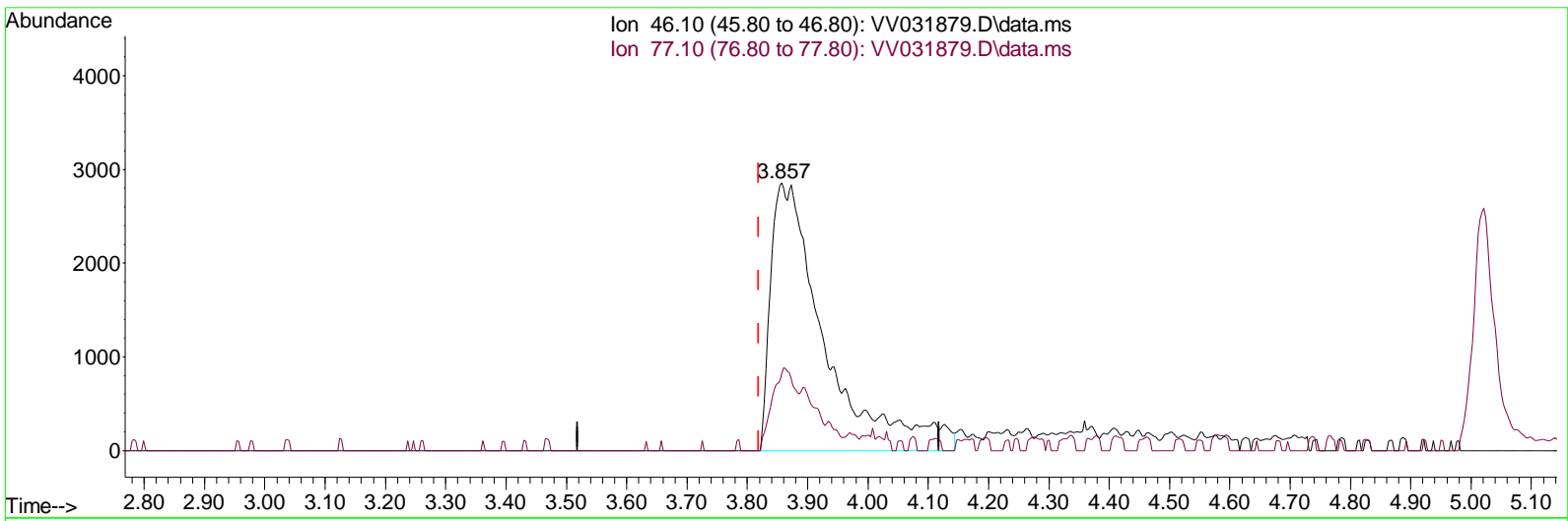
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TIC: VV031879.D\data.ms

(20) 2-Butanone-d5 (S)

3.857min (+ 0.039) 8.92 ug/L m

response	17773	
Ion	Exp%	Act%
46.10	100.00	100.00
77.10	18.70	13.40
0.00	0.00	0.00
0.00	0.00	0.00

