

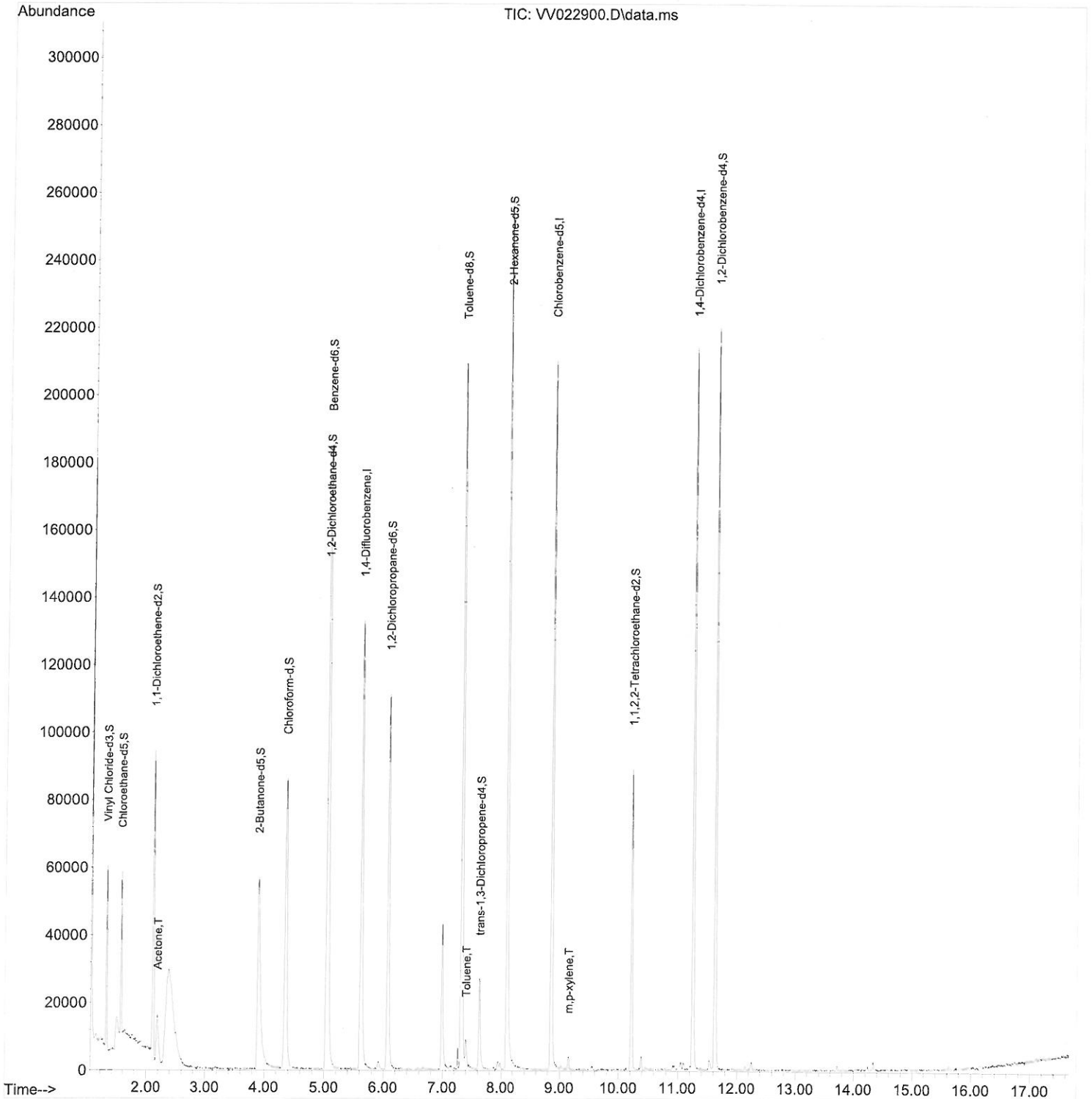
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101821\
Data File : VV022900.D
Acq On : 19 Oct 2021 05:05
Operator : SY/MD
Sample : M4137-15
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BFGD3

Manual Integrations
APPROVED

MMDadoda
10/19/2021 3:48:36 PM

Quant Time: Oct 19 09:00:24 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100721WMA.M
Quant Title : TRACE VOA SFAM1.0
QLast Update : Tue Oct 19 04:07:06 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

