

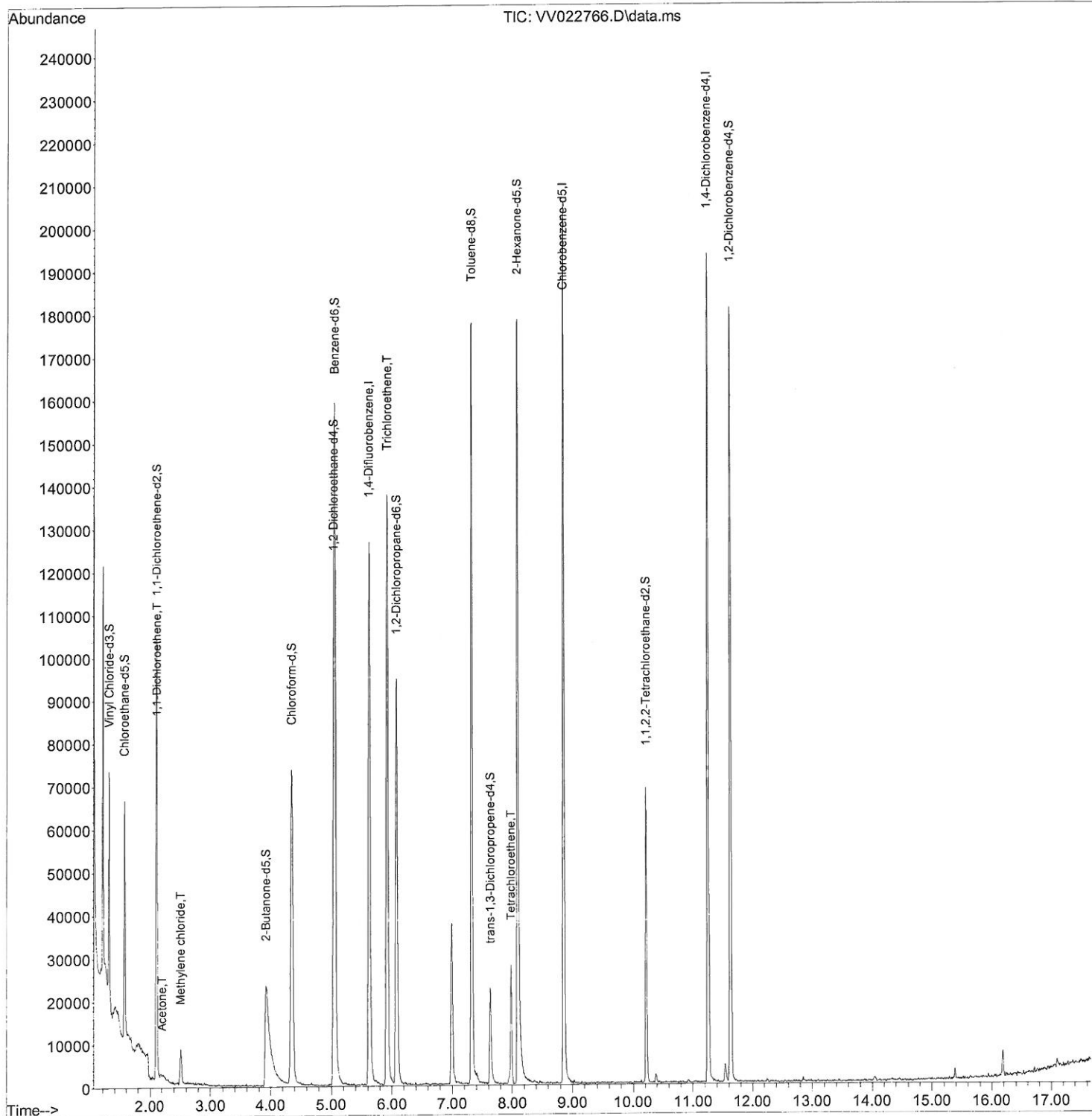
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101321\
Data File : VV022766.D
Acq On : 13 Oct 2021 17:02
Operator : SY/MD
Sample : M4109-17DL 4X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
YATK1DL

Quant Time: Oct 14 02:15:20 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100721WMA.M
Quant Title : TRACE VOA SFAM1.0
QLast Update : Thu Oct 14 02:12:50 2021
Response via : Initial Calibration

Manual Integrations
APPROVED

MMDadoda
10/14/2021 2:06:47 PM



Quantitation Report (Qedit)