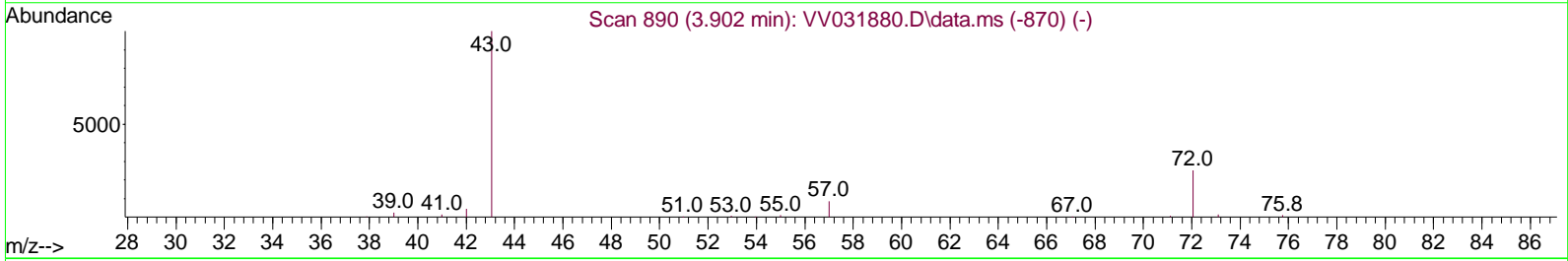
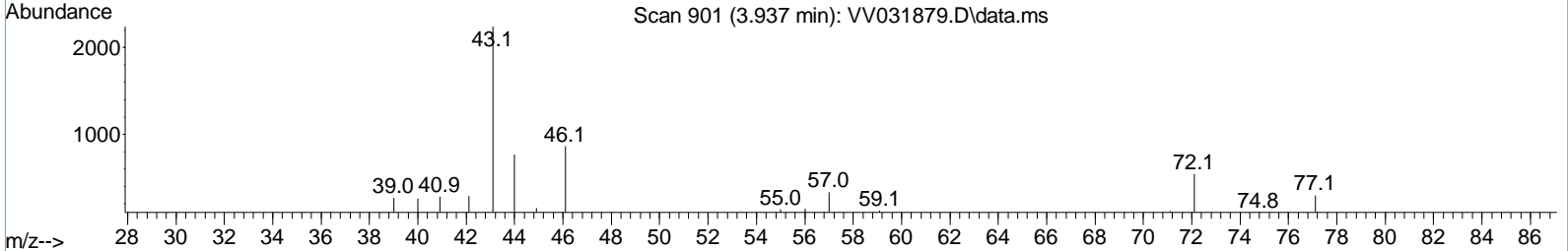
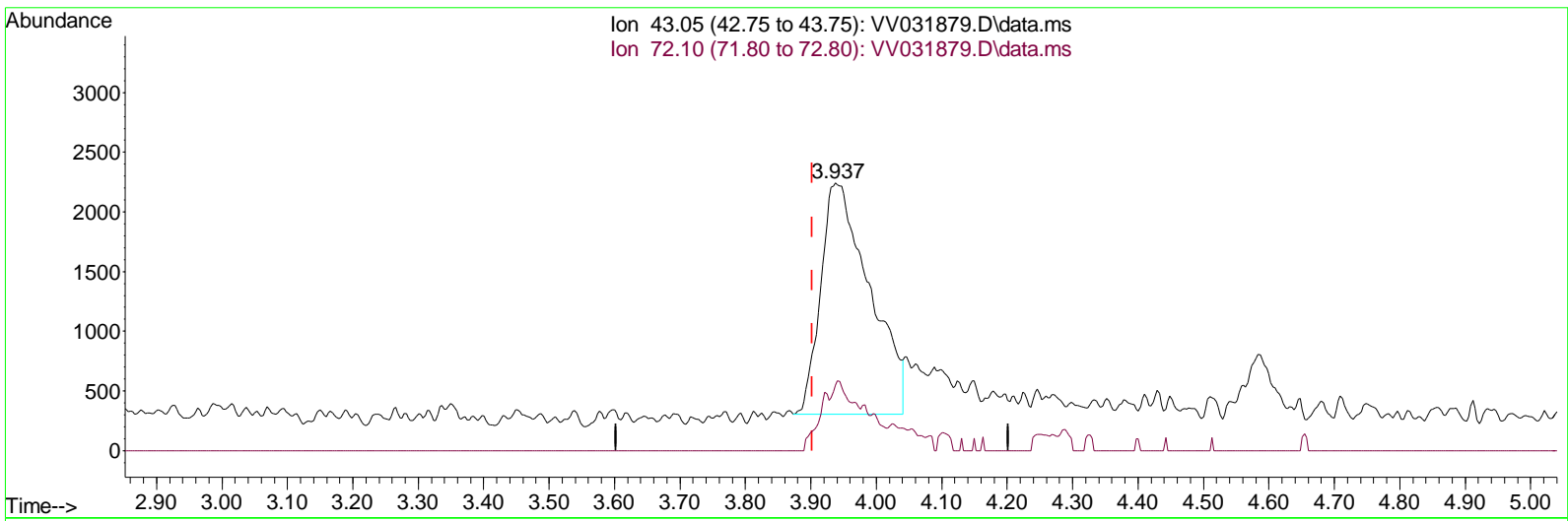
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(21) 2-Butanone (T)

3.937min (+ 0.035) 6.19 ug/L

response 10046

Ion	Exp%	Act%
43.05	100.00	100.00
72.10	18.00	5.83#
0.00	0.00	0.00
0.00	0.00	0.00

