

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032023\
 Data File : VU053304.D
 Acq On : 20 Mar 2023 13:50
 Operator : JC/MD
 Sample : 01956-05DL 20X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 E0288DL

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM031523WMA.M
 Title : VOC Analysis

Signal : TIC: VU053304.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.359	3	6	19	rVB	124690	110723	24.64%	2.630%
2	1.597	73	80	86	rBV	56790	63606	14.16%	1.511%
3	1.633	87	91	102	rVB	33293	35129	7.82%	0.835%
4	1.912	170	178	190	rBV	47403	62077	13.82%	1.475%
5	2.565	370	381	395	rBV	92714	147454	32.82%	3.503%
6	2.684	410	418	429	rVB5	2260	3802	0.85%	0.090%
7	3.137	551	559	564	rBV4	1323	2087	0.46%	0.050%
8	3.231	584	588	592	rBV2	491	444	0.10%	0.011%
9	3.465	657	661	664	rBV	234	188	0.04%	0.004%
10	3.983	812	822	838	rBV7	2323	6017	1.34%	0.143%
11	4.112	860	862	867	rVV2	308	182	0.04%	0.004%
12	4.533	980	993	1007	rVB4	4234	9612	2.14%	0.228%
13	4.613	1007	1018	1044	rBV	35352	87610	19.50%	2.081%
14	4.899	1103	1107	1110	rBV	163	192	0.04%	0.005%
15	5.057	1141	1156	1175	rBV	87165	189419	42.16%	4.500%
16	5.169	1189	1191	1194	rVB2	284	199	0.04%	0.005%
17	5.221	1204	1207	1210	rVB2	310	223	0.05%	0.005%
18	5.247	1211	1215	1218	rBV	245	209	0.05%	0.005%
19	5.385	1246	1258	1273	rBV3	15946	34740	7.73%	0.825%
20	5.562	1310	1313	1317	rVB	298	229	0.05%	0.005%
21	5.587	1318	1321	1325	rBV	161	179	0.04%	0.004%
22	5.722	1342	1363	1369	rBV2	167956	449310	100.00%	10.674%
23	5.957	1427	1436	1456	rVB3	12485	27697	6.16%	0.658%
24	6.134	1486	1491	1493	rBV	185	199	0.04%	0.005%
25	6.243	1512	1525	1548	rVV	146919	277250	61.71%	6.587%
26	6.472	1588	1596	1610	rBV4	3819	6826	1.52%	0.162%
27	6.526	1610	1613	1618	rVB2	471	323	0.07%	0.008%
28	6.684	1649	1662	1676	rBV	108905	211115	46.99%	5.015%
29	6.835	1704	1709	1712	rBV	282	282	0.06%	0.007%
30	6.973	1749	1752	1754	rBV	459	331	0.07%	0.008%
31	7.054	1775	1777	1782	rBV2	299	256	0.06%	0.006%
32	7.082	1783	1786	1791	rVV	205	177	0.04%	0.004%
33	7.179	1812	1816	1826	rBV4	988	1168	0.26%	0.028%
34	7.375	1875	1877	1883	rBV2	311	327	0.07%	0.008%
35	7.500	1914	1916	1920	rVV	327	195	0.04%	0.005%
36	7.562	1922	1935	1956	rVV2	62951	107617	23.95%	2.557%

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Integrator: RTE

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM031523WMA.M
 Title : VOC Analysis

