

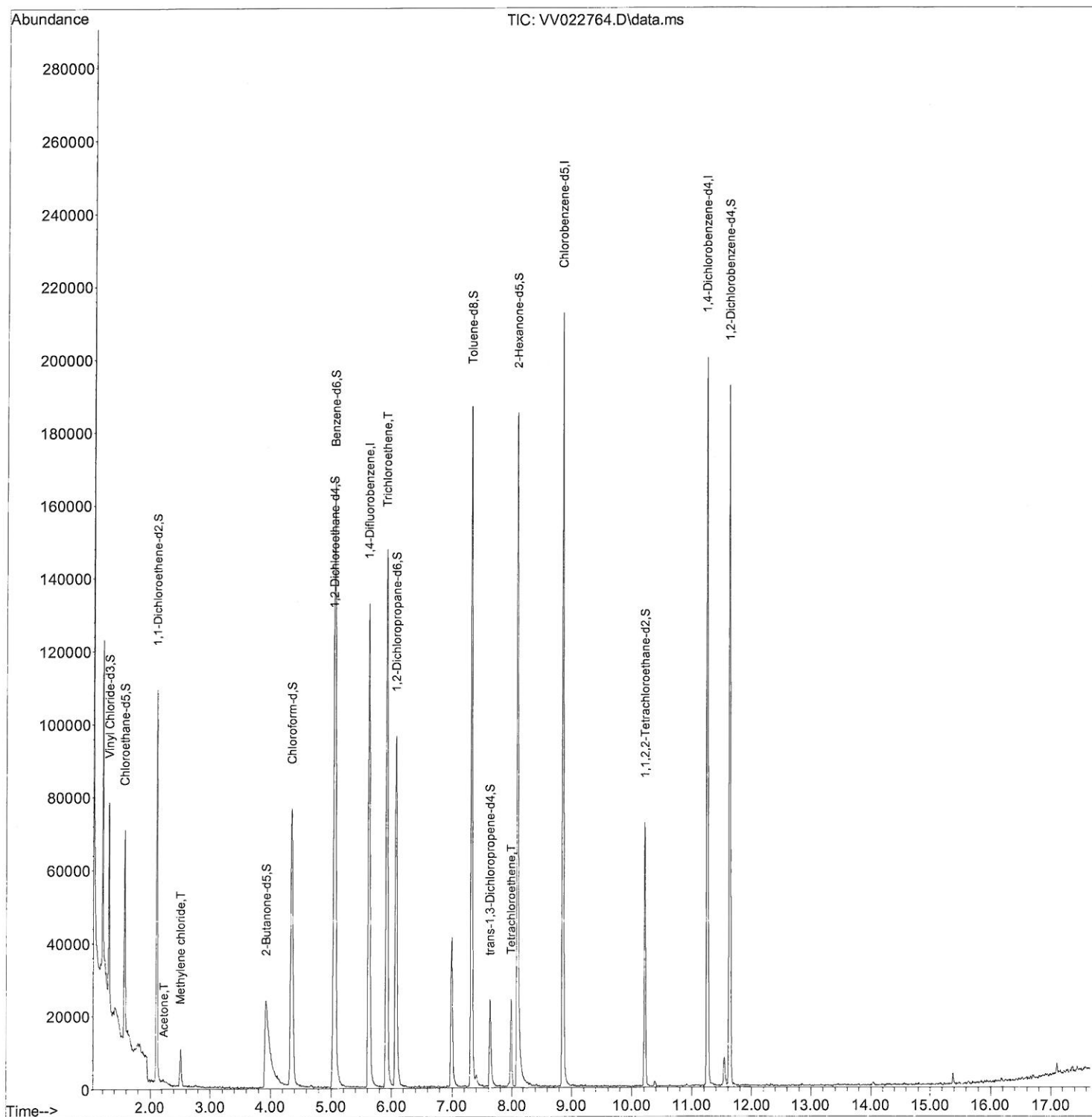
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101321\
Data File : VV022764.D
Acq On : 13 Oct 2021 16:14
Operator : SY/MD
Sample : M4109-10DL 5X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
YATJ6DL

Quant Time: Oct 14 02:14:54 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 2:06:43 PM



Quantitation Report (Qedit)

