

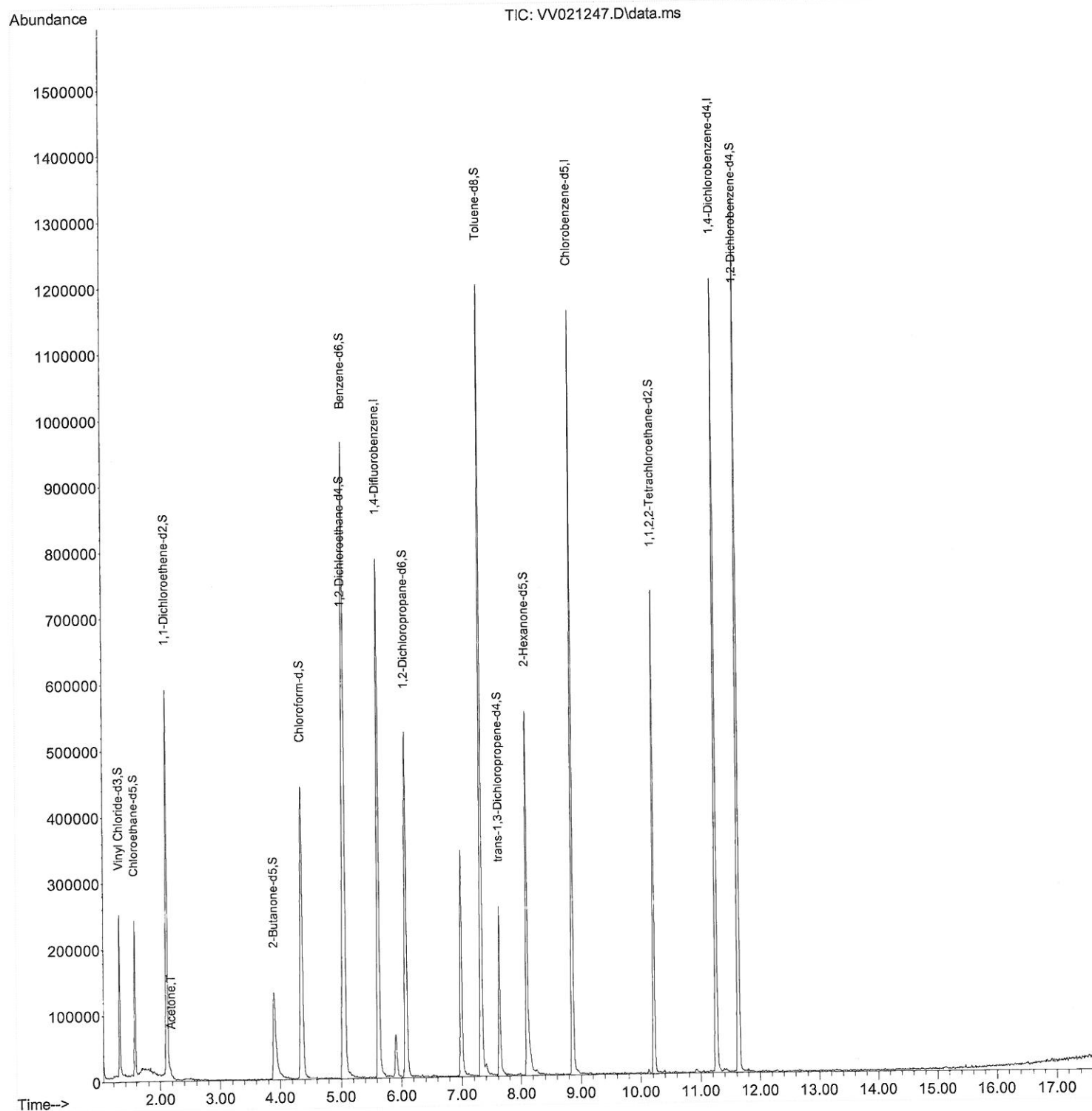
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050321\
Data File : VV021247.D
Acq On : 03 May 2021 15:51
Operator : SY/MD
Sample : M2209-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG7C3

Manual Integrations
APPROVED

MMDadoda
5/4/2021 11:49:08 AM

Quant Time: May 04 05:25:50 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM050321WMA.M
Quant Title : VOC Analysis
QLast Update : Mon May 03 18:54:51 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