37	7.690	1970	1975	1979	rBV4	720	669	0.15%	0.016%
38	7.848	2019	2024	2025	rVV2	1266	1022	0.23%	0.024%
39	7.893	2027	2038	2053	rVV	209359	361136	80.38%	8.579%
40	7.967	2054	2061	2076	rVB	12303	21593	4.81%	0.513%
41	8.176	2115	2126	2141	rBV2	42579	71633	15.94%	1.702%
42	8.237	2141	2145	2155	rVB6	2013	2735	0.61%	0.065%
43	8.571	2242	2249	2250	rBV3	1124	1284	0.29%	0.031%
44	8.626	2256	2266	2296	rBV2	104138	203413	45.27%	4.832%
45	8.857	2333	2338	2346	rVB5	1084	1725	0.38%	0.041%
46	8.980	2371	2376	2378	rBV2	271	267	0.06%	0.006%
47	9.240	2453	2457	2460	rBV	199	208	0.05%	0.005%
48	9.414	2497	2511	2515	rBV	208652	335041	74.57%	7.959%
49	9.565	2548	2558	2572	rVB	35940	57891	12.88%	1.375%
50	9.690	2582	2597	2615	rBV	32344	53349	11.87%	1.267%
51	9.857	2645	2649	2654	rVB2	367	388	0.09%	0.009%
52	9.941	2665	2675	2684	rBV6	4219	6752	1.50%	0.160%
53	9.979	2684	2687	2689	rBV2	372	221	0.05%	0.005%
54	10.005	2689	2695	2703	rBV4	3599	5467	1.22%	0.130%
55	10.111	2717	2728	2740	rBV3	4956	8250	1.84%	0.196%
56	10.266	2773	2776	2780	rBV2	371	198	0.04%	0.005%
57	10.352	2796	2803	2807	rBV3	336	410	0.09%	0.010%
58	10.481	2835	2843	2853	rVB2	5224	6831	1.52%	0.162%
59	10.642	2887	2893	2902	rVB5	1281	1594	0.35%	0.038%
60	10.751	2914	2927	2945	rVB	149893	235028	52.31%	5.583%
61	10.819	2945	2948	2951	rBV3	659	519	0.12%	0.012%
62	10.902	2964	2974	2984	rBV	5674	8246	1.84%	0.196%
63	10.992	2996	3002	3017	rVB2	5492	10430	2.32%	0.248%
64	11.336	3103	3109	3116	rBV3	1031	1223	0.27%	0.029%
65	11.397	3124	3128	3134	rVB2	999	970	0.22%	0.023%
66	11.465	3139	3149	3163	rBV	15758	23446	5.22%	0.557%
67	11.706	3221	3224	3228	rVB3	351	267	0.06%	0.006%
68	11.738	3228	3234	3242	rBV4	1643	2072	0.46%	0.049%
69	11.806	3242	3255	3274	rVV	212553	337427	75.10%	8.016%
70	11.893	3277	3282	3289	rVB4	1600	2119	0.47%	0.050%
71	12.092	3336	3344	3350	rBV3	3963	5553	1.24%	0.132%
72	12.188	3361	3374	3389	rBV	206510	337216	75.05%	8.011%
73	12.301	3403	3409	3415	rBV5	820	918	0.20%	0.022%
74	12.462	3453	3459	3465	rBV4	1733	2256	0.50%	0.054%
75	12.568	3479	3492	3501	rBV	5488	7658	1.70%	0.182%
76	12.680	3519	3527	3534	rBV4	2631	4305	0.96%	0.102%

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77	12.864	3577	3584	3590	rVB3	613	852	0.19%	0.020%
78	12.915	3596	3600	3602	rBV	559	431	0.10%	0.010%
79	12.970	3612	3617	3625	rVB5	4618	5987	1.33%	0.142%
80	13.021	3625	3633	3643	rBV2	6867	9621	2.14%	0.229%
81	13.163	3671	3677	3682	rBV	403	471	0.10%	0.011%
82	13.323	3723	3727	3735	rVB4	1714	1974	0.44%	0.047%
83	13.442	3755	3764	3782	rVB5	18831	30395	6.76%	0.722%
84	13.552	3793	3798	3808	rBV3	802	1137	0.25%	0.027%
85	13.619	3808	3819	3830	rVV	45332	72208	16.07%	1.715%
86	13.725	3849	3852	3854	rBV	340	242	0.05%	0.006%
87	13.777	3864	3868	3875	rVB3	738	741	0.16%	0.018%
88	13.915	3902	3911	3914	rBV4	547	979	0.22%	0.023%
89	13.934	3914	3917	3927	rVB5	1688	1945	0.43%	0.046%
90	14.085	3956	3964	3979	rVB	18196	30127	6.71%	0.716%
91	14.143	3979	3982	3988	rVV2	280	299	0.07%	0.007%
92	14.188	3988	3996	4004	rVB4	4112	5706	1.27%	0.136%
93	14.253	4014	4016	4018	rBV	349	196	0.04%	0.005%
94	14.278	4018	4024	4034	rVV6	3362	4741	1.06%	0.113%
95	14.683	4140	4150	4158	rVB4	12220	18519	4.12%	0.440%
96	14.873	4207	4209	4212	rVB3	573	370	0.08%	0.009%
97	15.011	4244	4252	4260	rBV4	12684	18055	4.02%	0.429%
98	15.220	4307	4317	4327	rVV	11667	18653	4.15%	0.443%
99	15.407	4365	4375	4386	rBV2	16747	26228	5.84%	0.623%
100	15.497	4400	4403	4406	rBV	447	345	0.08%	0.008%

