

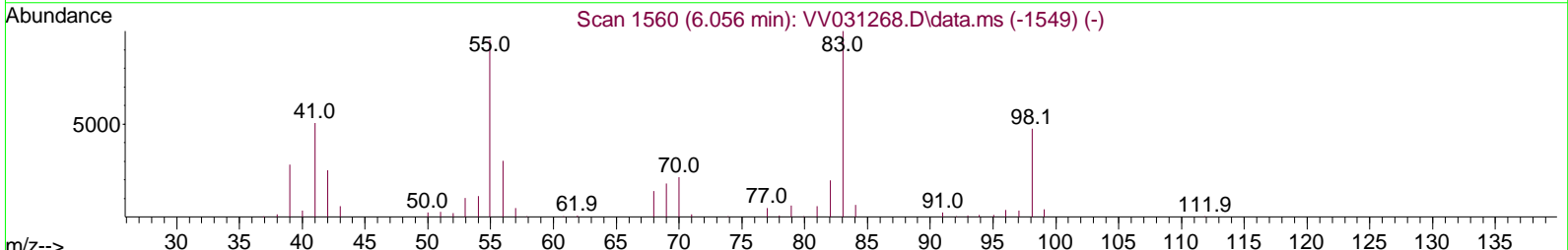
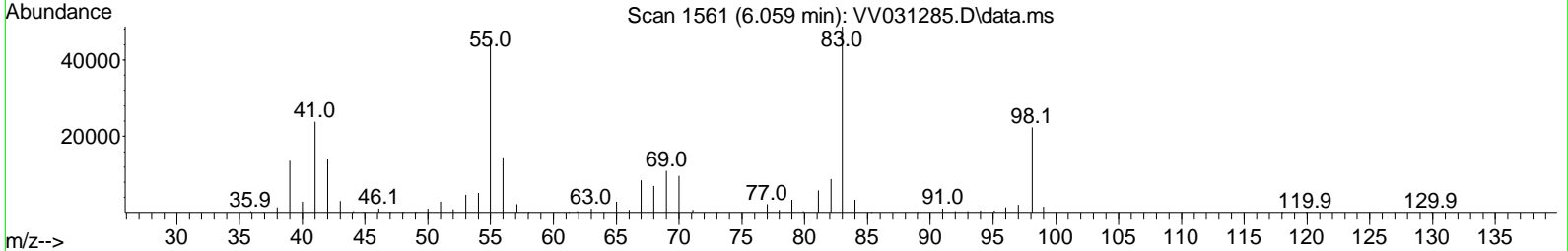
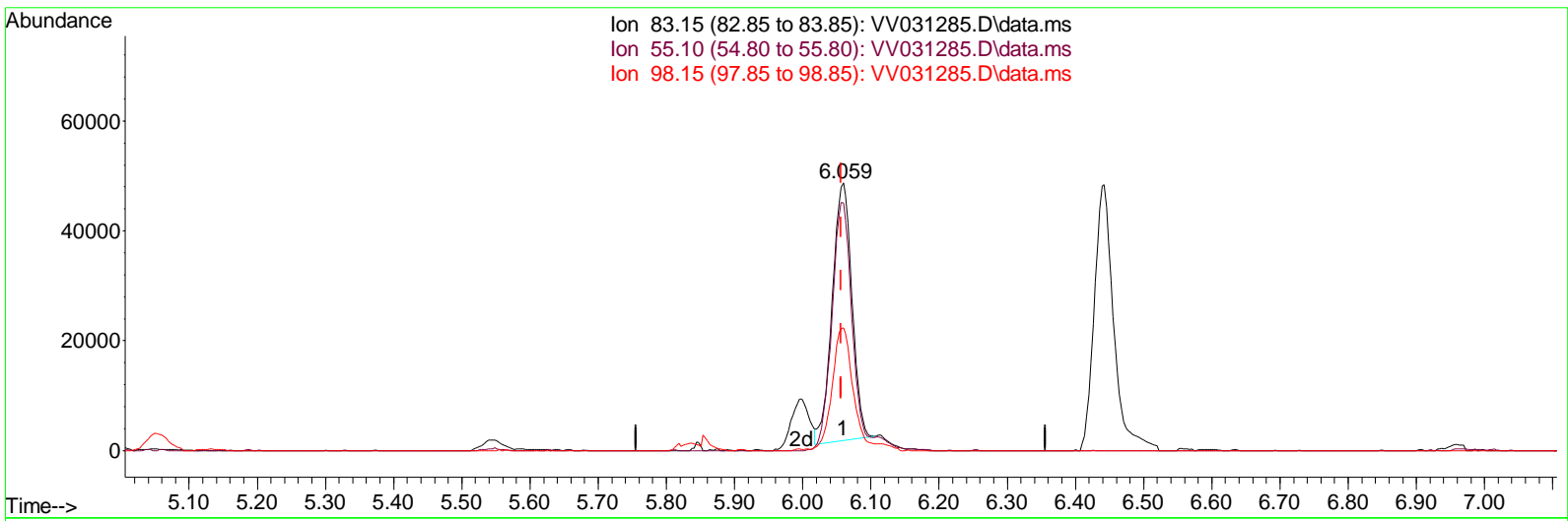
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV053023\
 Data File : VV031285.D
 Acq On : 30 May 2023 22: 30
 Operator : SY/MD
 Sample : VSTDCCC005EC
 Mi sc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_V
Lab Sampled :
 VSTDCCC005EC