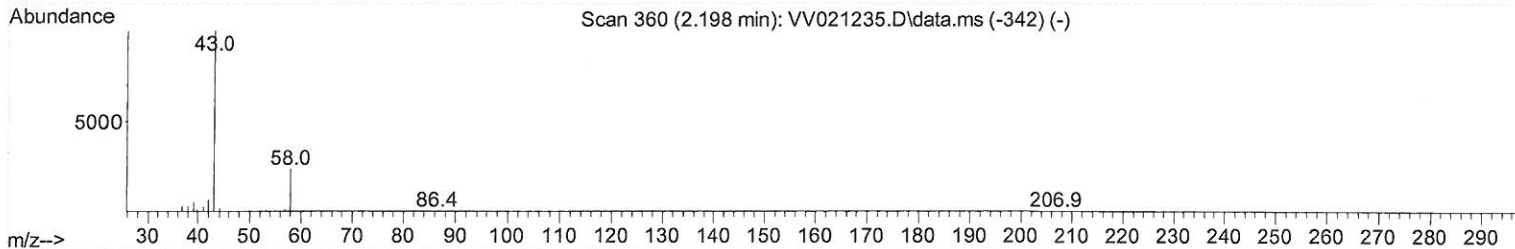
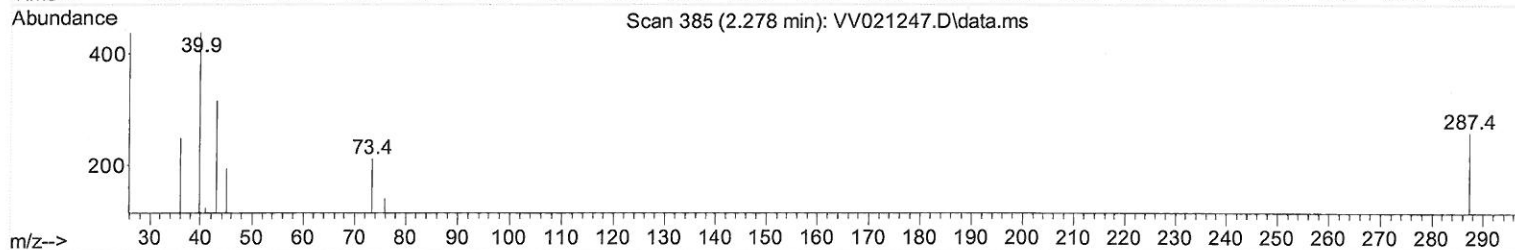
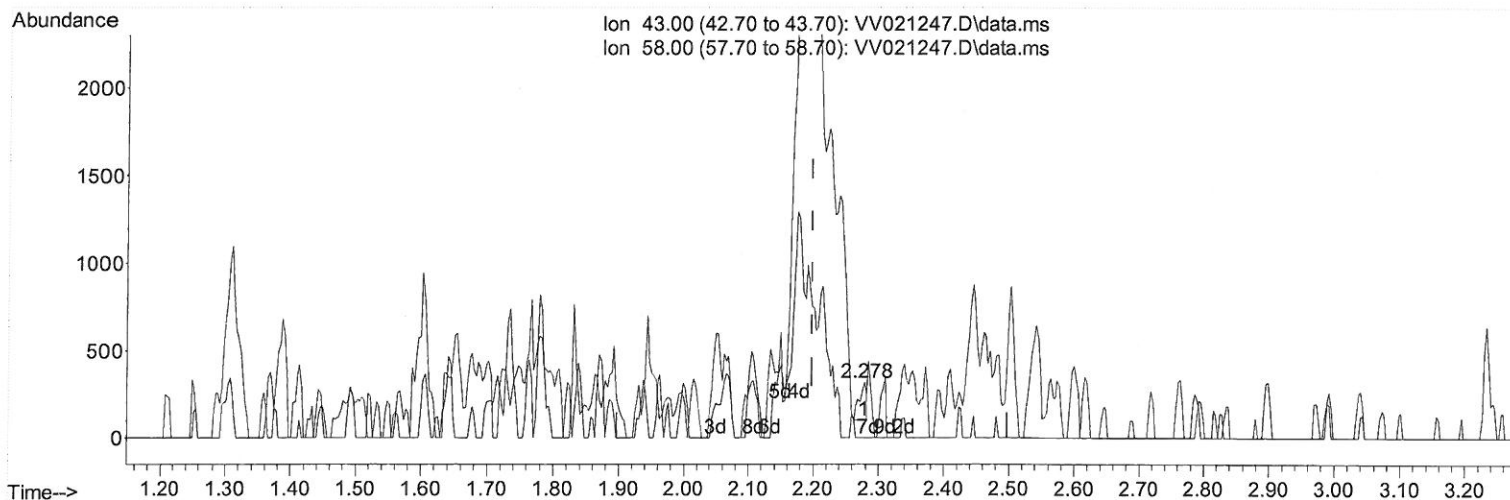
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TIC: VV021247.D\data.ms