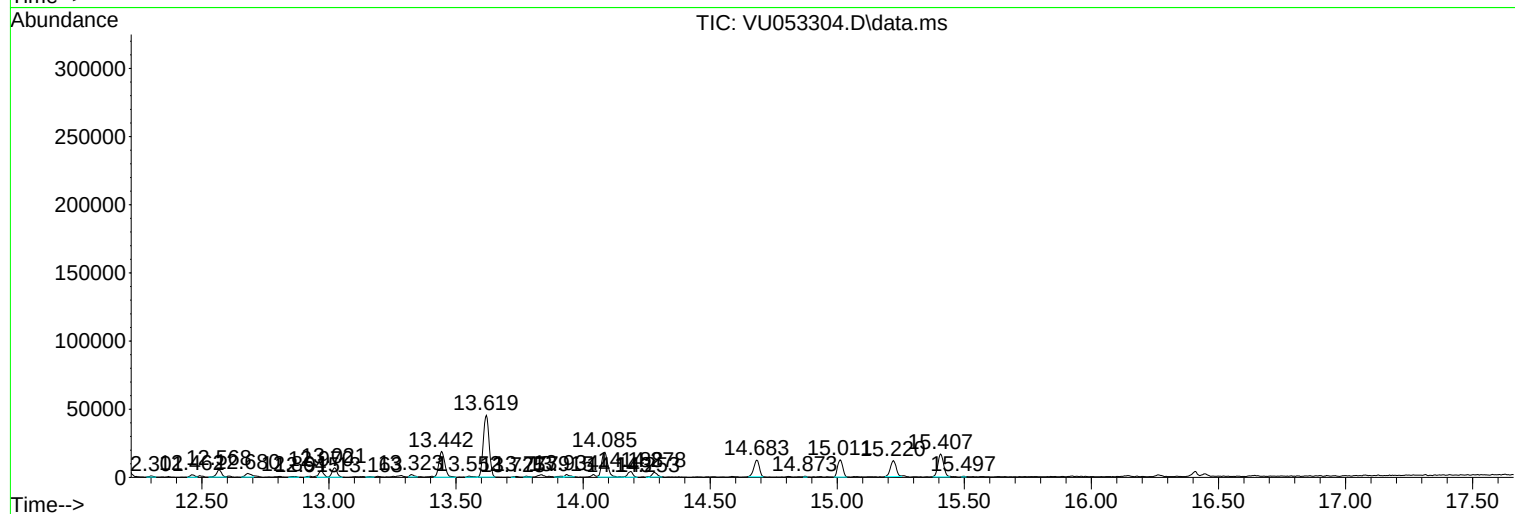
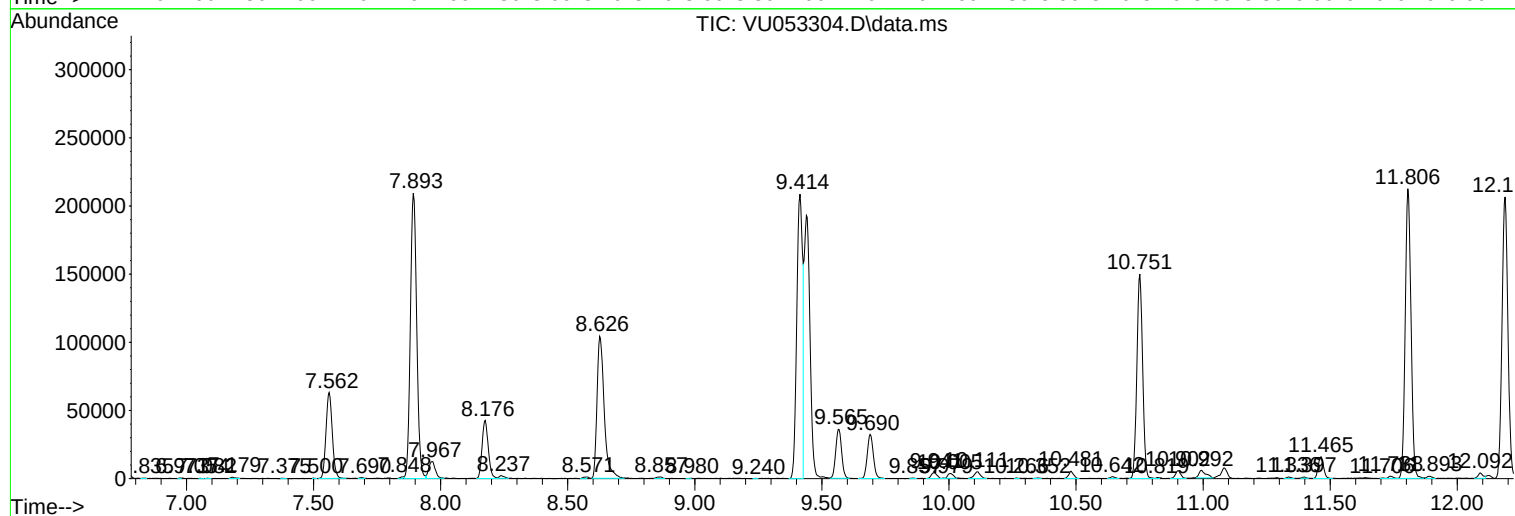
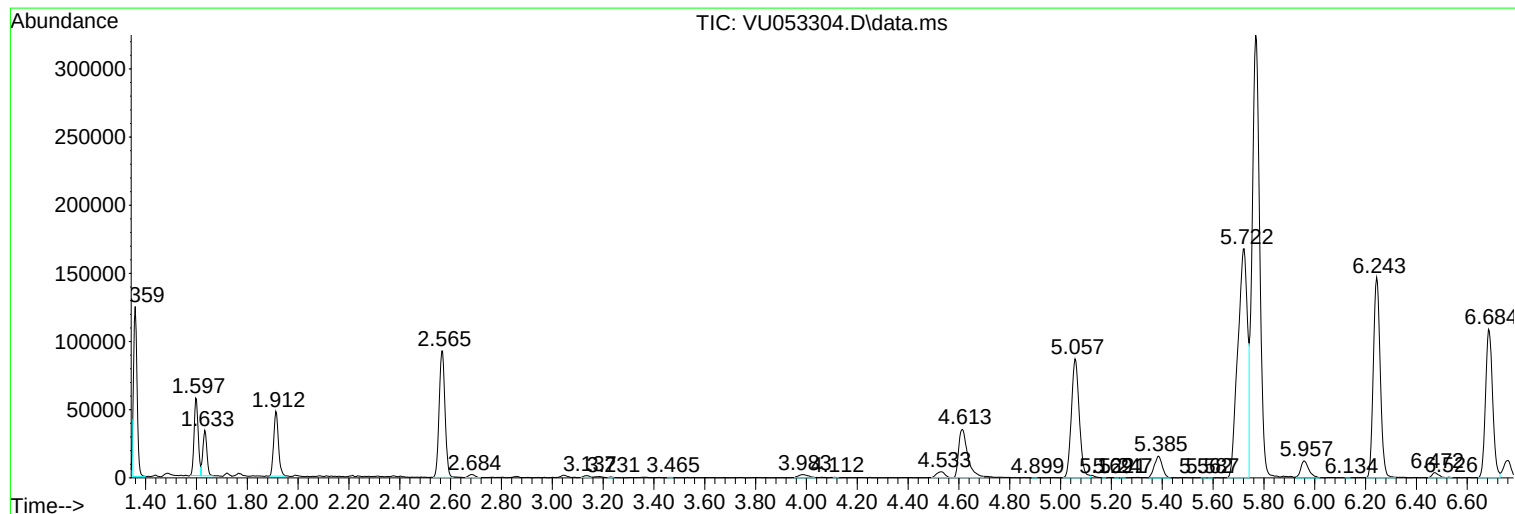
Sum of corrected areas: 4209346