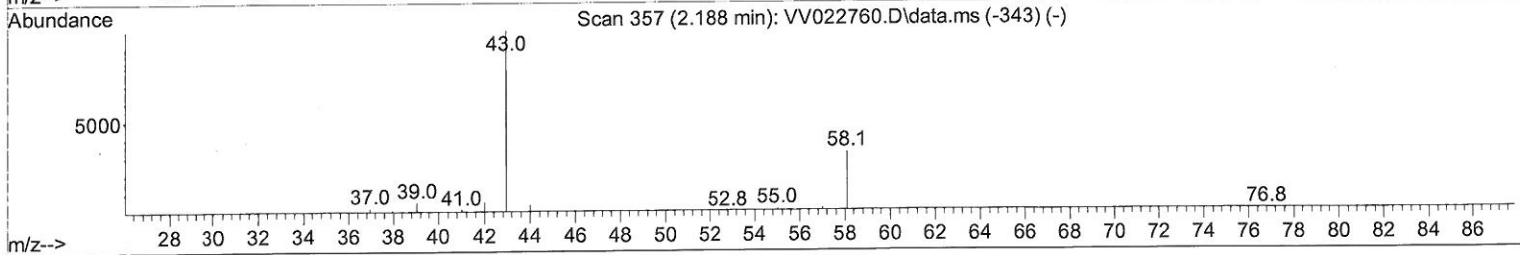
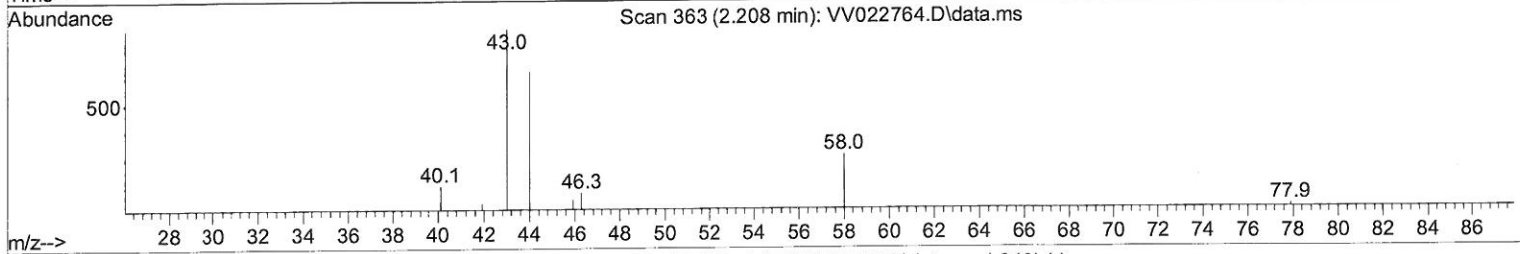
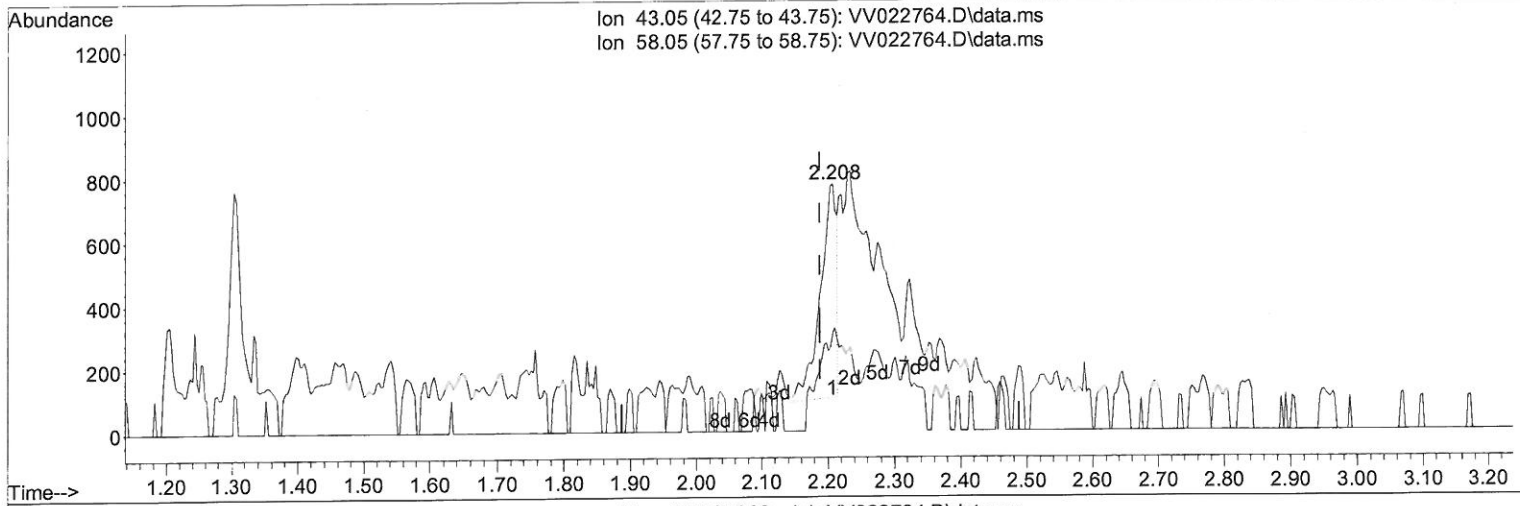
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TIC: VV022764.D\data.ms