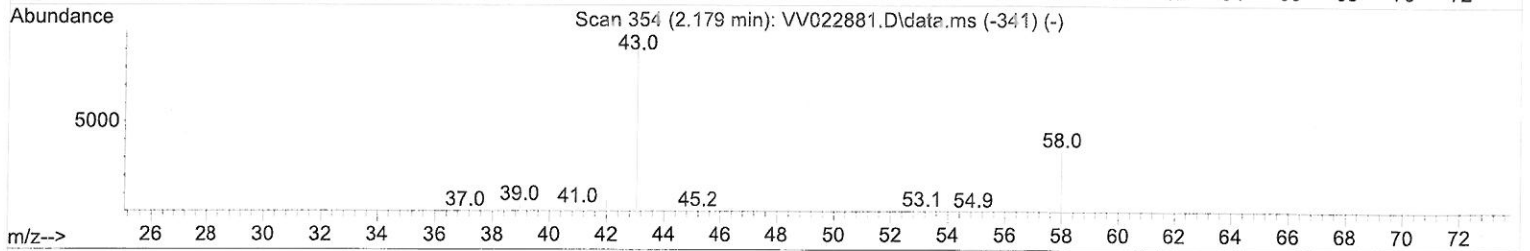
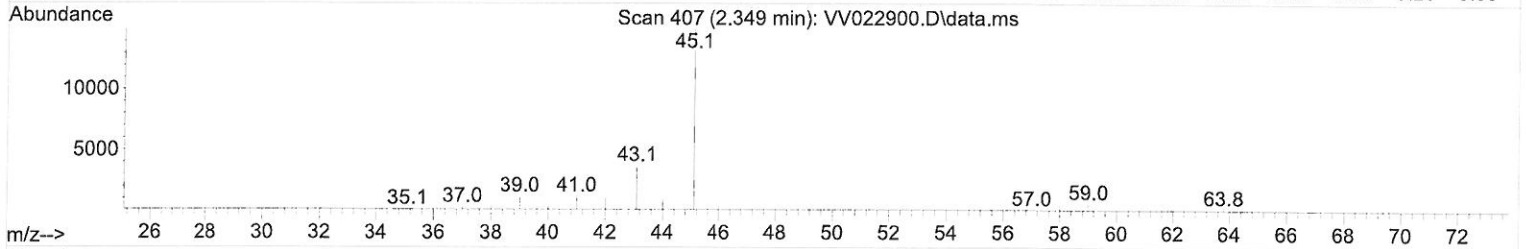
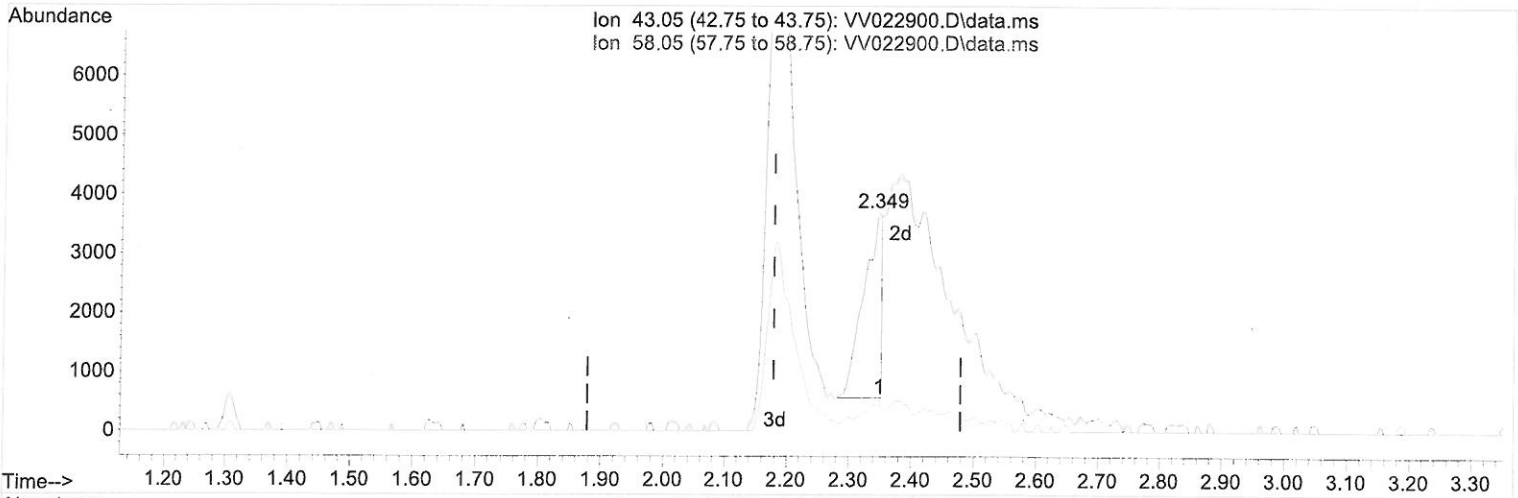
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TIC: VV022900.D\data.ms