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Instrument :
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ClientSampleId :
E0288DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM031523WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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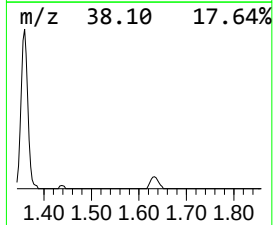
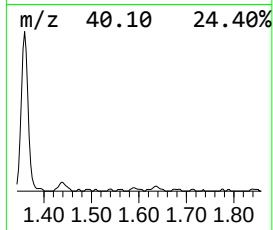
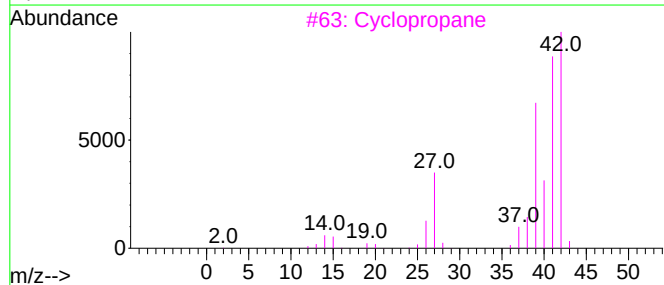
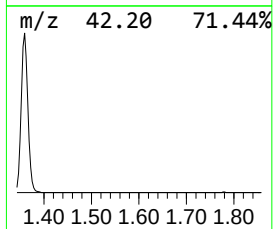
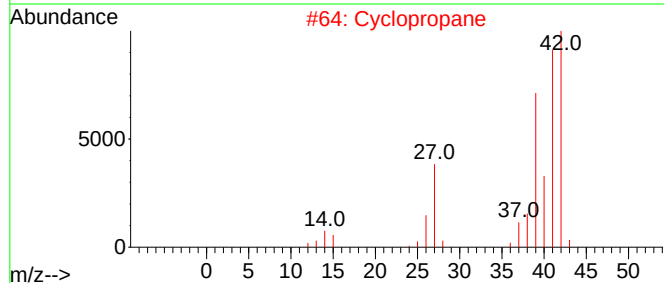
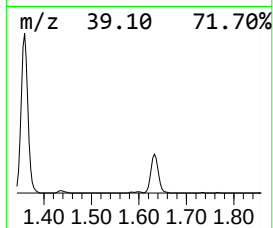
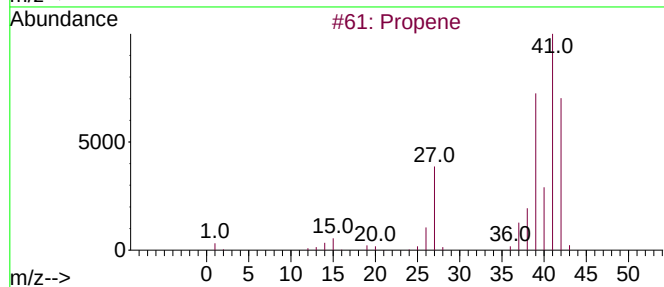
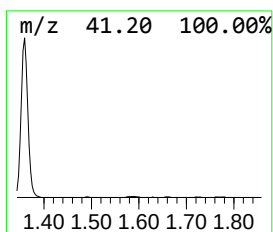
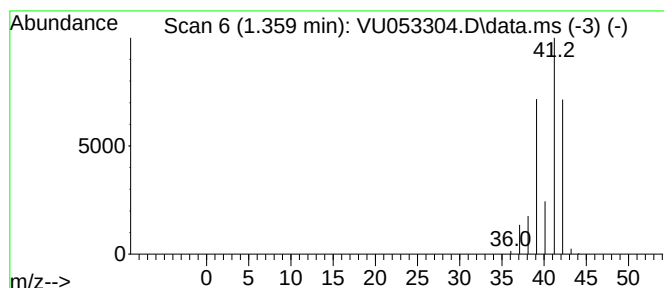
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM031523WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.359	19.97 ug/L	110723	1,4-Difluorobenzene	6.243