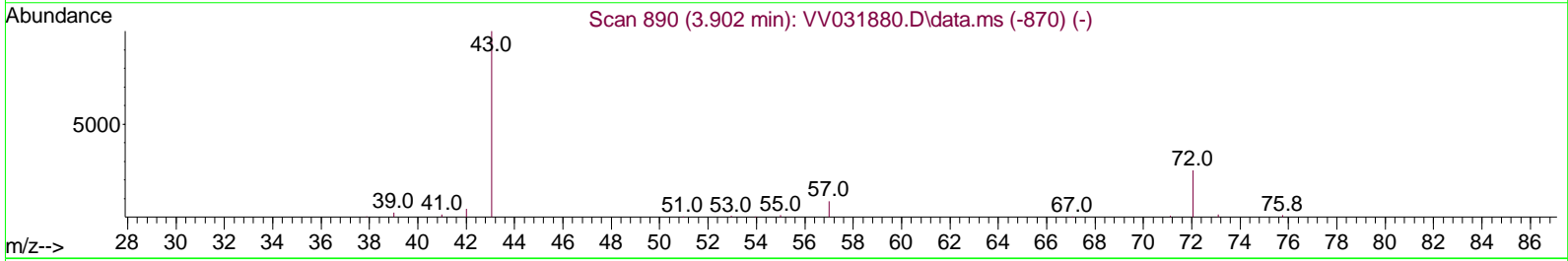
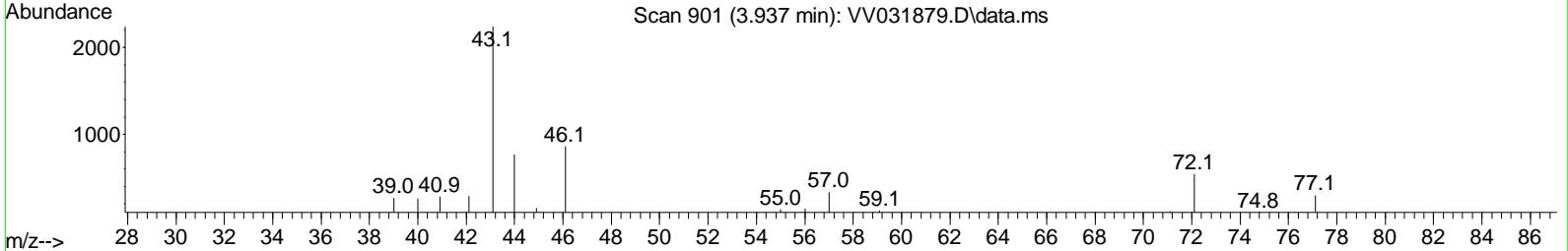
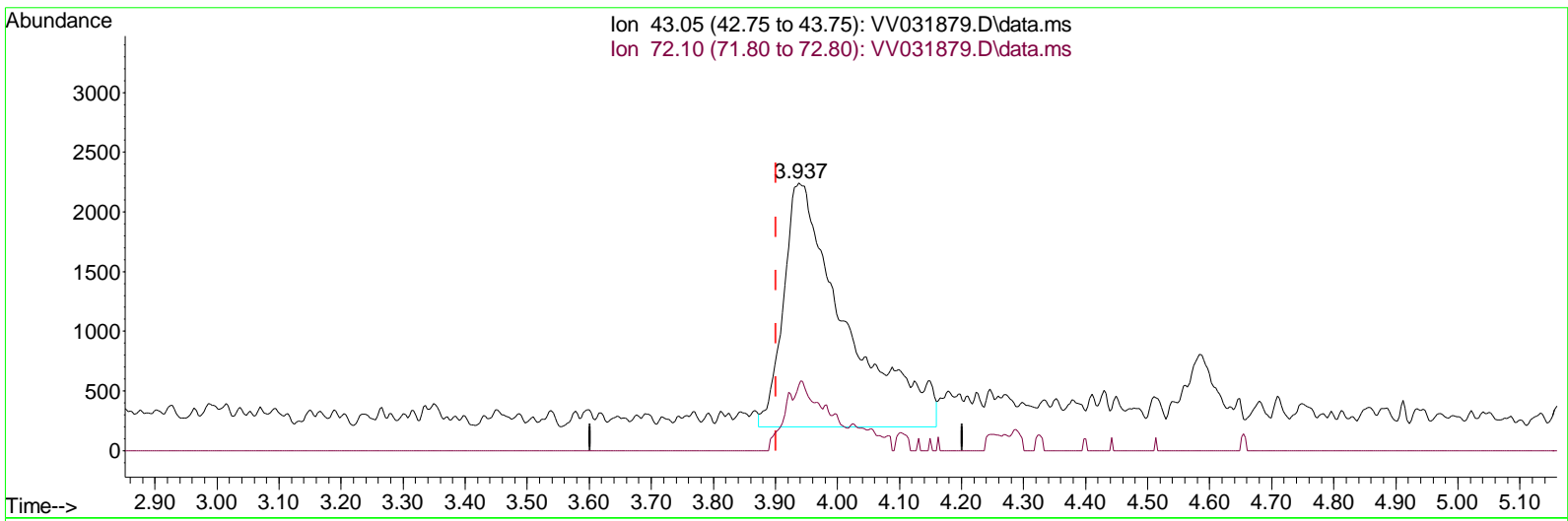
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TIC: VV031879.D\data.ms

(21) 2-Butanone (T)

3.937min (+ 0.035) 8.71 ug/L m

response	14122
Ion	Exp% Act%
43.05	100.00 100.00
72.10	18.00 4.15#
0.00	0.00 0.00
0.00	0.00 0.00