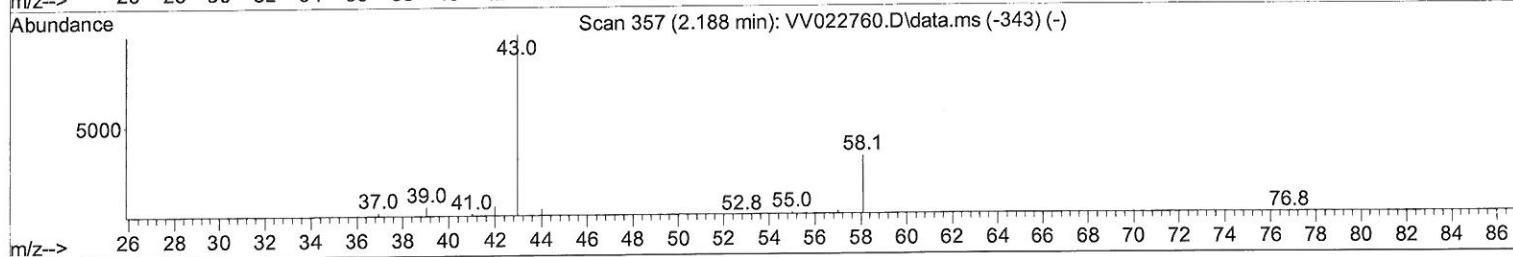
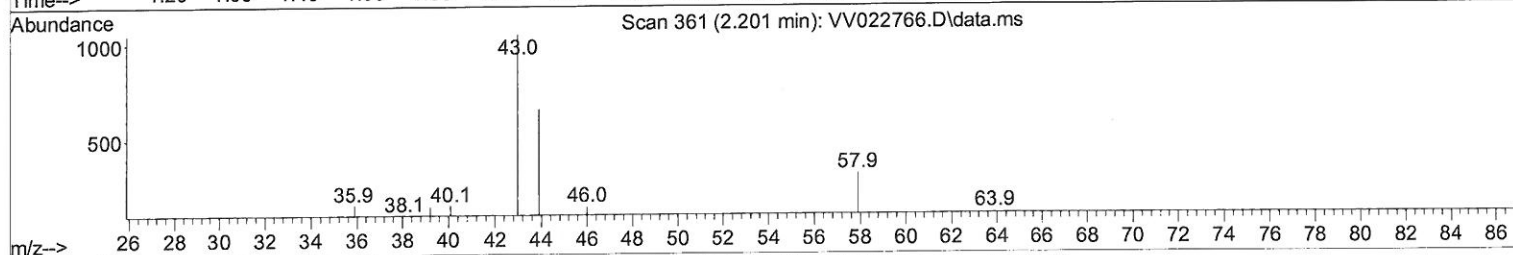
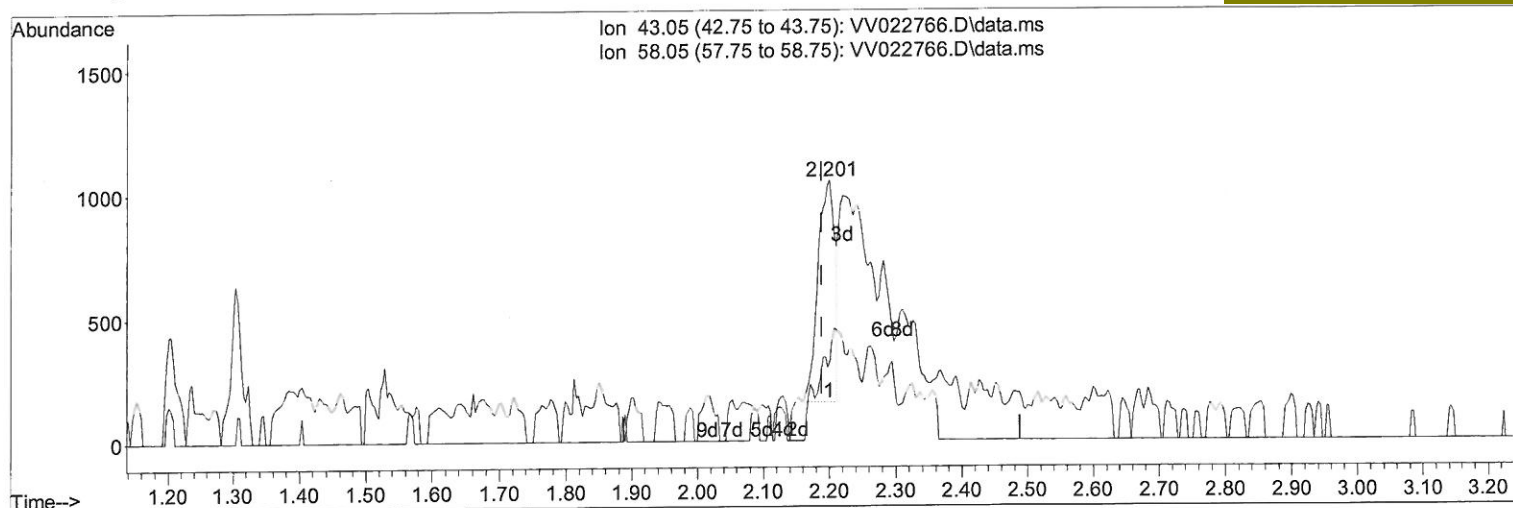
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TIC: VV022766.D\data.ms