(13) Acetone (T)

2.349min (+ 0.170) 5.32 ug/L

response 6378

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	12.10	7.96
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

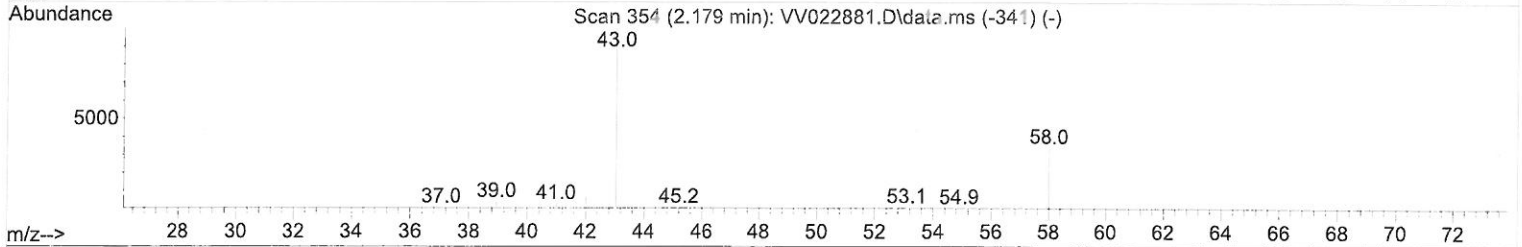
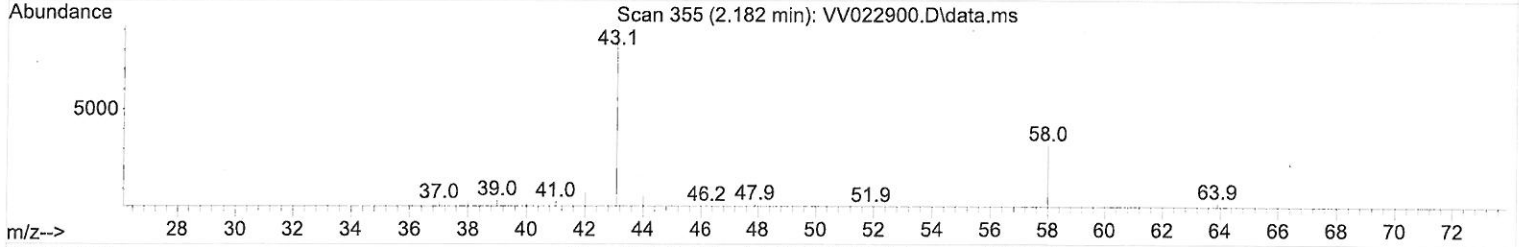
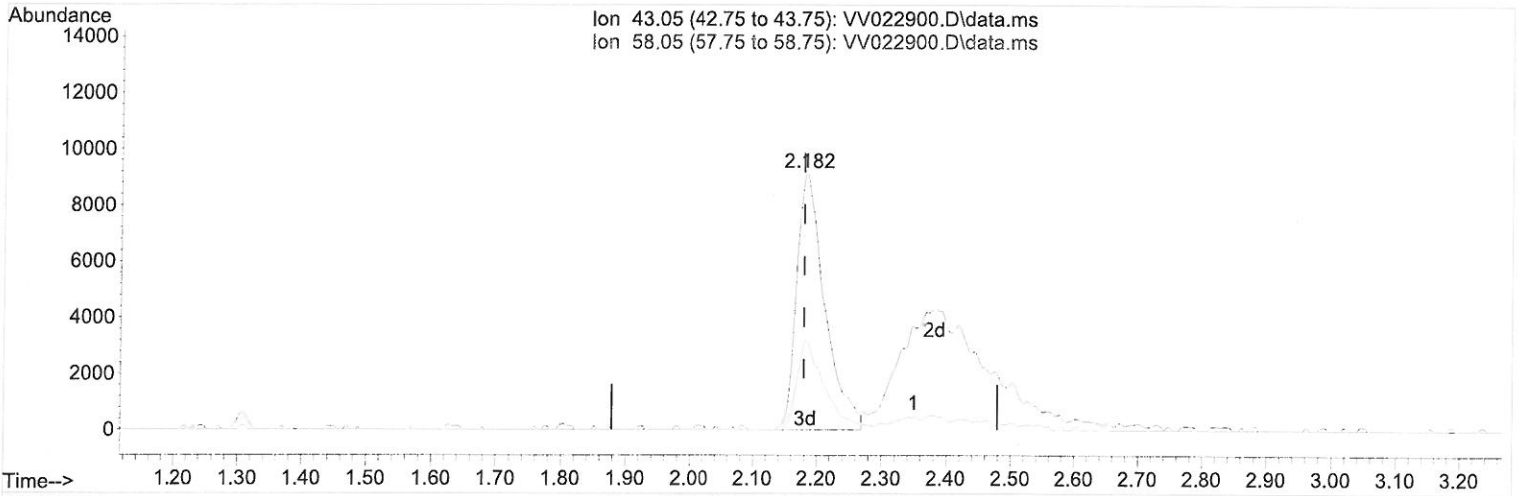
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 Quant Title : TRACE VOA SFAM1.0
 QLast Update : Tue Oct 19 04:07:06 2021
 Response via : Initial Calibration