Manual Integrations APPROVED

Reviewed By :Krupa Patel 05/31/2023
 Supervised By :Mahesh Dadoda 05/31/2023

Quant Time: May 31 05: 33: 11 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR053023WMA.M
 Quant Title : TRACE VOA SFAM1.0
 QLast Update : Wed May 31 05: 14: 27 2023
 Response via : Initial Calibration



TIC: VV031285.D\data.ms

(35) Methylcyclohexane (T)

6.059min (+ 0.003) 4.12 ug/L

response	94370	
Ion	Exp%	Act%
83.15	100.00	100.00
55.10	90.70	99.52
98.15	44.90	52.21
0.00	0.00	0.00

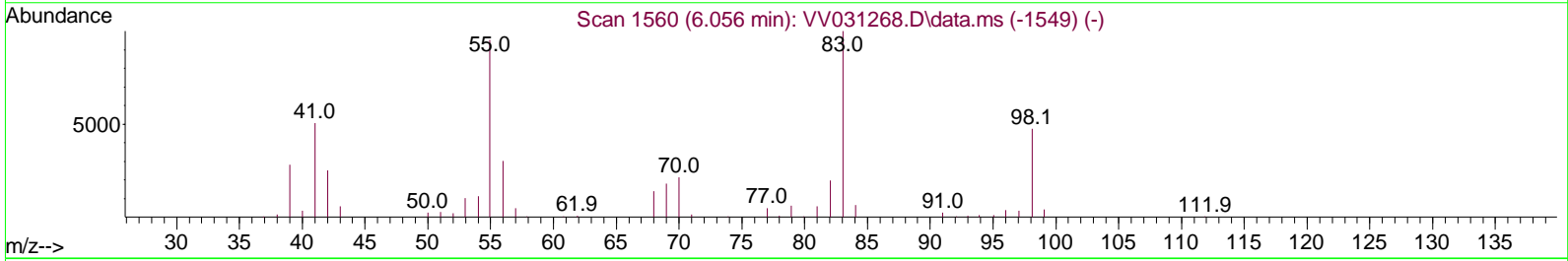
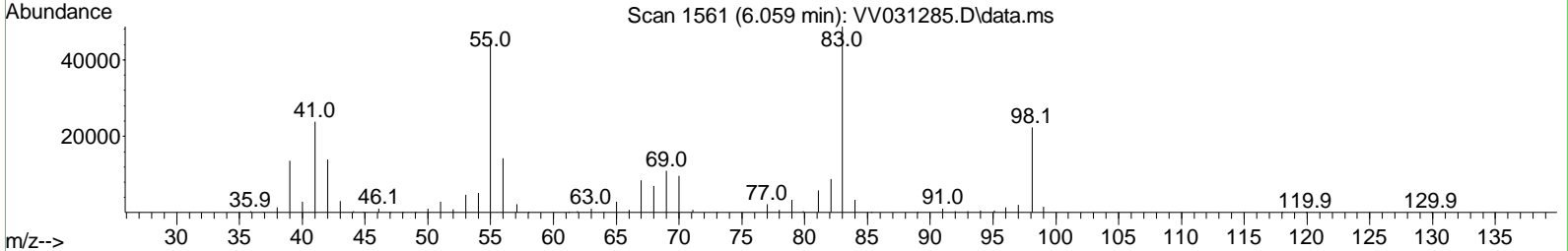
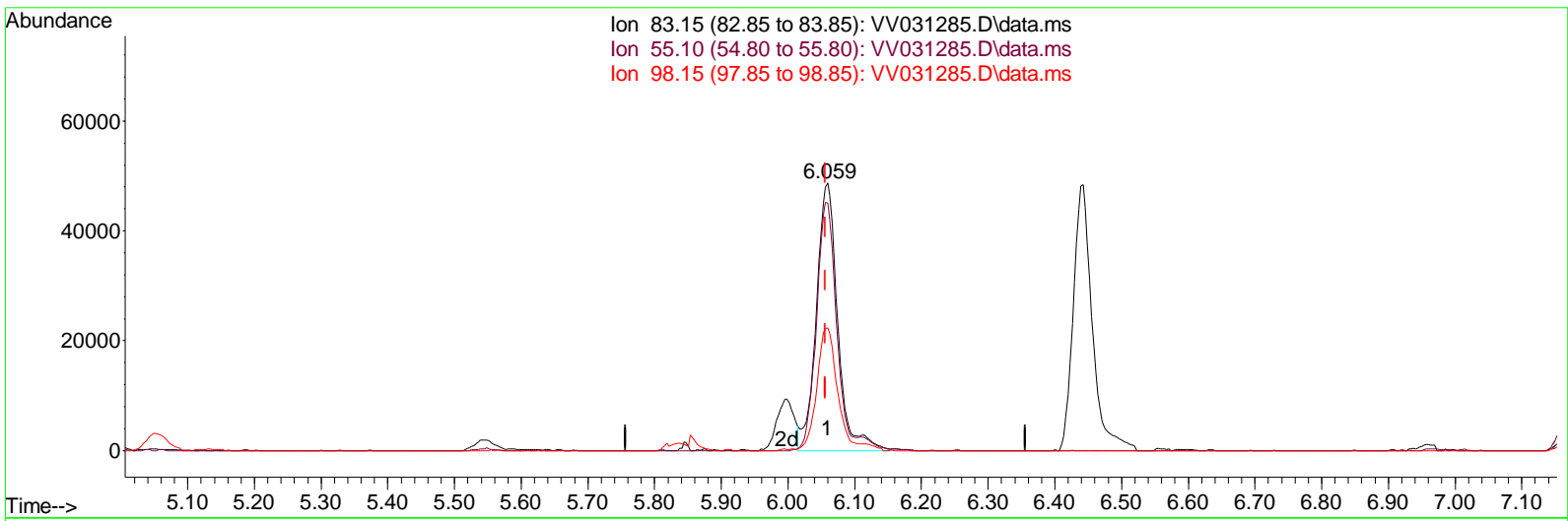
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TIC: VV031285.D\data.ms

(35) Methylcyclohexane (T)

6.059min (+ 0.003) 4.76 ug/L m

response 109032

Ion	Exp%	Act%
83.15	100.00	100.00
55.10	90.70	86.14
98.15	44.90	45.19
0.00	0.00	0.00

