

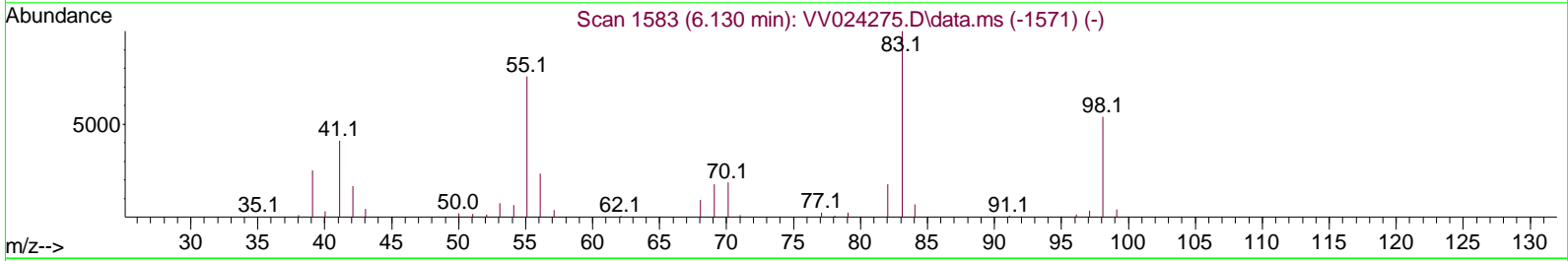
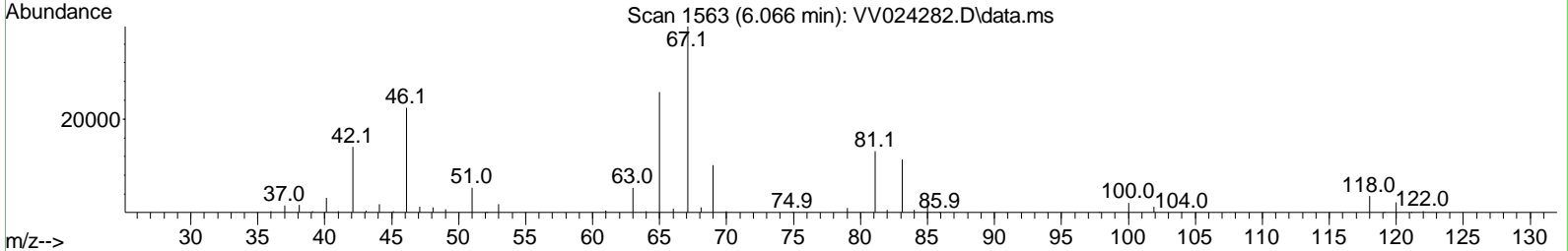
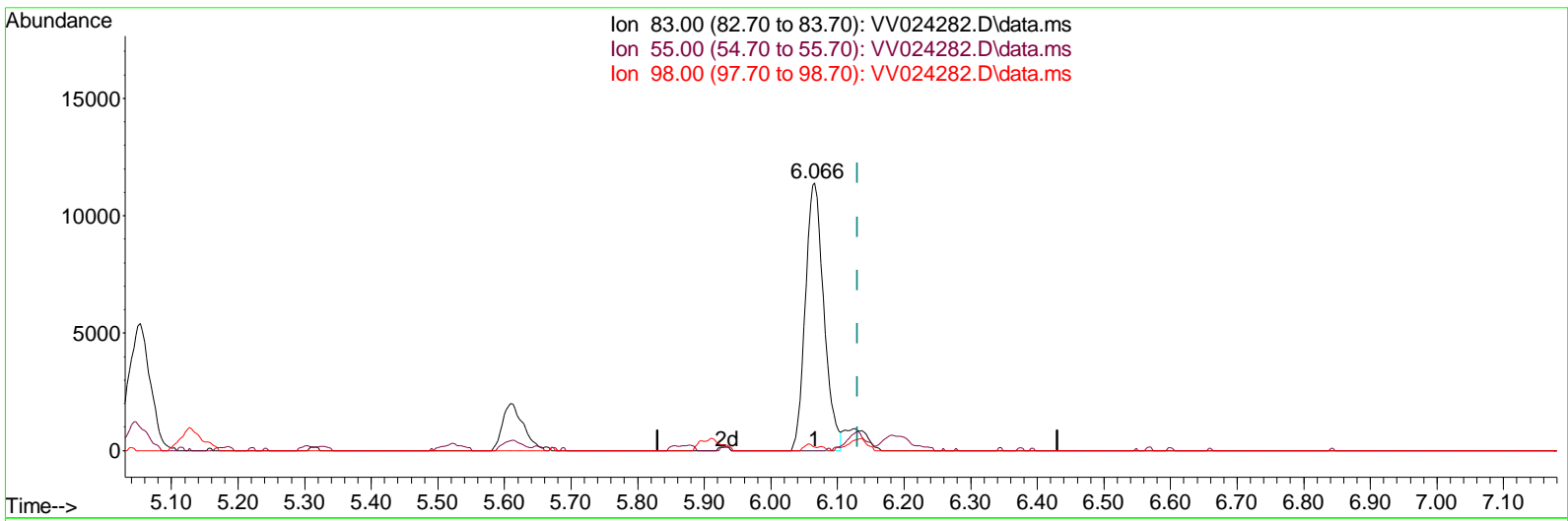
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
 Data File : VV024282.D
 Acq On : 05 Jan 2022 16:56
 Operator : SY/MD
 Sample : N1025-10
 Mi sc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_V
ClientSampleId :
 BGLH4