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propene	42	C3H6	000115-07-1	90
2			Cyclopropane	42	C3H6	000075-19-4	86
3			Cyclopropane	42	C3H6	000075-19-4	78
4			Propene	42	C3H6	000115-07-1	50
5			Cyclopropene	40	C3H4	002781-85-3	16



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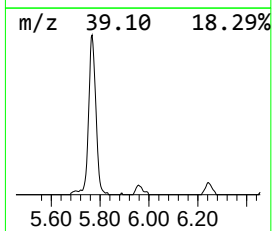
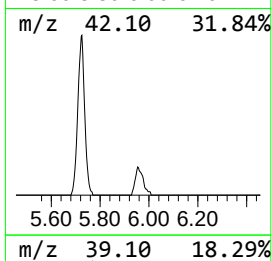
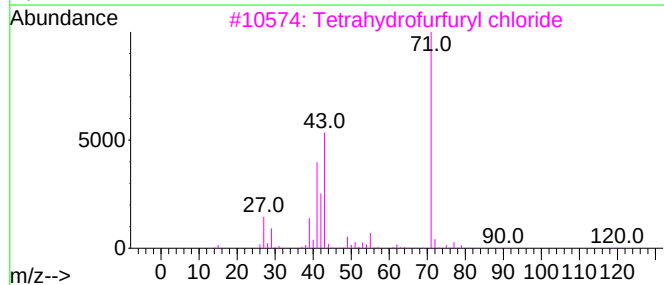
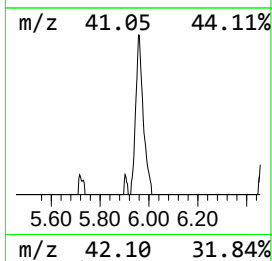
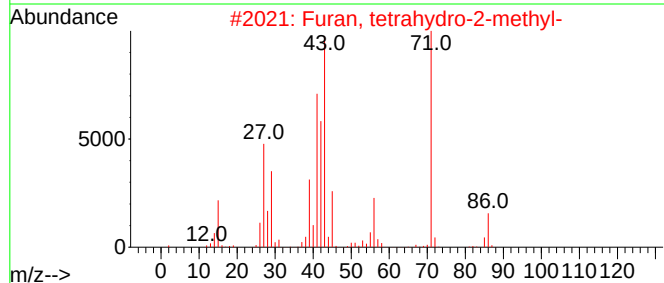
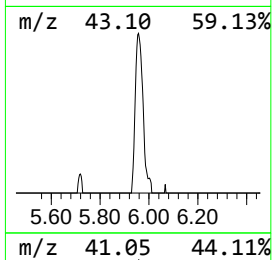
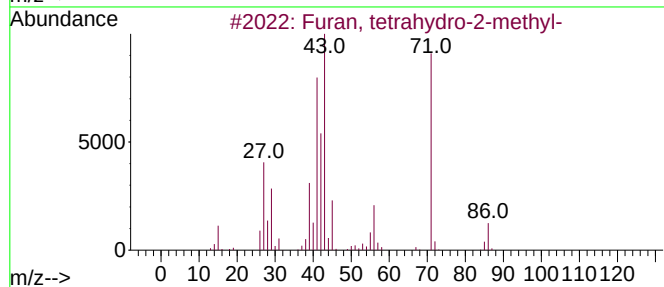
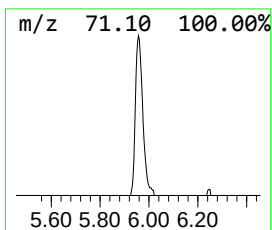
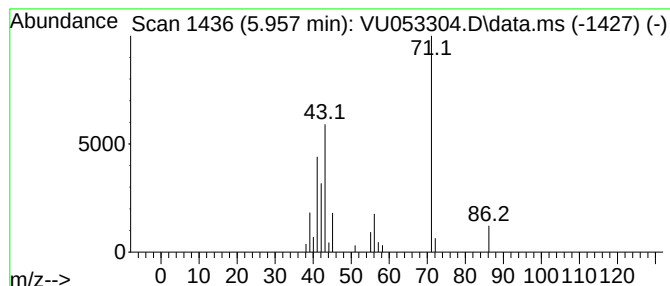
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM031523WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Furan, tetrahydro-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.957	4.99 ug/L	27697	1,4-Difluorobenzene	6.243