(13) Acetone (T)

2.208min (+ 0.019) 0.97 ug/L

response 1142

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	12.10	11.30
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

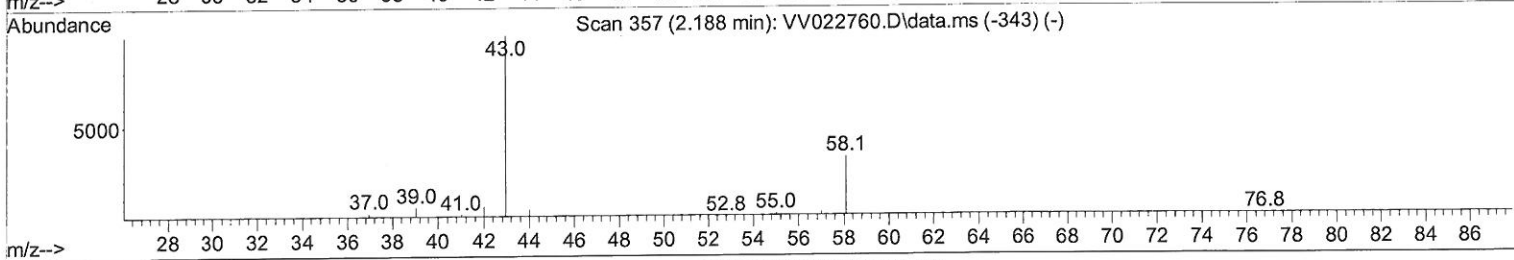
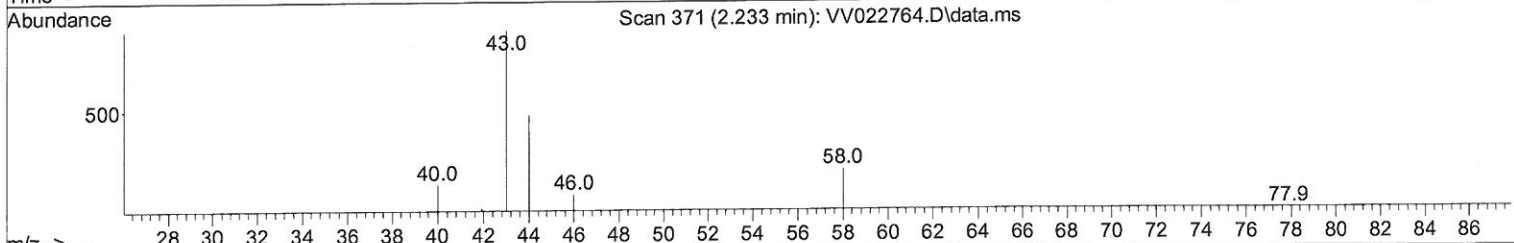
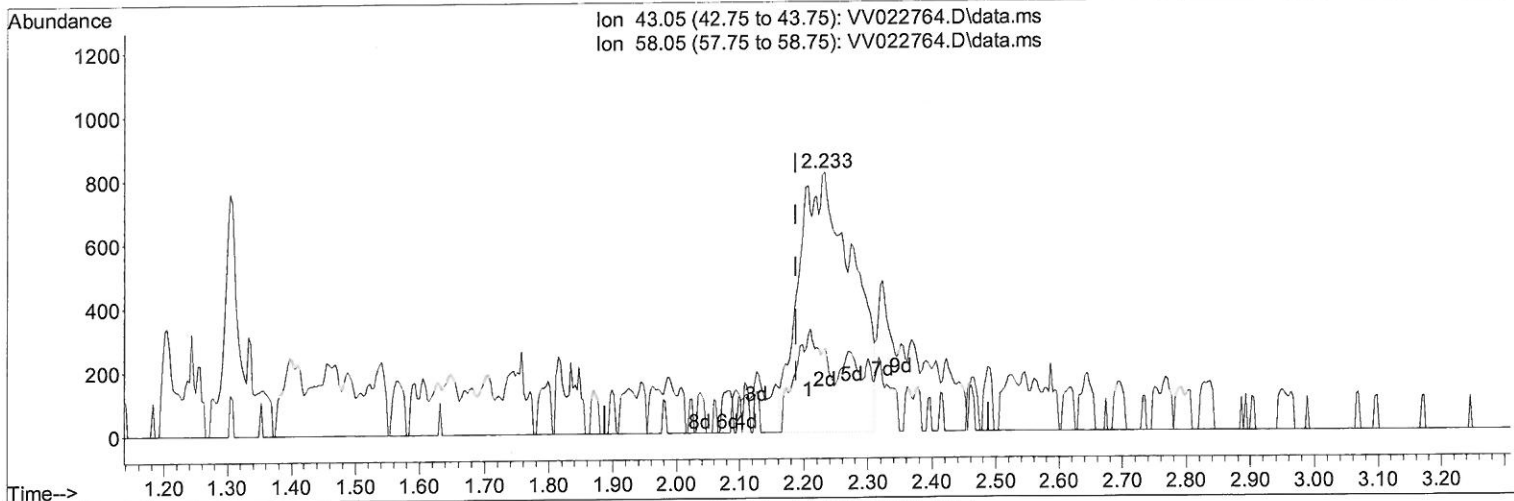
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Manual Integrations
 APPROVED