Manual IntegrationsAPPROVED

Reviewed By :John Carlone 01/06/2022
 Supervised By :Mahesh Dadoda 01/06/2022

Quant Time: Jan 06 01:01:22 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
 Quant Title : VOC Analysis
 QLast Update : Thu Jan 06 00:59:13 2022
 Response via : Initial Calibration



TIC: VV024282.D\data.ms

(35) Methylcyclohexane (T)

6.066min (-0.064) 8.95 ug/L

response	22111	
Ion	Exp%	Act%
83.00	100.00	100.00
55.00	69.50	0.18#
98.00	51.60	0.52#
0.00	0.00	0.00

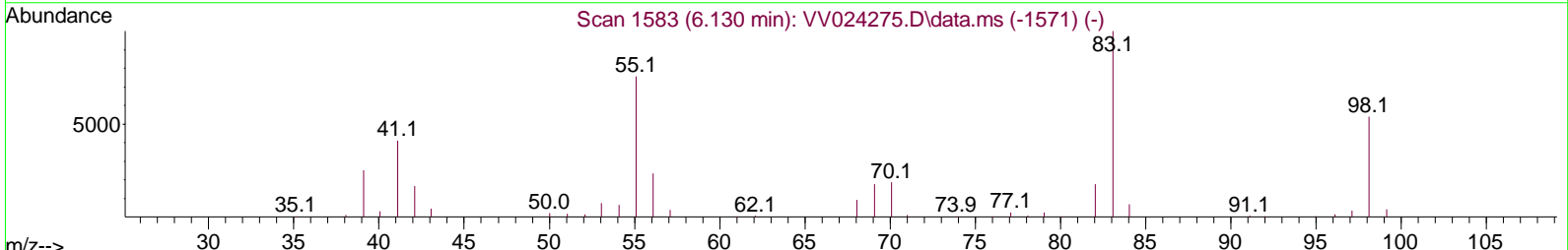
