

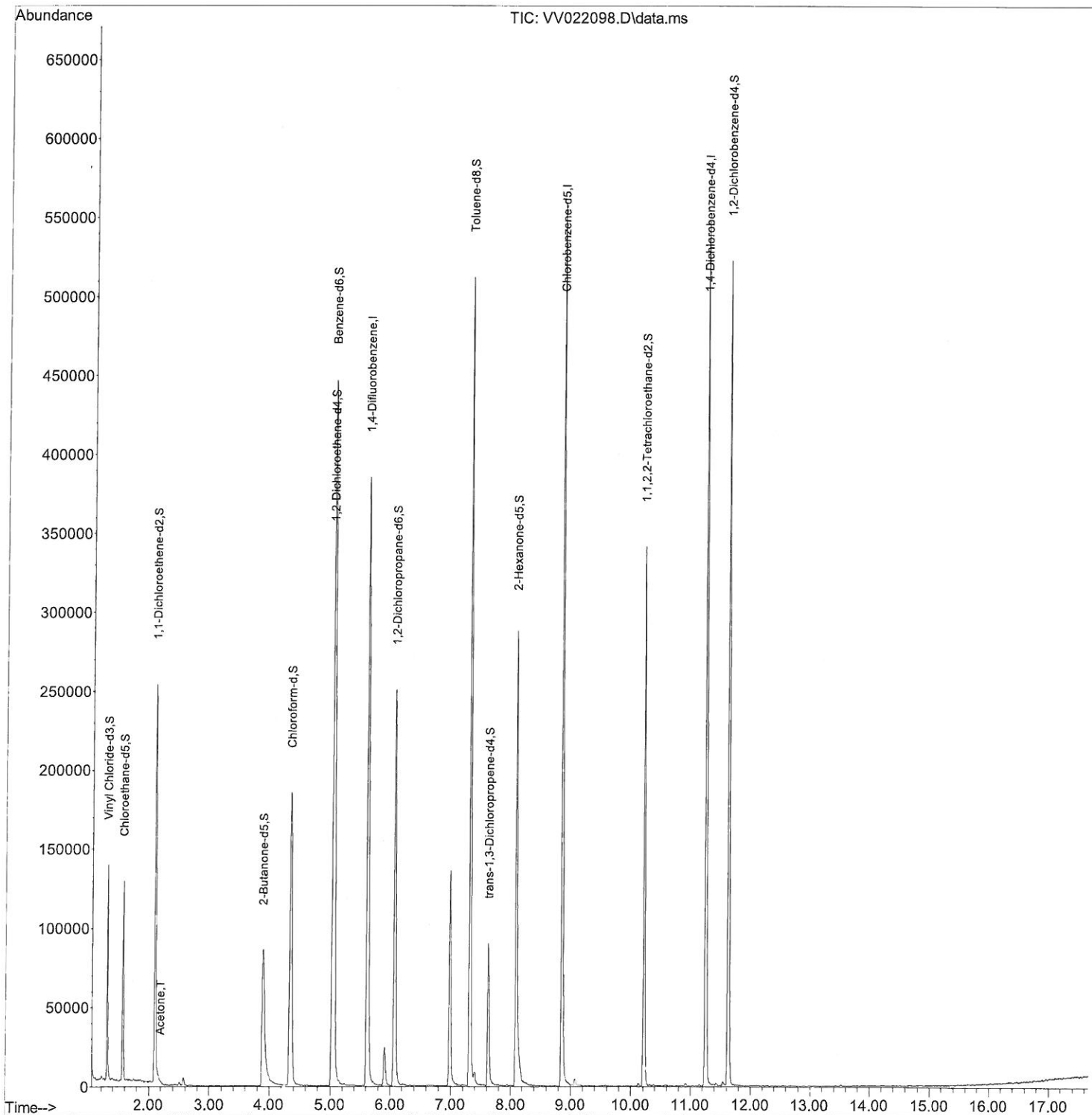
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090721\
Data File : VV022098.D
Acq On : 07 Sep 2021 13:23
Operator : SY/MD
Sample : M3651-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKL9

Manual Integrations
APPROVED

MMDadoda
9/8/2021 10:17:07 AM

Quant Time: Sep 08 01:16:27 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
Quant Title : VOC Analysis
QLast Update : Wed Sep 08 01:14:49 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