TIC: VV022900.D\data.ms

(13) Acetone (T)

2.182min (+ 0.003) 23.37 ug/L m

response 28028

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	12.10	1.81
0.00	0.00	0.00
0.00	0.00	0.00

7 MD
 10/27/21

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VW101821\
 Data File : VW022900.D
 Acq On : 19 Oct 2021 05:05
 Operator : SY/MD
 Sample : M4137-15
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BFGD3

Manual Integrations
 APPROVED

MMDadoda
 10/19/2021 3:48:36 PM

Quant Time: Oct 19 09:00:24 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100721WMA.M
 Quant Title : TRACE VOA SFAM1.0
 QLast Update : Tue Oct 19 04:07:06 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Difluorobenzene	5.619	114	119231	5.000	ug/L	0.00
28) Chlorobenzene-d5	8.854	117	122414	5.000	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.252	152	57154	5.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.307	65	32007	3.902	ug/L	0.00
Spiked Amount 5.000	Range 40 - 130		Recovery =	78.000%		
7) Chloroethane-d5	1.568	69	28255	3.833	ug/L	0.00
Spiked Amount 5.000	Range 65 - 130		Recovery =	76.600%		
11) 1,1-Dichloroethene-d2	2.108	63	48110	2.990	ug/L	0.00
Spiked Amount 5.000	Range 60 - 125		Recovery =	59.800%#		
20) 2-Butanone-d5	3.889	46	113599	52.814	ug/L	0.00
Spiked Amount 50.000	Range 40 - 130		Recovery =	105.620%		
24) Chloroform-d	4.352	84	91756	5.428	ug/L	0.00
Spiked Amount 5.000	Range 70 - 125		Recovery =	108.600%		
26) 1,2-Dichloroethane-d4	5.037	65	42865	5.224	ug/L	0.00
Spiked Amount 5.000	Range 70 - 130		Recovery =	104.400%		
32) Benzene-d6	5.053	84	172189	4.725	ug/L	0.00
Spiked Amount 5.000	Range 70 - 125		Recovery =	94.600%		
36) 1,2-Dichloropropane-d6	6.072	67	56406	5.227	ug/L	0.00
Spiked Amount 5.000	Range 60 - 140		Recovery =	104.600%		
41) Toluene-d8	7.317	98	142965	4.538	ug/L	0.00
Spiked Amount 5.000	Range 70 - 130		Recovery =	90.800%		
43) trans-1,3-Dichloroprop...	7.625	79	15912	4.445	ug/L	0.00
Spiked Amount 5.000	Range 55 - 130		Recovery =	88.800%		
46) 2-Hexanone-d5	8.091	63	88039	61.022	ug/L	0.00
Spiked Amount 50.000	Range 45 - 130		Recovery =	122.040%		
56) 1,1,2,2-Tetrachloroeth...	10.217	84	42077	5.755	ug/L	0.00
Spiked Amount 5.000	Range 65 - 120		Recovery =	115.200%		
66) 1,2-Dichlorobenzene-d4	11.625	152	58546	5.525	ug/L	0.00
Spiked Amount 5.000	Range 80 - 120		Recovery =	110.400%		
Target Compounds						
13) Acetone	2.182	43	28028m	23.368	ug/L	Qvalue
42) Toluene	7.397	91	6397	0.191	ug/L	89
53) m,p-xylene	9.149	106	1070	0.080	ug/L	90

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 10/27/21

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