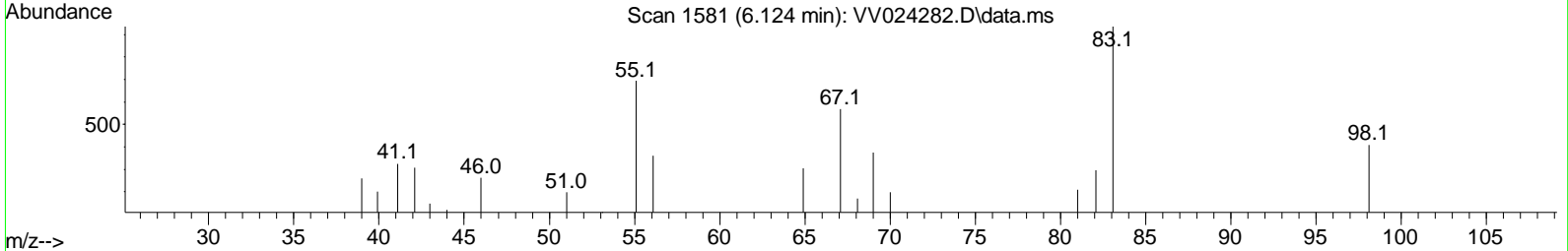
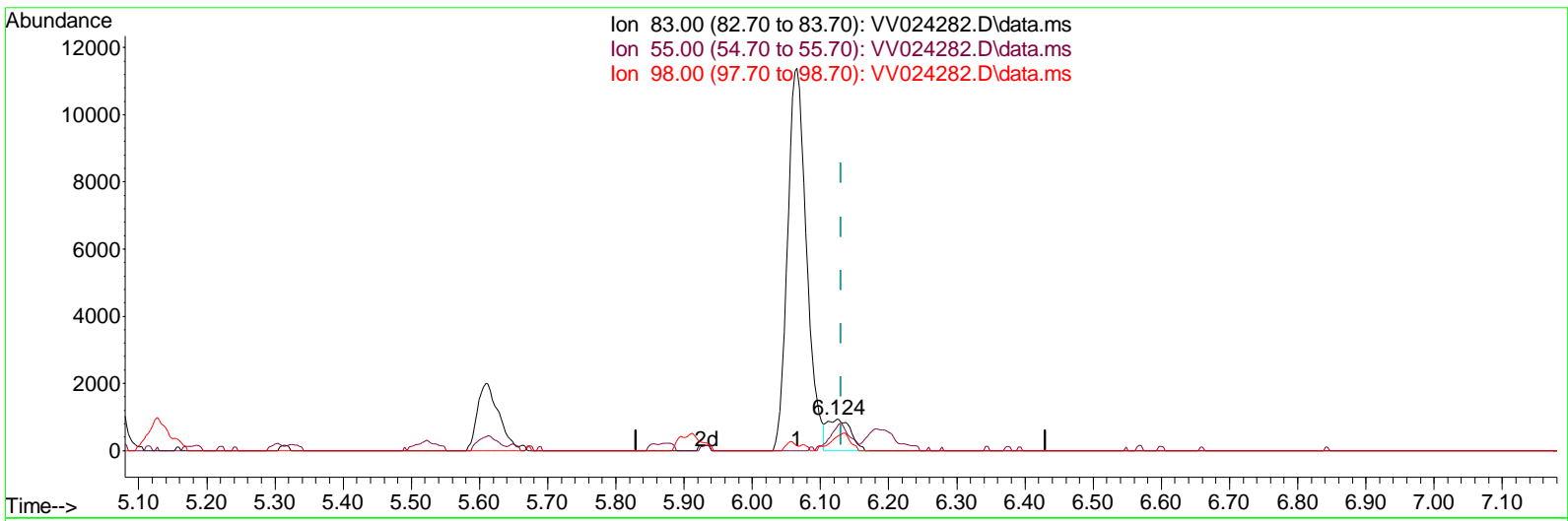
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TIC: VV024282.D\data.ms

(35) Methylcyclohexane (T)

6.124min (-0.006) 0.91 ug/L m

response	2239	
Ion	Exp%	Act%
83.00	100.00	100.00
55.00	69.50	1.79#
98.00	51.60	5.09#
0.00	0.00	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\WV010522\
 Data File : WV024282.D
 Acq On : 05 Jan 2022 16:56
 Operator : SY/MD
 Sample : N1025-10
 Mi sc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 9 Sample Multi plier: 1