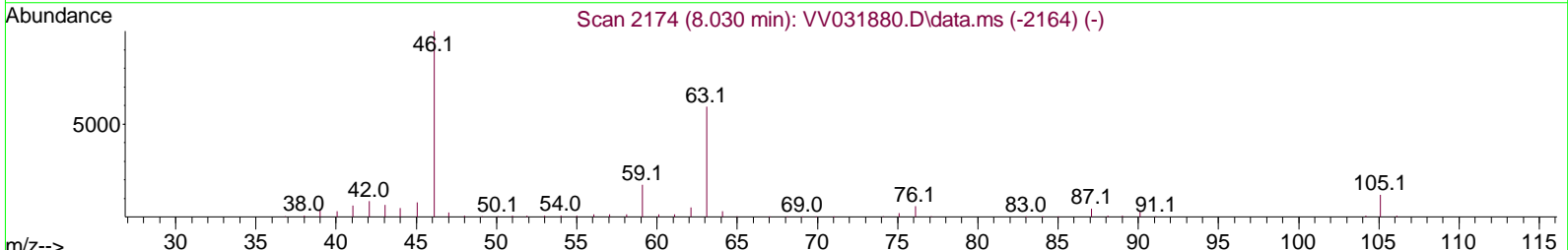
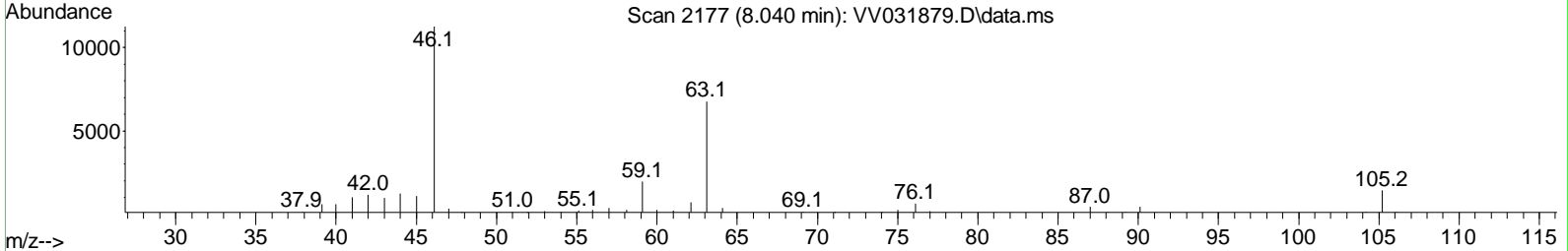
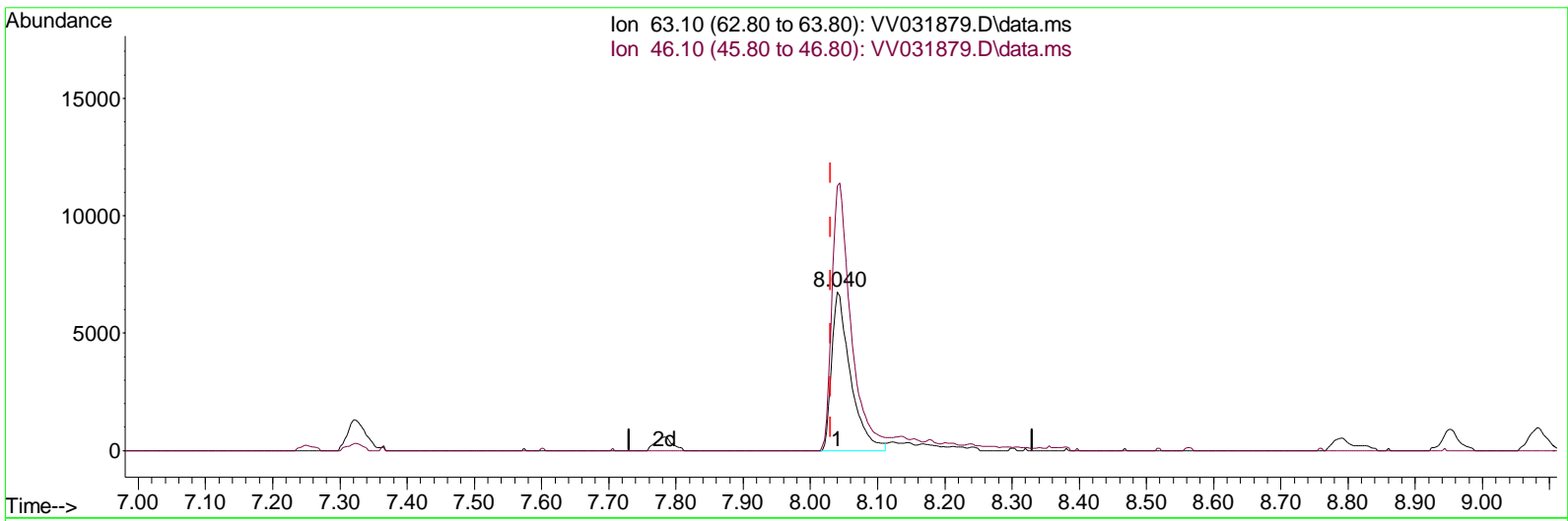
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TIC: VV031879.D\data.ms

(46) 2-Hexanone-d5 (S)

8.040min (+ 0.010) 7.37 ug/L

response	13113
Ion	Exp% Act%
63.10	100.00 100.00
46.10	172.40 173.80
0.00	0.00 0.00
0.00	0.00 0.00