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 QLast Update : Wed May 31 05:14:27 2023
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Compound	R.T.	QI on	Response	Conc	Units	Dev(Mi n)
Internal Standards						
1) 1,4-Di fl uorobenzene	5.545	114	120530	5.000	ug/L	0.00
28) Chl orobenzene-d5	8.793	117	128268	5.000	ug/L	0.00
58) 1,4-Di chl orobenzene-d4	11.191	152	69417	5.000	ug/L	0.00
System Moni toring Compounds						
4) Vi nyl Chl ori de-d3	1.282	65	53020	5.310	ug/L	0.00
Spi ked Amount 5.000	Range 40	- 130	Recovery	= 106.200%		
7) Chl oroethane-d5	1.536	69	57617	5.413	ug/L	0.00
Spi ked Amount 5.000	Range 65	- 130	Recovery	= 108.200%		
11) 1,1-Di chl oroethene-d2	2.063	65	27826	5.339	ug/L	0.00
Spi ked Amount 5.000	Range 60	- 125	Recovery	= 106.800%		
20) 2-Butanone-d5	3.815	46	187404	54.334	ug/L	0.00
Spi ked Amount 50.000	Range 40	- 130	Recovery	= 108.660%		
24) Chl oroform-d	4.259	84	127347	5.311	ug/L	0.00
Spi ked Amount 5.000	Range 70	- 125	Recovery	= 106.200%		
26) 1,2-Di chl oroethane-d4	4.950	65	70494	5.500	ug/L	0.00
Spi ked Amount 5.000	Range 70	- 130	Recovery	= 110.000%		
32) Benzene-d6	4.966	84	230789	4.975	ug/L	0.00
Spi ked Amount 5.000	Range 70	- 125	Recovery	= 99.600%		
36) 1,2-Di chl oropropane-d6	5.995	67	81319	4.968	ug/L	0.00
Spi ked Amount 5.000	Range 60	- 140	Recovery	= 99.400%		
41) Tol uene-d8	7.249	98	220974	5.349	ug/L	0.00
Spi ked Amount 5.000	Range 70	- 130	Recovery	= 107.000%		
43) trans-1,3-Di chl oroprop. . .	7.561	79	27043	5.054	ug/L	0.00
Spi ked Amount 5.000	Range 55	- 130	Recovery	= 101.000%		
46) 2-Hexanone-d5	8.034	63	134784	53.331	ug/L	0.00
Spi ked Amount 50.000	Range 45	- 130	Recovery	= 106.660%		
56) 1,1,2,2-Tetrachl oroeth. . .	10.159	84	66471	5.295	ug/L	0.00
Spi ked Amount 5.000	Range 65	- 120	Recovery	= 105.800%		
66) 1,2-Di chl orobenzene-d4	11.567	152	72430	5.021	ug/L	0.00
Spi ked Amount 5.000	Range 80	- 120	Recovery	= 100.400%		
Target Compounds						
2) Di chl orodi fl uoromethane	1.108	85	98022	5.066	ug/L	98
3) Chl oromethane	1.217	50	92281	5.028	ug/L	97
5) Vi nyl chl ori de	1.285	62	93134	5.069	ug/L	99
6) Bromomethane	1.491	94	52564	5.086	ug/L	96
8) Chl oroethane	1.552	64	57096	4.956	ug/L	99
9) Tri chl orofl uoromethane	1.716	101	131732	4.748	ug/L	100
10) 1,1,2-Tri chl oro-1,2,2-. . .	2.073	101	72910	4.727	ug/L	100
12) 1,1-Di chl oroethene	2.073	96	63261	4.845	ug/L	97
13) Acetone	2.143	43	116641	41.165	ug/L	95
14) Carbon di sul fi de	2.246	76	188517	4.781	ug/L	100
15) Methyl Acetate	2.388	43	27949	5.215	ug/L	97
16) Methyl ene chl ori de	2.452	84	71199	4.013	ug/L	99
17) Methyl tert-butyl Ether	2.712	73	137663	4.865	ug/L	95
18) trans-1,2-Di chl oroethene	2.700	96	59235	4.770	ug/L	97
19) 1,1-Di chl oroethane	3.117	63	136681	4.972	ug/L	97
21) 2-Butanone	3.896	43	176735	48.091	ug/L	99
22) ci s-1,2-Di chl oroethene	3.822	96	63405	4.864	ug/L	100
23) Bromochl oromethane	4.153	128	28647	4.963	ug/L	92