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ClientSampleId :
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Compound	R. T.	QI on	Response	Conc	Units	Dev(Mi n)
Internal Standards						
1) 1,4-Di fluorobenzene	5.612	114	259159	50.000	ug/L	0.00
28) Chlorobenzene-d5	8.850	117	237469	50.000	ug/L	0.00
58) 1,4-Di chlorobenzene-d4	11.246	152	136899	50.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl chloride-d3	1.307	65	82767	44.376	ug/L	0.00
Spi ked Amount 50.000	Range 60 - 135		Recovery =	88.760%		
7) Chloroethane-d5	1.568	69	65152	45.983	ug/L	0.00
Spi ked Amount 50.000	Range 70 - 130		Recovery =	91.960%		
11) 1,1-Di chloroethene-d2	2.108	63	120621	34.603	ug/L	0.00
Spi ked Amount 50.000	Range 60 - 125		Recovery =	69.200%		
21) 2-Butanone-d5	3.876	46	101547	97.721	ug/L	0.00
Spi ked Amount 100.000	Range 40 - 130		Recovery =	97.720%		
24) Chloroform-d	4.342	84	151091	44.324	ug/L	0.00
Spi ked Amount 50.000	Range 70 - 125		Recovery =	88.640%		
26) 1,2-Di chloroethane-d4	5.027	65	104180	45.647	ug/L	0.00
Spi ked Amount 50.000	Range 70 - 125		Recovery =	91.300%		
32) Benzene-d6	5.047	84	294112	47.461	ug/L	0.00
Spi ked Amount 50.000	Range 70 - 125		Recovery =	94.920%		
36) 1,2-Di chloropropane-d6	6.066	67	81893	47.154	ug/L	0.00
Spi ked Amount 50.000	Range 70 - 120		Recovery =	94.300%		
41) Tol uene-d8	7.313	98	282726	47.121	ug/L	0.00
Spi ked Amount 50.000	Range 80 - 120		Recovery =	94.240%		
43) trans-1,3-Di chloroprop. . .	7.619	79	45728	45.726	ug/L	0.00
Spi ked Amount 50.000	Range 60 - 125		Recovery =	91.460%		
47) 2-Hexanone-d5	8.082	63	78877	104.538	ug/L	0.00
Spi ked Amount 100.000	Range 45 - 130		Recovery =	104.540%		
56) 1,1,2,2-Tetrachloroeth. . .	10.214	84	115924	46.708	ug/L	0.00
Spi ked Amount 50.000	Range 65 - 120		Recovery =	93.420%		
66) 1,2-Di chlorobenzene-d4	11.622	152	131123	48.053	ug/L	0.00
Spi ked Amount 50.000	Range 80 - 120		Recovery =	96.100%		
Target Compounds						
5) Vinyl chloride	1.314	62	65461	35.220	ug/L	98
12) 1,1-Di chloroethene	2.114	96	2240	1.419	ug/L #	1
13) Acetone	2.172	43	152683	148.817	ug/L	99
17) trans-1,2-Di chloroethene	2.764	96	5875	3.612	ug/L	92
19) 1,1-Di chloroethane	3.188	63	56897	21.088	ug/L	96
20) ci s-1,2-Di chloroethene	3.905	96	547727	315.280	ug/L	99
22) 2-Butanone	3.963	43	76185	69.904	ug/L	94
27) 1,2-Di chloroethane	5.127	62	24100	9.393	ug/L	96
29) Cycl ohexane	4.677	56	2235	1.053	ug/L #	95
33) Benzene	5.095	78	856521	141.850	ug/L	100
35) Methyl cycl ohexane	6.124	83	2239m	0.907	ug/L	
40) 4-Methyl -2-pentanone	7.230	43	12266	6.623	ug/L	97
42) Tol uene	7.384	91	2680104	395.409	ug/L	99
51) Chlorobenzene	8.879	112	65720	14.395	ug/L	99
52) Ethyl benzene	9.008	91	1630923	225.231	ug/L	100
53) m, p-Xyl ene	9.136	106	957296	329.818	ug/L	97
54) o-Xyl ene	9.542	106	566054	204.301	ug/L	98
60) I sopropyl benzene	9.931	105	216654	28.600	ug/L	99

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Compound	R. T.	QI on	Response	Conc	Units	Dev(Min)
62) 1,3,5-Tri methyl benzene	10.535	105	132135	20.278	ug/L	100
63) 1,2,4-Tri methyl benzene	10.911	105	1135087	174.483	ug/L	100
64) 1,3-Di chl orobenzene	11.181	146	3874	0.992	ug/L	94
65) 1,4-Di chl orobenzene	11.271	146	9720	2.413	ug/L	99
67) 1,2-Di chl orobenzene	11.641	146	93278	23.172	ug/L	98
70) 1,2,4-tri chl orobenzene	13.262	180	12263	4.773	ug/L	96
72) 1,2,3-Tri chl orobenzene	13.744	180	3249	1.223	ug/L #	92

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

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