(13) Acetone (T)

2.278min (+ 0.080) 0.06 ug/L

response 162

Ion	Exp%	Act%
43.00	100.00	100.00
58.00	15.40	28.40
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

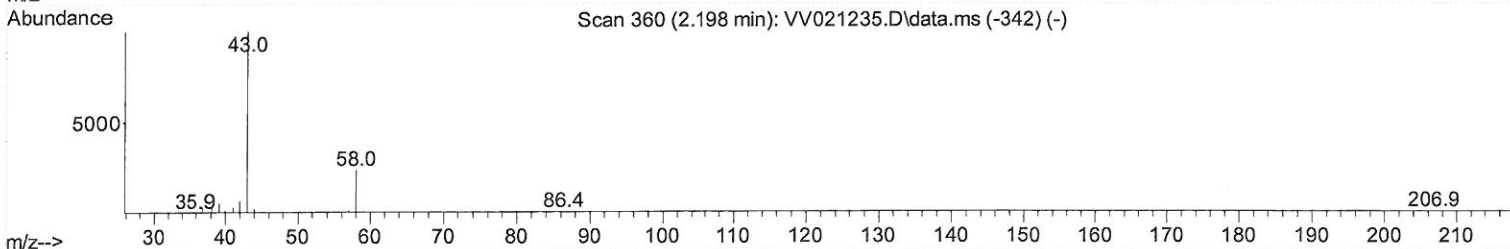
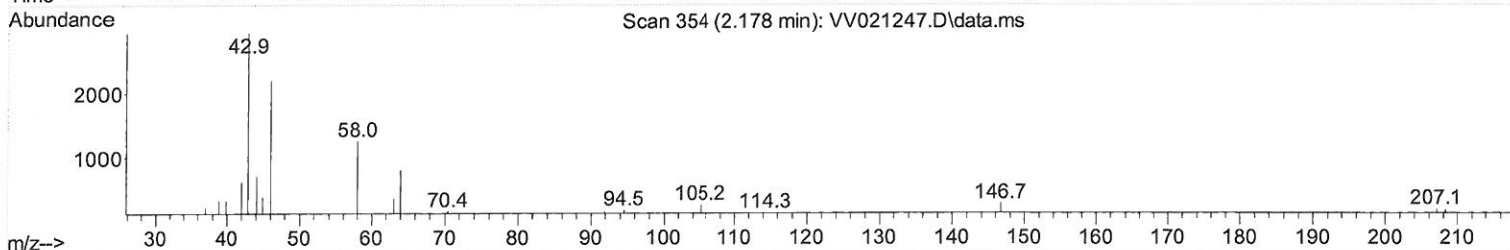
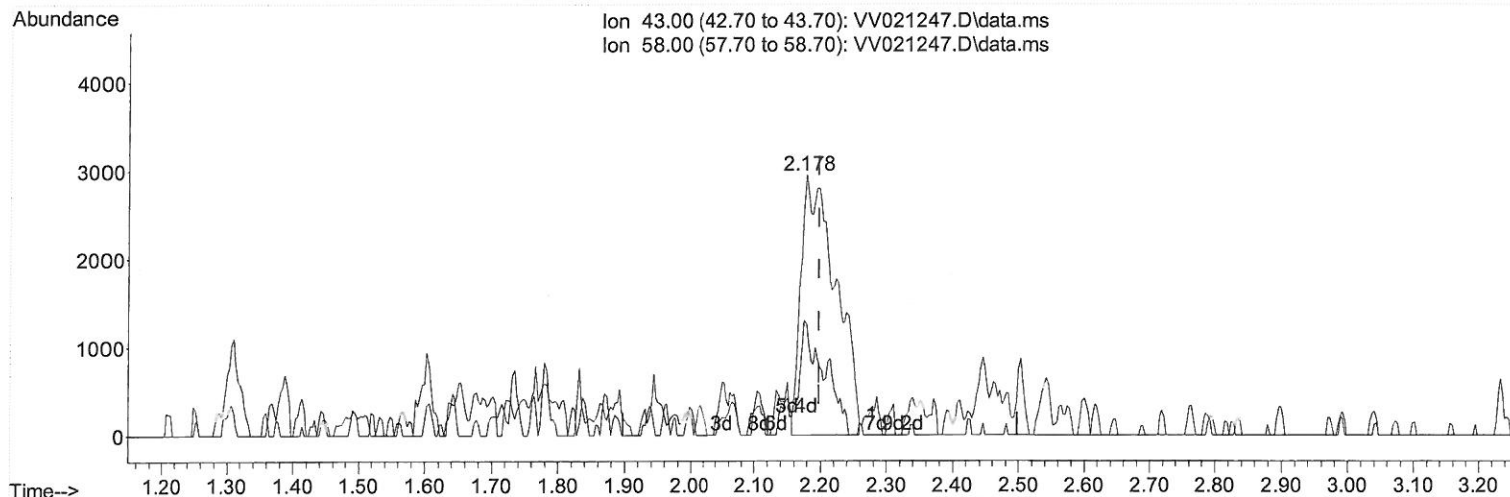
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 QLast Update : Mon May 03 18:54:51 2021
 Response via : Initial Calibration