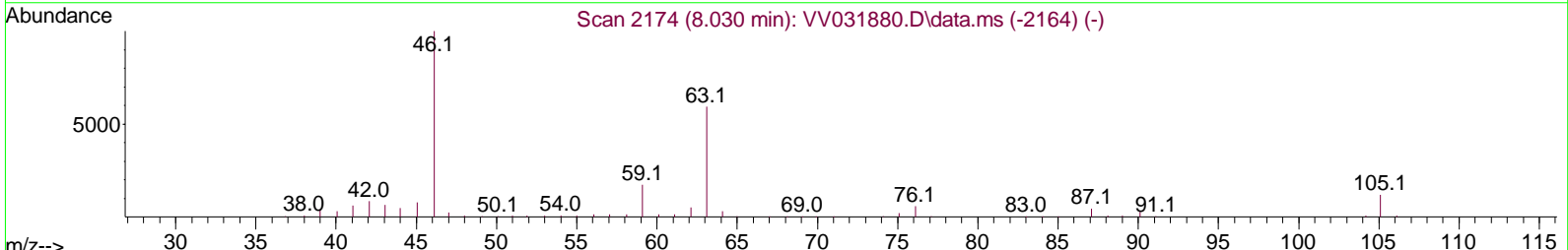
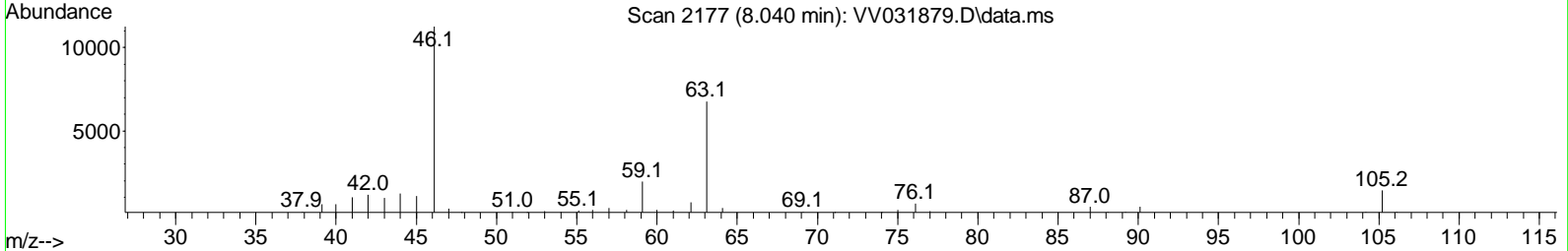
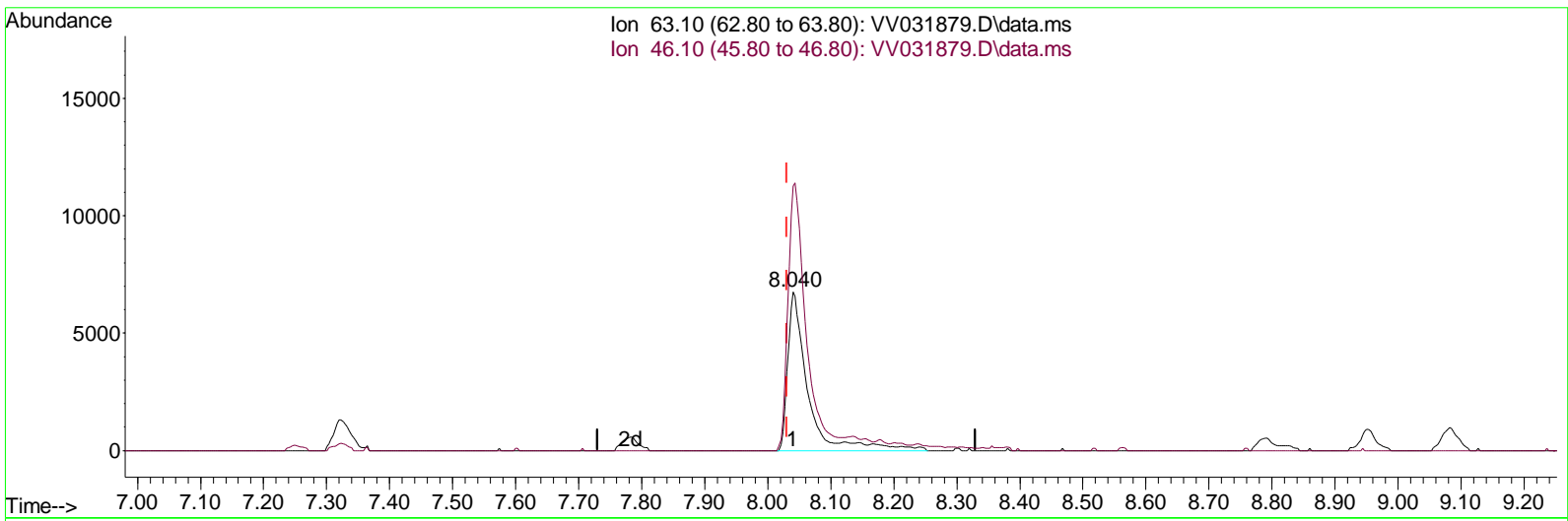
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 Data File : VV031879.D
 Acq On : 23 Aug 2023 19:59
 Operator : SY/MD
 Sample : VSTD00173
 Mi sc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_V
ClientSampleId :
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Manual IntegrationsAPPROVED

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 Response via : Initial Calibration



TIC: VV031879.D\data.ms

(46) 2-Hexanone-d5 (S)

8.040min (+ 0.010) 8.42 ug/L m

response 14982

Ion	Exp%	Act%
63.10	100.00	100.00
46.10	172.40	152.12
0.00	0.00	0.00
0.00	0.00	0.00

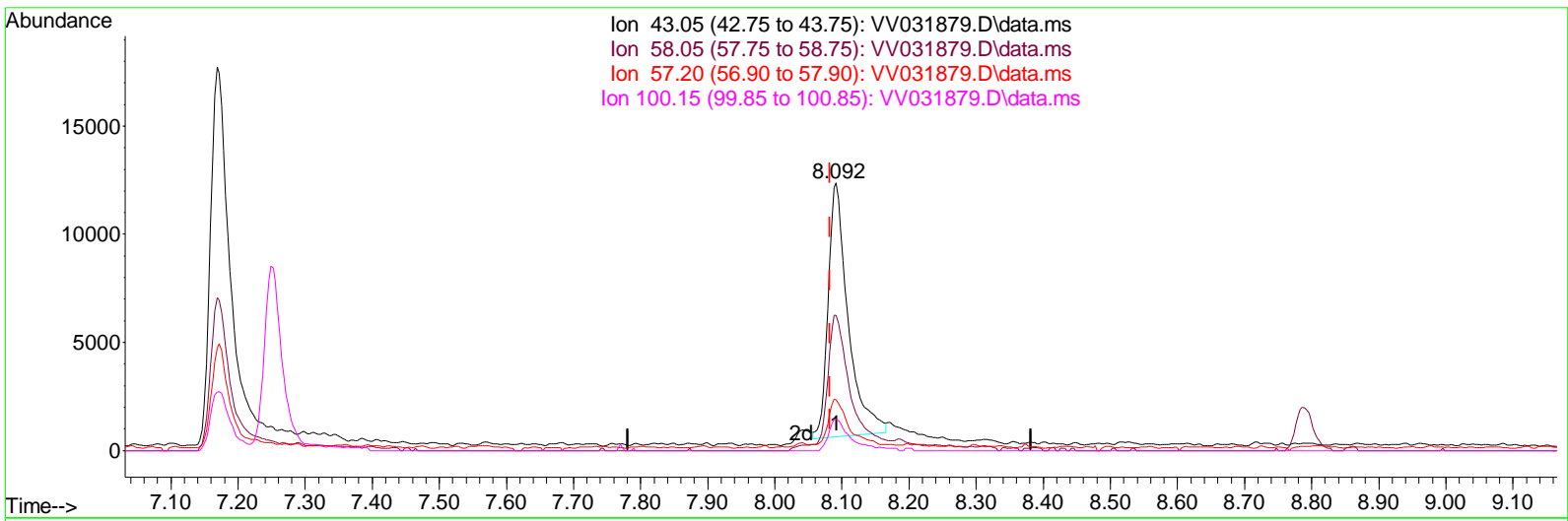
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 Data File : VV031879.D
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 ALS Vial : 27 Sample Multiplier: 1