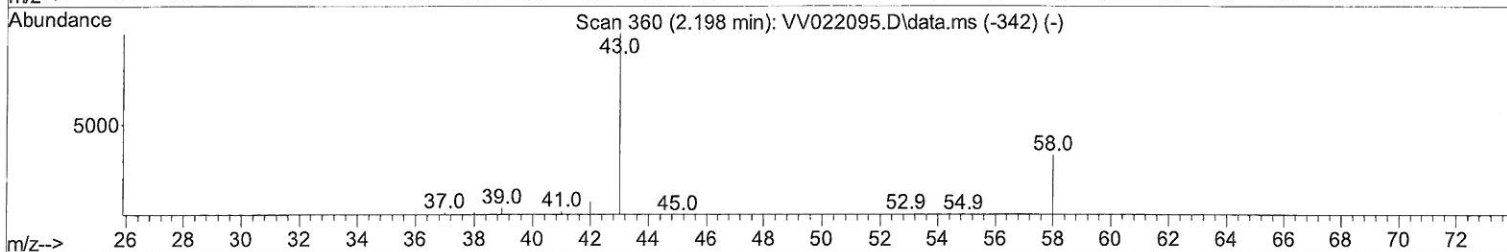
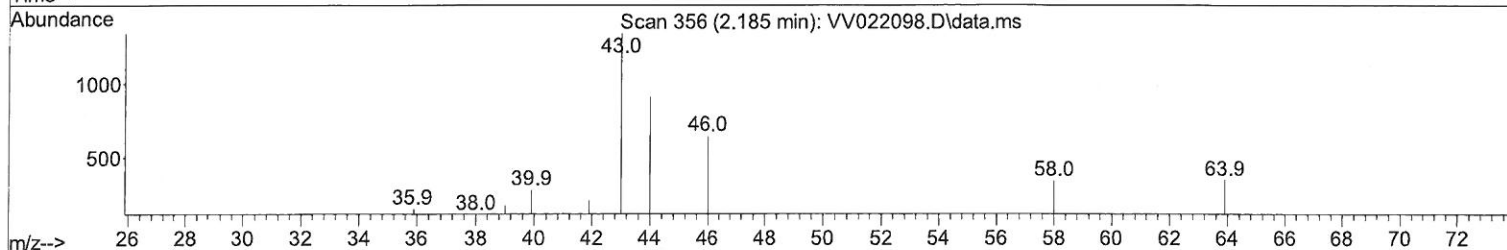
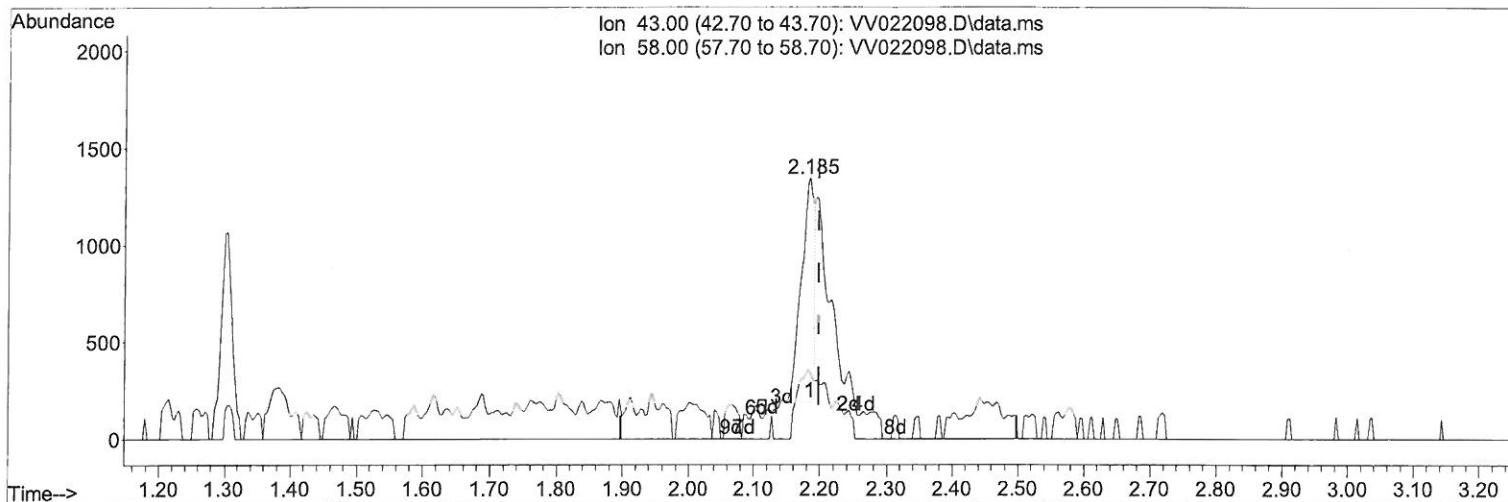
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TIC: VV022098.D\data.ms