TIC: VV021247.D\data.ms

(13) Acetone (T)

2.178min (-0.019) 3.94 ug/L m

response 10767

Ion	Exp%	Act%
43.00	100.00	100.00
58.00	15.40	0.43
0.00	0.00	0.00
0.00	0.00	0.00

Handwritten signature: MMD
 5/10/21

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VW050321\
 Data File : VW021247.D
 Acq On : 03 May 2021 15:51
 Operator : SY/MD
 Sample : M2209-10
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BG7C3

Manual Integrations
 APPROVED

MMDadoda
 5/4/2021 11:49:08 AM

Quant Time: May 04 05:25:50 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM050321WMA.M
 Quant Title : VOC Analysis
 QLast Update : Mon May 03 18:54:51 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Difluorobenzene	5.622	114	648889	50.00	ug/L	0.00
28) Chlorobenzene-d5	8.857	117	618376	50.00	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.255	152	318902	50.00	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.307	65	144586	50.23	ug/L	0.00
Spiked Amount	50.000	Range 60 - 135	Recovery	= 100.46%		
7) Chloroethane-d5	1.568	69	133607	51.94	ug/L	0.00
Spiked Amount	50.000	Range 70 - 130	Recovery	= 103.88%		
11) 1,1-Dichloroethene-d2	2.108	63	289808	36.09	ug/L	0.00
Spiked Amount	50.000	Range 60 - 125	Recovery	= 72.18%		
21) 2-Butanone-d5	3.895	46	236593	102.43	ug/L	0.00
Spiked Amount	100.000	Range 40 - 130	Recovery	= 102.43%		
24) Chloroform-d	4.352	84	449988	49.12	ug/L	0.00
Spiked Amount	50.000	Range 70 - 125	Recovery	= 98.24%		
26) 1,2-Dichloroethane-d4	5.034	65	347563	50.68	ug/L	0.00
Spiked Amount	50.000	Range 70 - 125	Recovery	= 101.36%		
32) Benzene-d6	5.053	84	745332	52.34	ug/L	0.00
Spiked Amount	50.000	Range 70 - 125	Recovery	= 104.68%		
36) 1,2-Dichloropropane-d6	6.072	67	192099	52.51	ug/L	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	= 105.02%		
41) Toluene-d8	7.320	98	756058	51.23	ug/L	0.00
Spiked Amount	50.000	Range 80 - 120	Recovery	= 102.46%		
43) trans-1,3-Dichloroprop...	7.625	79	136743	50.26	ug/L	0.00
Spiked Amount	50.000	Range 60 - 125	Recovery	= 100.52%		
47) 2-Hexanone-d5	8.095	63	166971	98.84	ug/L	0.00
Spiked Amount	100.000	Range 45 - 130	Recovery	= 98.84%		
56) 1,1,2,2-Tetrachloroeth...	10.220	84	278146	47.51	ug/L	0.00
Spiked Amount	50.000	Range 65 - 120	Recovery	= 95.02%		
66) 1,2-Dichlorobenzene-d4	11.631	152	330970	52.66	ug/L	0.00
Spiked Amount	50.000	Range 80 - 120	Recovery	= 105.32%		
Target Compounds						
13) Acetone	2.178	43	10767m	3.94	ug/L	Qvalue

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