Instrument :
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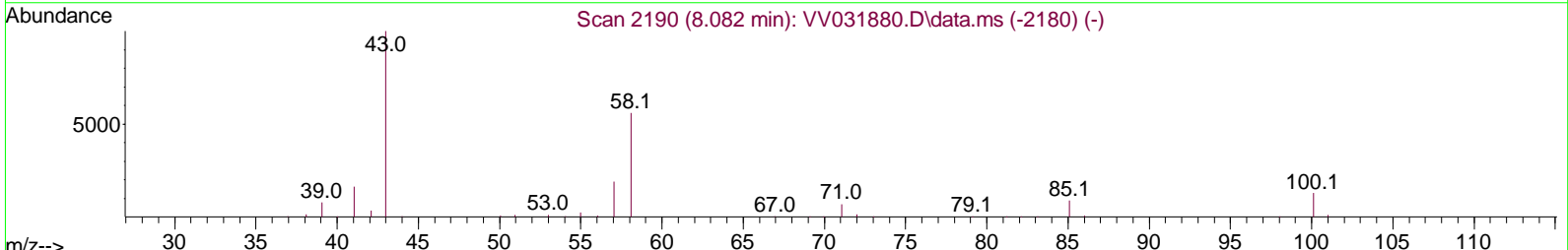
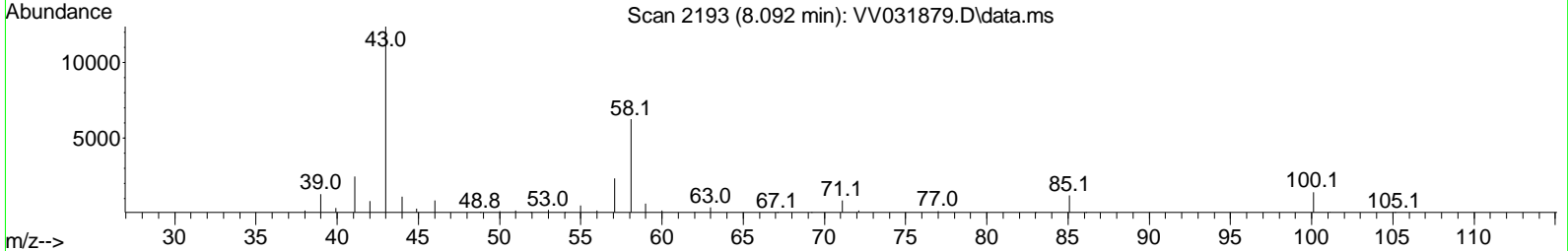
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Ion 43.05 (42.75 to 43.75): VV031879.D\data.ms
 Ion 58.05 (57.75 to 58.75): VV031879.D\data.ms
 Ion 57.20 (56.90 to 57.90): VV031879.D\data.ms
 Ion 100.15 (99.85 to 100.85): VV031879.D\data.ms



TIC: VV031879.D\data.ms

(48) 2-Hexanone (T)

8.092min (+ 0.010) 7.20 ug/L

response	24271	
Ion	Exp%	Act%
43.05	100.00	100.00
58.05	55.40	63.15
57.20	17.90	19.17
100.15	12.00	12.86