(13) Acetone (T)

2.185min (-0.013) 1.15 ug/L

response 1582

Ion	Exp%	Act%
43.00	100.00	100.00
58.00	0.20	48.55#
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

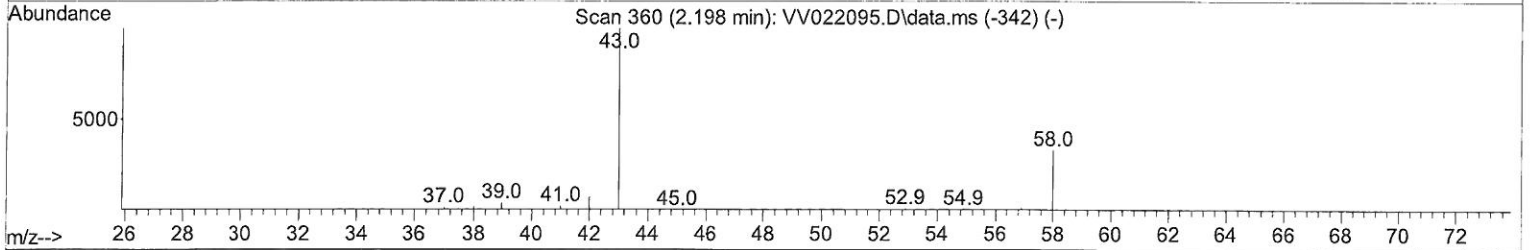
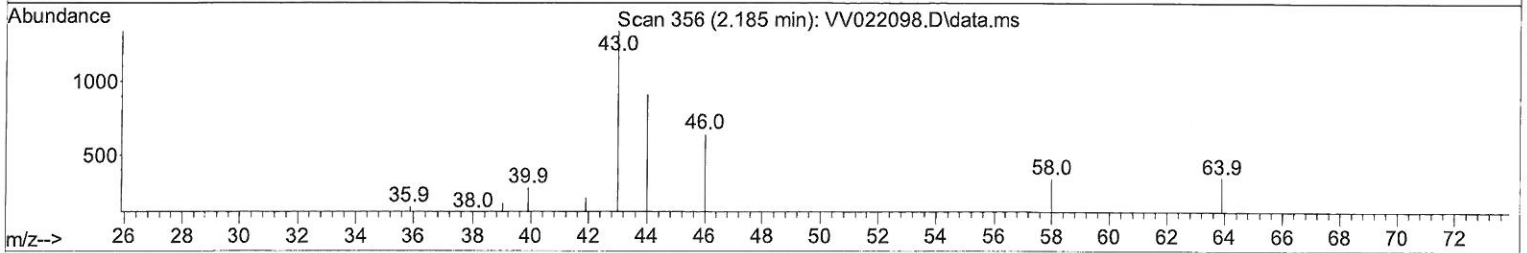
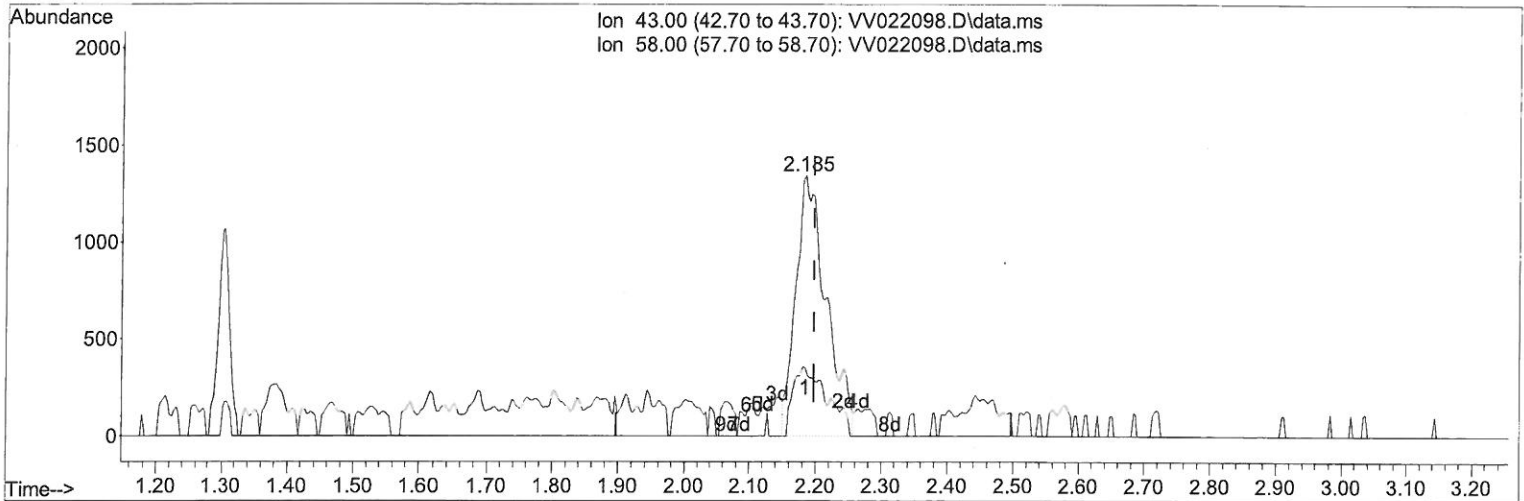
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Manual Integrations
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 QLast Update : Wed Sep 08 01:14:49 2021
 Response via : Initial Calibration