MMDadoda
 10/14/2021 2:06:43 PM



TIC: VV022764.D\data.ms

(13) Acetone (T)

2.233min (+ 0.045) 4.02 ug/L m 7 MD
 10/20/21

response 4742

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	12.10	2.72
0.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VW101321\
 Data File : VW022764.D
 Acq On : 13 Oct 2021 16:14
 Operator : SY/MD
 Sample : M4109-10DL 5X
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_V
 Client Sample ID :
 YATJ6DL

Quant Time: Oct 14 02:14:54 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100721WMA.M
 Quant Title : TRACE VOA SFAM1.0
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Manual Integrations
 APPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Difluorobenzene	5.619	114	117218	5.000	ug/L	0.00
28) Chlorobenzene-d5	8.854	117	122529	5.000	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.252	152	55151	5.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.304	65	34477	4.275	ug/L	0.00
Spiked Amount	5.000	Range 40 - 130	Recovery	=	85.600%	
7) Chloroethane-d5	1.568	69	34693	4.787	ug/L	0.00
Spiked Amount	5.000	Range 65 - 130	Recovery	=	95.800%	
11) 1,1-Dichloroethene-d2	2.108	63	54550	3.448	ug/L	0.00
Spiked Amount	5.000	Range 60 - 125	Recovery	=	69.000%	
20) 2-Butanone-d5	3.928	46	82547	39.036	ug/L	0.01
Spiked Amount	50.000	Range 40 - 130	Recovery	=	78.080%	
24) Chloroform-d	4.352	84	81022	4.875	ug/L	0.00
Spiked Amount	5.000	Range 70 - 125	Recovery	=	97.600%	
26) 1,2-Dichloroethane-d4	5.037	65	36619	4.540	ug/L	0.00
Spiked Amount	5.000	Range 70 - 130	Recovery	=	90.800%	
32) Benzene-d6	5.053	84	150948	4.138	ug/L	0.00
Spiked Amount	5.000	Range 70 - 125	Recovery	=	82.800%	
36) 1,2-Dichloropropane-d6	6.072	67	48953	4.532	ug/L	0.00
Spiked Amount	5.000	Range 60 - 140	Recovery	=	90.600%	
41) Toluene-d8	7.320	98	127074	4.030	ug/L	0.00
Spiked Amount	5.000	Range 70 - 130	Recovery	=	80.600%	
43) trans-1,3-Dichloroprop...	7.625	79	14808	4.132	ug/L	0.00
Spiked Amount	5.000	Range 55 - 130	Recovery	=	82.600%	
46) 2-Hexanone-d5	8.095	63	74797	51.795	ug/L	0.00
Spiked Amount	50.000	Range 45 - 130	Recovery	=	103.580%	
56) 1,1,2,2-Tetrachloroeth...	10.220	84	34422	4.704	ug/L	0.00
Spiked Amount	5.000	Range 65 - 120	Recovery	=	94.000%	
66) 1,2-Dichlorobenzene-d4	11.625	152	51912	5.077	ug/L	0.00
Spiked Amount	5.000	Range 80 - 120	Recovery	=	101.600%	
Target Compounds						
13) Acetone	2.233	43	4742m	4.022	ug/L	Qvalue
16) Methylene chloride	2.510	84	4458	0.457	ug/L	87
34) Trichloroethene	5.915	95	51531	5.742	ug/L	99
47) Tetrachloroethene	7.979	164	5365	0.739	ug/L	97

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