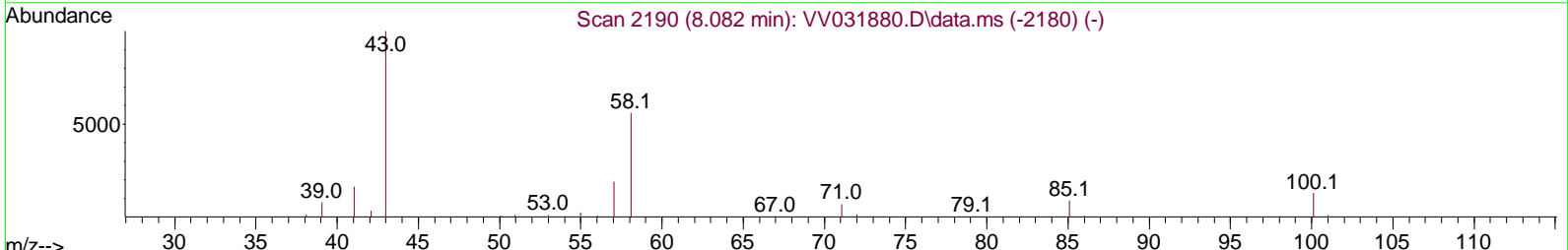
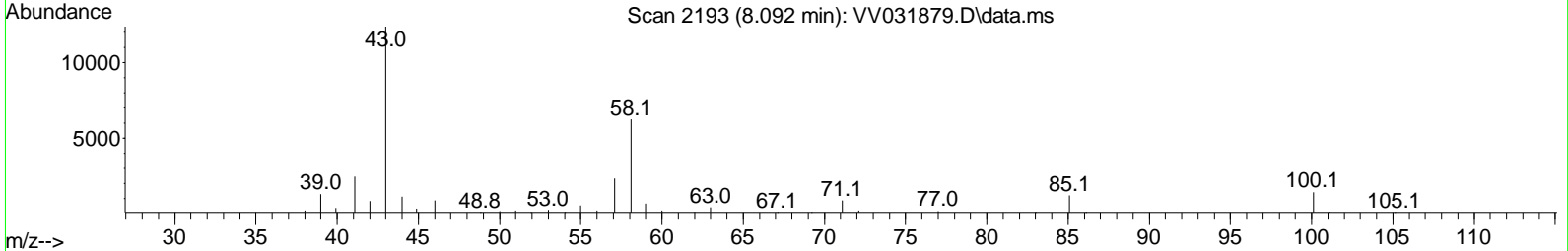
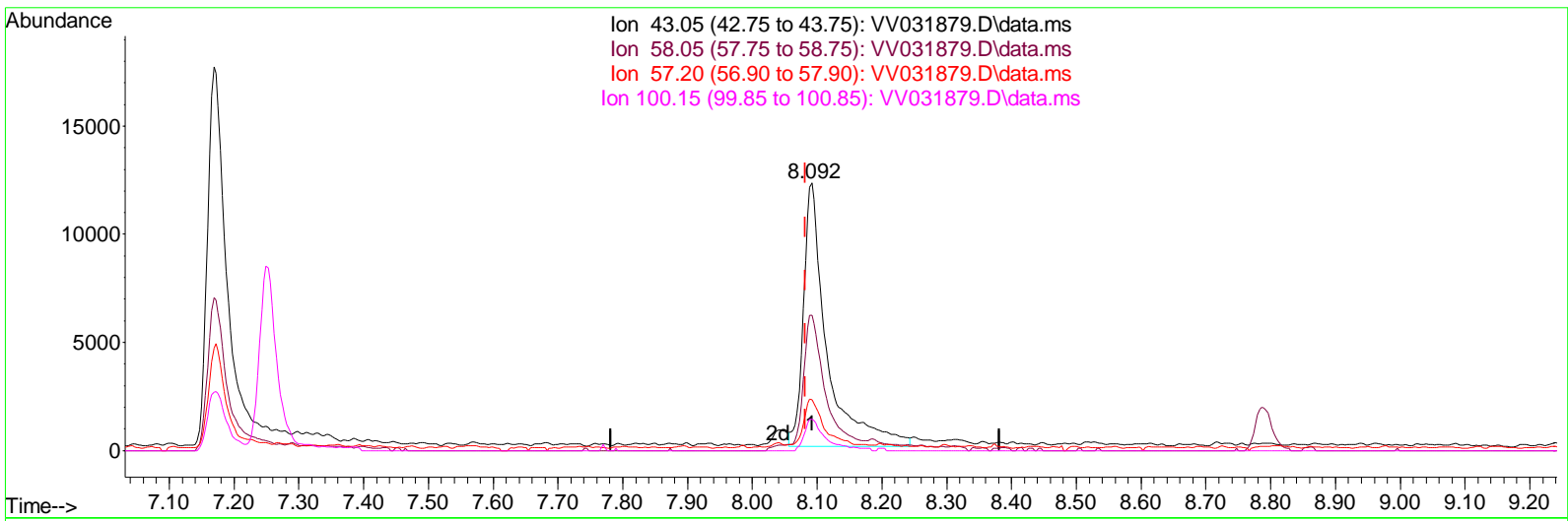
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TIC: VV031879.D\data.ms

(48) 2-Hexanone (T)

8.092min (+ 0.010) 9.16 ug/L m

response 30900

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	55.40	49.60
57.20	17.90	15.06
100.15	12.00	10.10

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 ALS Vial : 27 Sample Multi plier: 1