TIC: VV022098.D\data.ms

(13) Acetone (T)

2.185min (-0.013) 3.13 ug/L m

response 4290

Ion	Exp%	Act%
43.00	100.00	100.00
58.00	0.20	17.90#
0.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV090721\
 Data File : VV022098.D
 Acq On : 07 Sep 2021 13:23
 Operator : SY/MD
 Sample : M3651-03
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 DBKL9

Manual Integrations
 APPROVED

MMDadoda
 9/8/2021 10:17:07 AM

Quant Time: Sep 08 01:16:27 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082721WMA.M
 Quant Title : VOC Analysis
 QLast Update : Wed Sep 08 01:14:49 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Difluorobenzene	5.616	114	344802	50.000	ug/L	0.00
28) Chlorobenzene-d5	8.854	117	324529	50.000	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.249	152	149854	50.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.301	65	86852	36.373	ug/L	0.00
Spiked Amount	50.000	Range 60 - 135	Recovery	=	72.740%	
7) Chloroethane-d5	1.564	69	75981	40.747	ug/L	0.00
Spiked Amount	50.000	Range 70 - 130	Recovery	=	81.500%	
11) 1,1-Dichloroethene-d2	2.101	63	129114	32.695	ug/L	0.00
Spiked Amount	50.000	Range 60 - 125	Recovery	=	65.400%	
21) 2-Butanone-d5	3.889	46	169684	103.618	ug/L	0.00
Spiked Amount	100.000	Range 40 - 130	Recovery	=	103.620%	
24) Chloroform-d	4.346	84	194317	44.630	ug/L	0.00
Spiked Amount	50.000	Range 70 - 125	Recovery	=	89.260%	
26) 1,2-Dichloroethane-d4	5.031	65	121414	47.343	ug/L	0.00
Spiked Amount	50.000	Range 70 - 125	Recovery	=	94.680%	
32) Benzene-d6	5.047	84	393425	44.696	ug/L	0.00
Spiked Amount	50.000	Range 70 - 125	Recovery	=	89.400%	
36) 1,2-Dichloropropane-d6	6.069	67	127692	46.128	ug/L	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	=	92.260%	
41) Toluene-d8	7.317	98	351477	42.300	ug/L	0.00
Spiked Amount	50.000	Range 80 - 120	Recovery	=	84.600%	
43) trans-1,3-Dichloroprop...	7.622	79	53233	40.403	ug/L	0.00
Spiked Amount	50.000	Range 60 - 125	Recovery	=	80.800%	
47) 2-Hexanone-d5	8.088	63	101965	95.998	ug/L	0.00
Spiked Amount	100.000	Range 45 - 130	Recovery	=	96.000%	
56) 1,1,2,2-Tetrachloroeth...	10.217	84	162863	45.938	ug/L	0.00
Spiked Amount	50.000	Range 65 - 120	Recovery	=	91.880%	
66) 1,2-Dichlorobenzene-d4	11.625	152	140712	47.764	ug/L	0.00
Spiked Amount	50.000	Range 80 - 120	Recovery	=	95.520%	
Target Compounds						
13) Acetone	2.185	43	4290m	3.127	ug/L	Qvalue

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