(13) Acetone (T)

2.201min (+ 0.013) 1.35 ug/L

response 1551

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	12.10	17.15
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

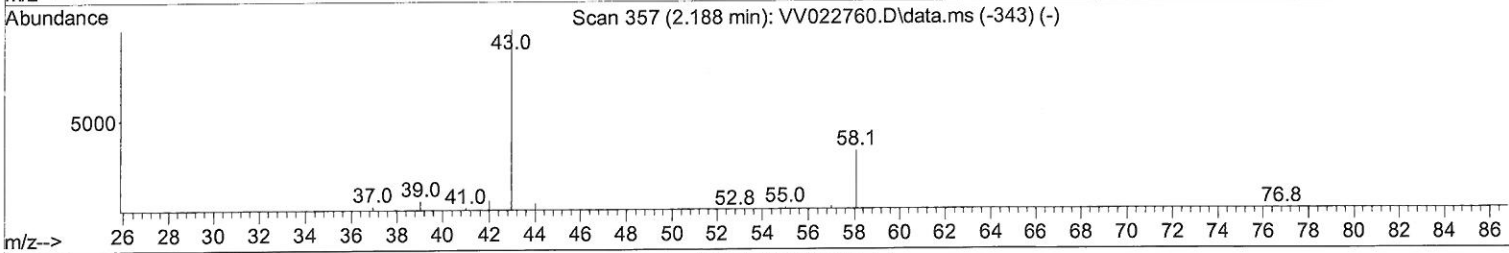
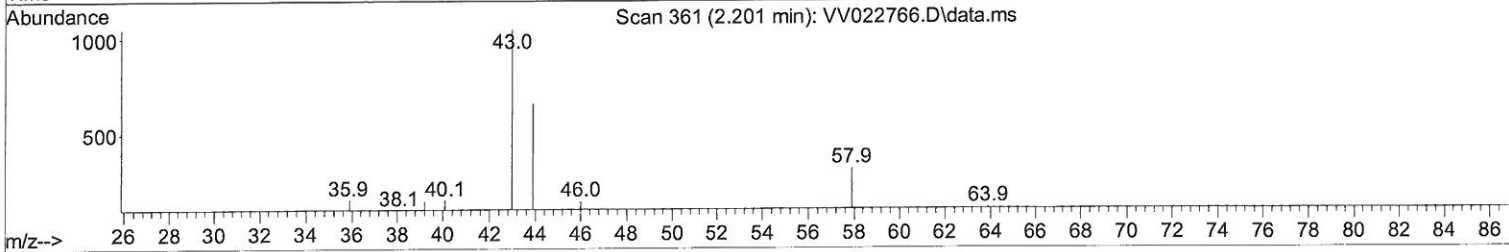
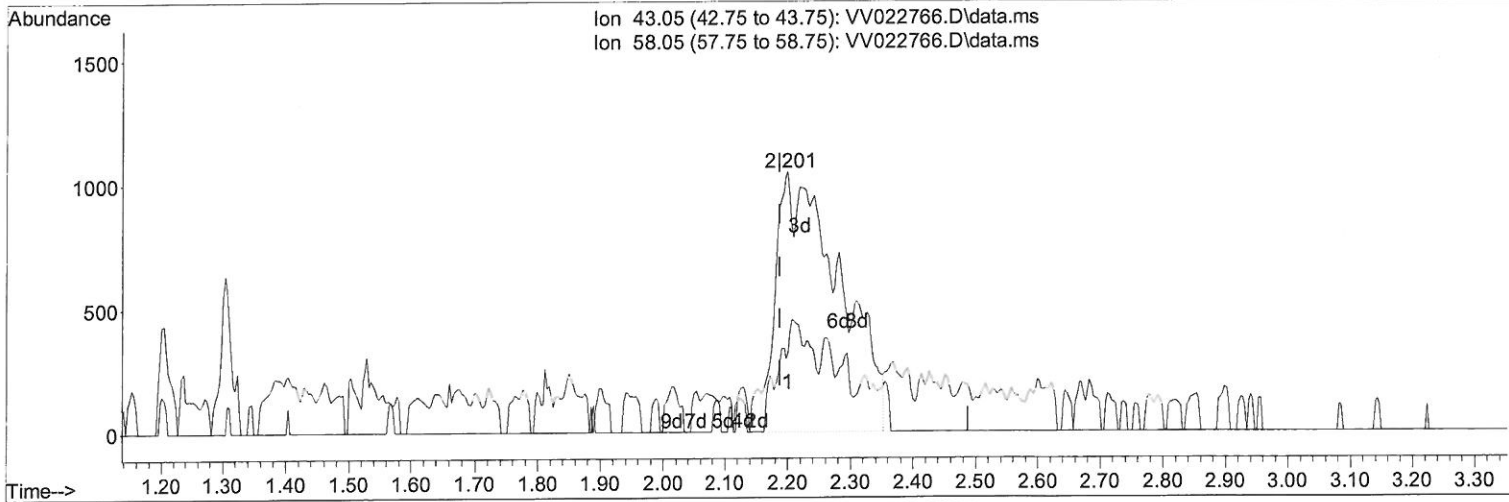
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Quant Time: Oct 14 02:15:20 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100721WMA.M
 Quant Title : TRACE VOA SFAM1.0
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Manual Integrations
 APPROVED