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Compound	R. T.	QI on	Response	Conc	Units	Dev(Mi n)	
Internal Standards							
1) 1,4-Di fl uorobenzene	5.539	114	120660	5.000	ug/L	0.00	
28) Chl orobenzene-d5	8.789	117	114043	5.000	ug/L	0.00	
58) 1,4-Di chl orobenzene-d4	11.188	152	60706	5.000	ug/L	0.00	
System Monitoring Compounds							
4) Vi nyl Chl ori de-d3	1.278	65	4346	1.066	ug/L	0.00	
7) Chl oroethane-d5	1.536	69	4509	0.966	ug/L	0.00	
11) 1,1-Di chl oroethene-d2	2.060	65	2474	1.032	ug/L	0.00	
20) 2-Butanone-d5	3.857	46	17773m	8.923	ug/L	0.04	
24) Chl oroform-d	4.259	84	14275	1.035	ug/L	0.00	
26) 1,2-Di chl oroethane-d4	4.957	65	7174	1.009	ug/L	0.00	
32) Benzene-d6	4.970	84	26626	1.026	ug/L	0.00	
36) 1,2-Di chl oropropane-d6	5.998	67	9027	0.986	ug/L	0.00	
41) Tol uene-d8	7.252	98	24106	1.001	ug/L	0.00	
43) trans-1,3-Di chl oroprop. . .	7.567	79	3142	0.942	ug/L	0.00	
46) 2-Hexanone-d5	8.040	63	14982m	8.421	ug/L	0.00	
56) 1,1,2,2-Tetrchl oroeth. . .	10.159	84	6257	0.953	ug/L	0.00	
66) 1,2-Di chl orobenzene-d4	11.567	152	9749	0.992	ug/L	0.00	
Target Compounds							
							Qval ue
2) Di chl orodi fl uoromethane	1.108	85	6796	0.986	ug/L		100
3) Chl oromethane	1.217	50	7408	1.003	ug/L		97
5) Vi nyl chl ori de	1.285	62	7606	1.005	ug/L		98
6) Bromomethane	1.491	94	4752	1.005	ug/L		94
8) Chl oroethane	1.552	64	4807	0.995	ug/L		91
9) Tri chl orofl uoromethane	1.716	101	9823	1.006	ug/L		94
10) 1,1,2-Tri chl oro-1,2,2- . . .	2.069	101	6170	1.060	ug/L		96
12) 1,1-Di chl oroethene	2.072	96	5386	0.992	ug/L		91
13) Acetone	2.146	43	14127m	11.301	ug/L		
14) Carbon di sul fi de	2.243	76	14427	0.944	ug/L		100
15) Methyl Acetate	2.394	43	1935	0.817	ug/L #		82
16) Methyl ene chl ori de	2.449	84	10153	1.149	ug/L		96
17) Methyl tert-butyl Ether	2.712	73	15723	0.933	ug/L		97
18) trans-1,2-Di chl oroethene	2.699	96	5929	0.963	ug/L		94
19) 1,1-Di chl oroethane	3.117	63	12351	0.974	ug/L		97
21) 2-Butanone	3.937	43	14122m	8.706	ug/L		
22) ci s-1,2-Di chl oroethene	3.828	96	7084	0.999	ug/L		96
23) Bromochl oromethane	4.162	128	2643	0.939	ug/L		96
25) Chl oroform	4.288	83	12604	0.975	ug/L		96
27) 1,2-Di chl oroethane	5.056	62	7223	0.934	ug/L		98
29) 1,1,1-Tri chl oroethane	4.523	97	11407	1.036	ug/L		96
30) Cycl ohexane	4.587	56	11452	1.012	ug/L		97
31) Carbon tetrchl ori de	4.741	117	9134	1.025	ug/L		95
33) Benzene	5.021	78	25997	0.990	ug/L		100
34) Tri chl oroethene	5.844	95	7343	0.956	ug/L		97
35) Methyl cycl ohexane	6.053	83	11859	1.061	ug/L		96
37) 1,2-Di chl oropropane	6.101	63	7335	1.013	ug/L		98
38) Bromodi chl oromethane	6.439	83	9039	0.965	ug/L		94
39) ci s-1,3-Di chl oropropene	6.969	75	9679	0.879	ug/L		100
40) 4-Methyl -2-pentanone	7.169	43	36936	8.142	ug/L		92
42) Tol uene	7.323	91	28325	0.971	ug/L		100

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV082423\
 Data File : VV031879.D
 Acq On : 23 Aug 2023 19:59
 Operator : SY/MD
 Sample : VSTD00173
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_V
ClientSampleId :
 VSTD001273

Manual Integrations APPROVED

Reviewed By : Romaben Patel 08/24/2023
 Supervised By : Mahesh Dadoda 08/24/2023

Quant Time: Aug 24 04:24:35 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR082323WMA.M
 Quant Title : TRACE VOA SFAM1.0
 QLast Update : Thu Aug 24 04:23:32 2023
 Response via : Initial Calibration

Compound	R. T.	QI on	Response	Conc	Units	Dev(Min)
44) trans-1,3-Dichloropropene	7.593	75	8809	0.958	ug/L	97
45) 1,1,2-Trichloroethane	7.780	97	4579	0.884	ug/L	97
47) Tetrachloroethene	7.908	164	4922	1.003	ug/L	97
48) 2-Hexanone	8.092	43	30900m	9.162	ug/L	
49) Dibromochloromethane	8.182	129	4990	0.924	ug/L	95
50) 1,2-Dibromoethane	8.294	107	4278	0.913	ug/L #	99
51) Chlorobenzene	8.821	112	18201	0.985	ug/L	98
52) Ethyl benzene	8.953	91	33743	1.006	ug/L	98
53) m,p-Xylene	9.082	106	12752	1.009	ug/L	99
54) o-Xylene	9.484	106	12029	0.985	ug/L	97
55) Styrene	9.503	104	20395	0.968	ug/L	99
57) 1,1,2,2-Tetrachloroethane	10.182	83	4901	0.934	ug/L #	89
59) Bromoform	9.673	173	2542	0.886	ug/L #	90
60) Isopropyl benzene	9.873	105	33689	0.998	ug/L	99
61) 1,2,3-Trichloropropane	10.220	75	4109	0.919	ug/L #	93
62) 1,3,5-Trimethyl benzene	10.481	105	28711	0.993	ug/L	100
63) 1,2,4-Trimethyl benzene	10.857	105	28562	1.005	ug/L	98
64) 1,3-Dichlorobenzene	11.124	146	14867	0.995	ug/L	97
65) 1,4-Dichlorobenzene	11.214	146	15497	1.009	ug/L	95
67) 1,2-Dichlorobenzene	11.587	146	13006	0.951	ug/L	99
68) 1,2-Dibromo-3-chloropropane	12.371	75	925	0.913	ug/L #	80
69) 1,3,5-Trimethyl chlorobenzene	12.590	180	11164	1.005	ug/L	100
70) 1,2,4-trimethyl chlorobenzene	13.204	180	9828	0.970	ug/L	98
71) Naphthalene	13.448	128	17768	0.931	ug/L	99
72) 1,2,3-Trimethyl chlorobenzene	13.686	180	7930	0.930	ug/L	100

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