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	91
2			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	91
3			Tetrahydrofurfuryl chloride	120	C5H9ClO	003003-84-7	59
4			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	53
5			3-Penten-2-ol	86	C5H10O	001569-50-2	53



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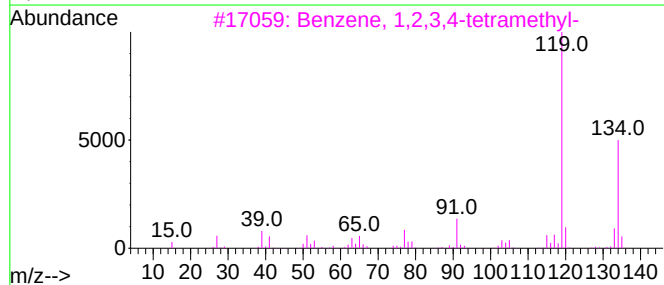
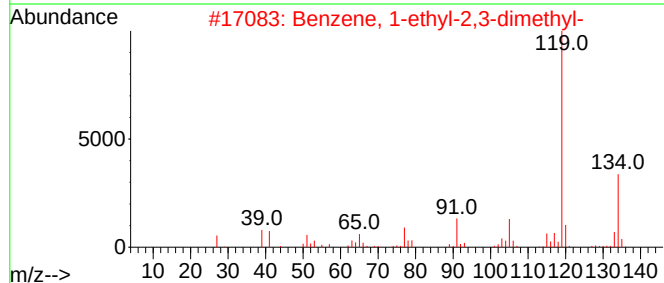
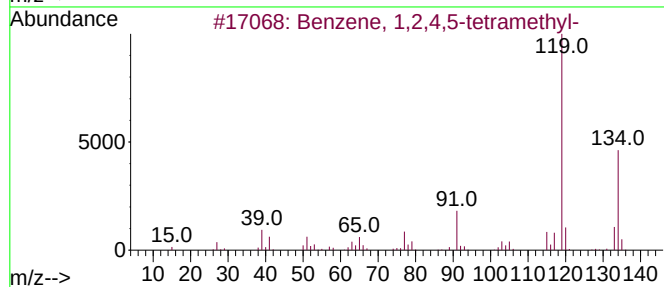
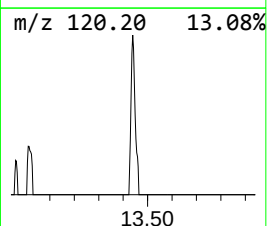
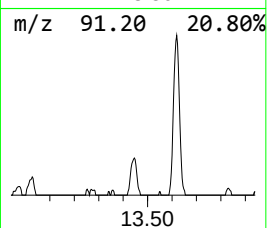
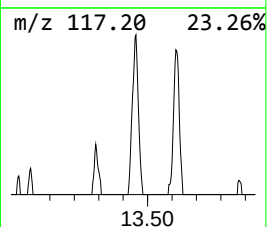
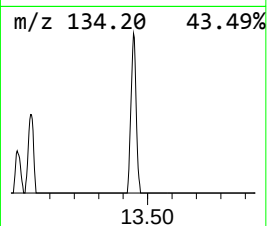
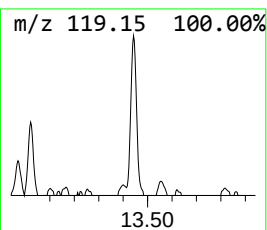
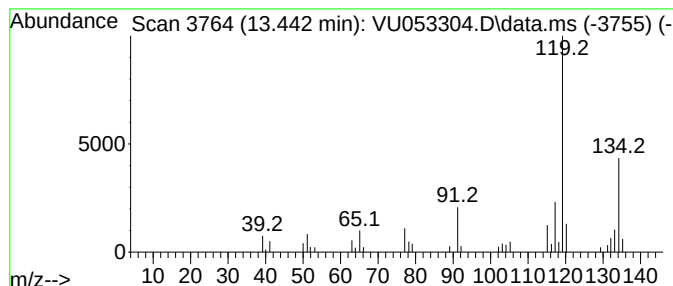
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM031523WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.442	4.50 ug/L	30395	1,4-Dichlorobenzene-d4	11.806

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	90
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	90
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	87
5			p-Cymene	134	C10H14	000099-87-6	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032023\
Data File : VU053304.D
Acq On : 20 Mar 2023 13:50
Operator : JC/MD
Sample : 01956-05DL 20X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E0288DL