MMDadoda
 10/14/2021 2:06:47 PM



TIC: VV022766.D\data.ms

(13) Acetone (T)

2.201min (+ 0.013) 6.34 ug/L m

response 7304

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	12.10	3.64
0.00	0.00	0.00
0.00	0.00	0.00

Handwritten: 7 mg / 10/20/21

Quantitation Report (QT Reviewed)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VW101321\
 Data File : VW022766.D
 Acq On : 13 Oct 2021 17:02
 Operator : SY/MD
 Sample : M4109-17DL 4X
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_V
Client Sample ID :
 YATK1DL

Manual Integrations
APPROVED

MMDadoda
 10/14/2021 2:06:47 PM

Quant Time: Oct 14 02:15:20 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100721WMA.M
 Quant Title : TRACE VOA SFAM1.0
 QLast Update : Thu Oct 14 02:12:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Difluorobenzene	5.619	114	114602	5.000	ug/L	0.00
28) Chlorobenzene-d5	8.853	117	117859	5.000	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.249	152	54032	5.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.304	65	33259	4.218	ug/L	0.00
Spiked Amount	5.000	Range 40 - 130	Recovery =	84.400%		
7) Chloroethane-d5	1.568	69	34612	4.885	ug/L	0.00
Spiked Amount	5.000	Range 65 - 130	Recovery =	97.600%		
11) 1,1-Dichloroethene-d2	2.105	63	52851	3.417	ug/L	0.00
Spiked Amount	5.000	Range 60 - 125	Recovery =	68.400%		
20) 2-Butanone-d5	3.921	46	77108	37.297	ug/L	0.00
Spiked Amount	50.000	Range 40 - 130	Recovery =	74.600%		
24) Chloroform-d	4.352	84	77863	4.792	ug/L	0.00
Spiked Amount	5.000	Range 70 - 125	Recovery =	95.800%		
26) 1,2-Dichloroethane-d4	5.037	65	35611	4.516	ug/L	0.00
Spiked Amount	5.000	Range 70 - 130	Recovery =	90.400%		
32) Benzene-d6	5.053	84	143854	4.100	ug/L	0.00
Spiked Amount	5.000	Range 70 - 125	Recovery =	82.000%		
36) 1,2-Dichloropropane-d6	6.072	67	46748	4.499	ug/L	0.00
Spiked Amount	5.000	Range 60 - 140	Recovery =	90.000%		
41) Toluene-d8	7.317	98	119388	3.936	ug/L	0.00
Spiked Amount	5.000	Range 70 - 130	Recovery =	78.800%		
43) trans-1,3-Dichloroprop...	7.628	79	13964	4.051	ug/L	0.00
Spiked Amount	5.000	Range 55 - 130	Recovery =	81.000%		
46) 2-Hexanone-d5	8.095	63	71279	51.315	ug/L	0.00
Spiked Amount	50.000	Range 45 - 130	Recovery =	102.620%		
56) 1,1,2,2-Tetrachloroeth...	10.220	84	33378	4.742	ug/L	0.00
Spiked Amount	5.000	Range 65 - 120	Recovery =	94.800%		
66) 1,2-Dichlorobenzene-d4	11.625	152	48668	4.858	ug/L	0.00
Spiked Amount	5.000	Range 80 - 120	Recovery =	97.200%		
Target Compounds						
12) 1,1-Dichloroethene	2.114	96	1685	0.243	ug/L #	1
13) Acetone	2.201	43	7304m	6.336	ug/L	
16) Methylene chloride	2.510	84	3582	0.376	ug/L	88
34) Trichloroethene	5.918	95	48541	5.623	ug/L	95
47) Tetrachloroethene	7.979	164	6615	0.948	ug/L	93

MD
10/20/21

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