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Compound	R. T.	QI on	Response	Conc	Units	Dev(Min)
25) Chloroform	4.285	83	144851	5.016	ug/L	97
27) 1,2-Dichloroethane	5.050	62	84238	5.303	ug/L	97
29) 1,1,1-Trichloroethane	4.519	97	119669	4.634	ug/L	100
30) Cyclohexane	4.593	56	107599	4.447	ug/L	97
31) Carbon tetrachloride	4.744	117	107760	4.655	ug/L	97
33) Benzene	5.018	78	269697	4.765	ug/L	100
34) Trichloroethene	5.841	95	68786	4.430	ug/L	95
35) Methylcyclohexane	6.059	83	109032m	4.761	ug/L	
37) 1,2-Dichloropropane	6.101	63	79764	5.205	ug/L	99
38) Bromodichloromethane	6.442	83	95995	4.744	ug/L #	98
39) cis-1,3-Dichloropropene	6.960	75	95536	4.671	ug/L	95
40) 4-Methyl-2-pentanone	7.166	43	463050	52.309	ug/L	99
42) Toluene	7.323	91	296593	5.058	ug/L	98
44) trans-1,3-Dichloropropene	7.590	75	83466	4.940	ug/L	97
45) 1,1,2-Trichloroethane	7.776	97	50437	4.834	ug/L	99
47) Tetrachloroethene	7.912	164	53132	4.555	ug/L	97
48) 2-Hexanone	8.085	43	340691	51.276	ug/L	99
49) Dibromochloromethane	8.185	129	57104	4.795	ug/L	99
50) 1,2-Dibromoethane	8.291	107	48193	4.821	ug/L	93
51) Chlorobenzene	8.821	112	171840	4.799	ug/L	97
52) Ethylbenzene	8.953	91	299601	4.861	ug/L	99
53) m,p-Xylene	9.079	106	108748	4.915	ug/L	99
54) o-Xylene	9.487	106	102594	4.921	ug/L	98
55) Styrene	9.503	104	187800	5.093	ug/L	99
57) 1,1,2,2-Tetrachloroethane	10.185	83	62240	5.008	ug/L	99
59) Bromoform	9.674	173	30314	4.558	ug/L	96
60) Isopropylbenzene	9.873	105	283482	4.786	ug/L	99
61) 1,2,3-Trichloropropane	10.217	75	44240	4.852	ug/L	96
62) 1,3,5-Trimethylbenzene	10.484	105	219526	4.560	ug/L	99
63) 1,2,4-Trimethylbenzene	10.860	105	234943	4.771	ug/L	100
64) 1,3-Dichlorobenzene	11.124	146	135473	4.764	ug/L	99
65) 1,4-Dichlorobenzene	11.217	146	136295	4.773	ug/L	99
67) 1,2-Dichlorobenzene	11.587	146	127381	4.864	ug/L	95
68) 1,2-Dibromo-3-chloropropane	12.374	75	9731	4.660	ug/L #	91
69) 1,3,5-Trimethylbenzene	12.590	180	92312	4.424	ug/L	99
70) 1,2,4-trimethylbenzene	13.207	180	69052	4.480	ug/L	99
71) Naphthalene	13.448	128	89998	4.186	ug/L	99
72) 1,2,3-Trimethylbenzene	13.686	180	63259	4.472	ug/L	99

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

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