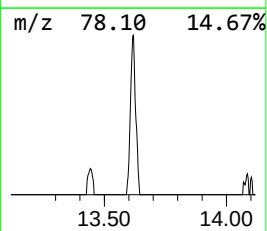
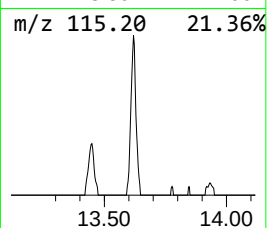
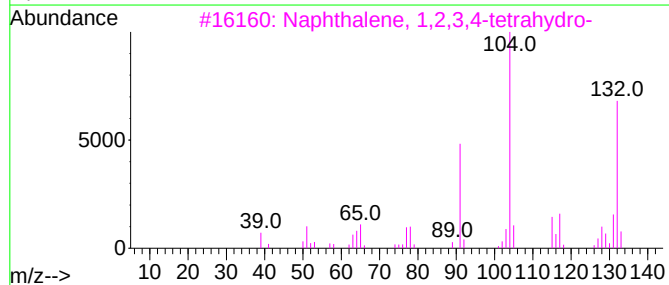
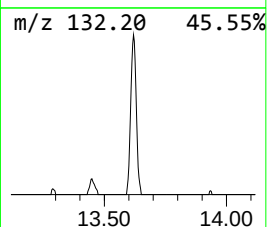
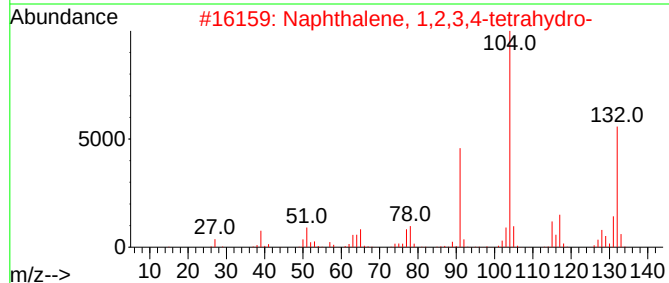
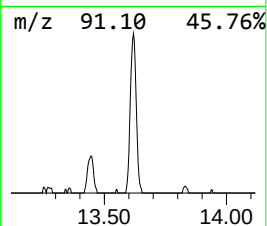
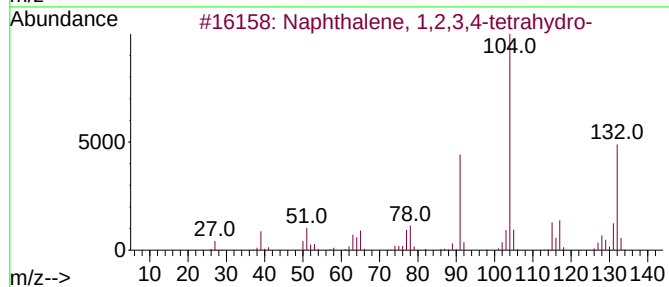
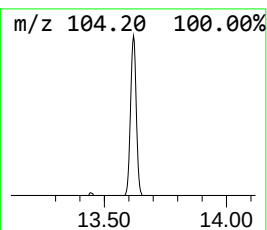
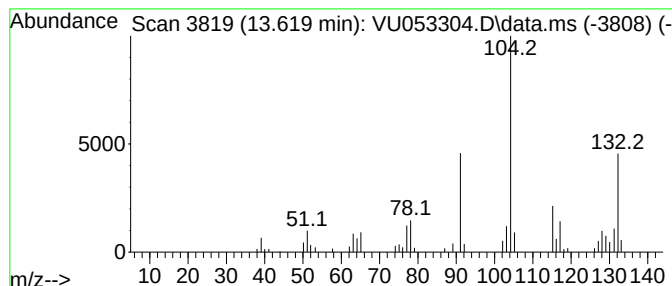
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.619	10.70 ug/L	72208	1,4-Dichlorobenzene-d4	11.806

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
3			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
4			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	93
5			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032023\
Data File : VU053304.D
Acq On : 20 Mar 2023 13:50
Operator : JC/MD
Sample : 01956-05DL 20X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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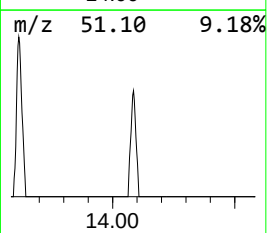
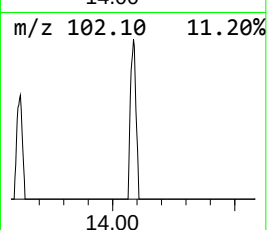
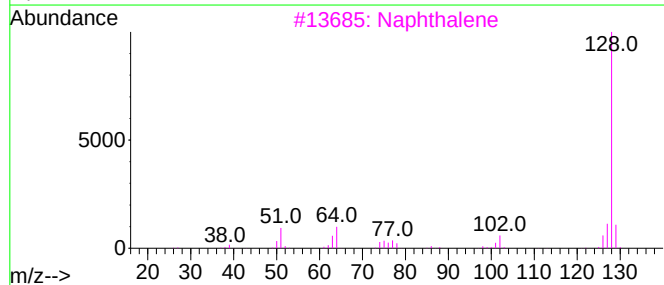
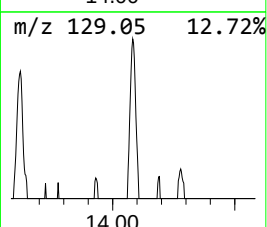
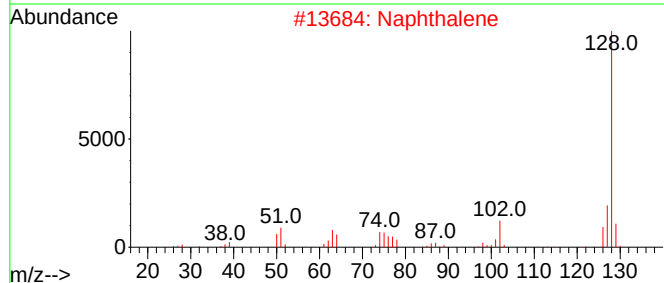
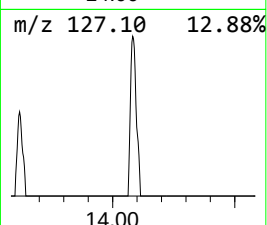
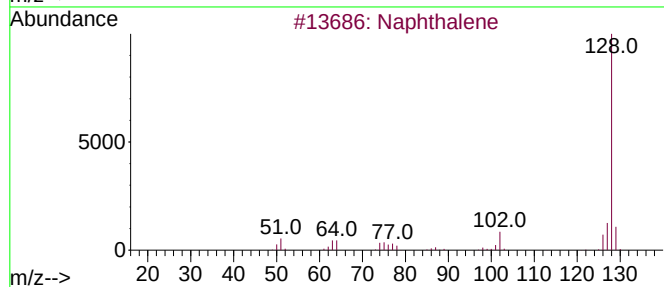
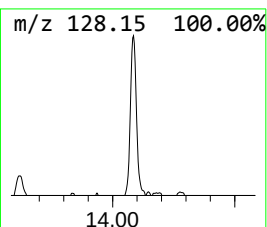
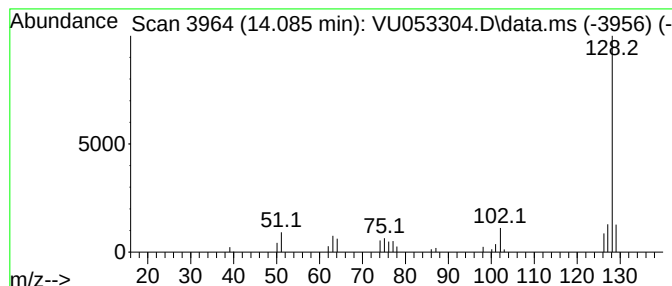
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Azulene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.085	4.46 ug/L	30127	1,4-Dichlorobenzene-d4	11.806

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Naphthalene	128	C10H8	000091-20-3	95
3			Naphthalene	128	C10H8	000091-20-3	94
4			Azulene	128	C10H8	000275-51-4	94
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032023\
Data File : VU053304.D
Acq On : 20 Mar 2023 13:50
Operator : JC/MD
Sample : 01956-05DL 20X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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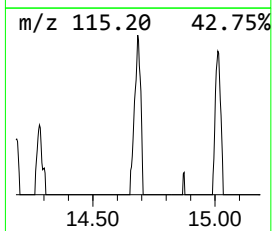
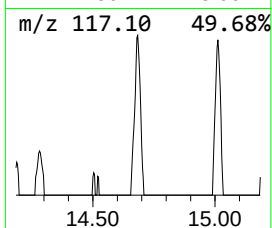
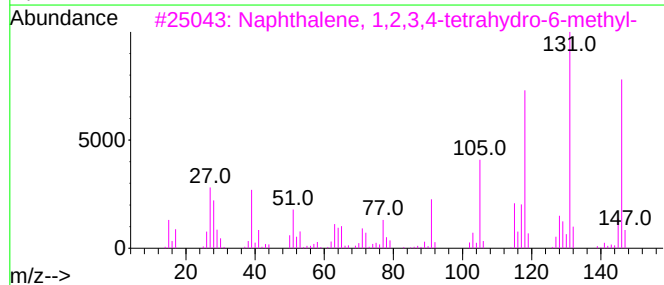
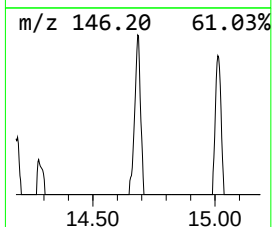
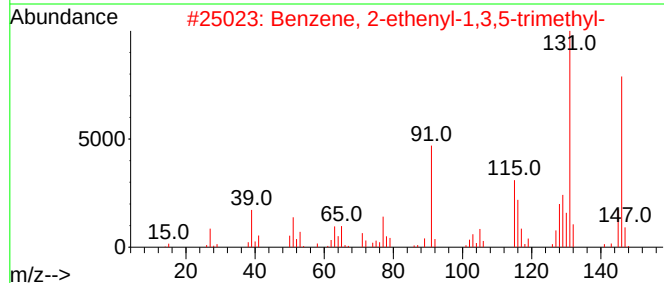
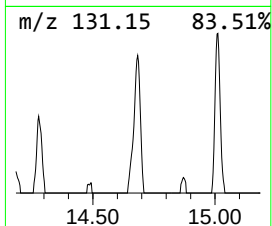
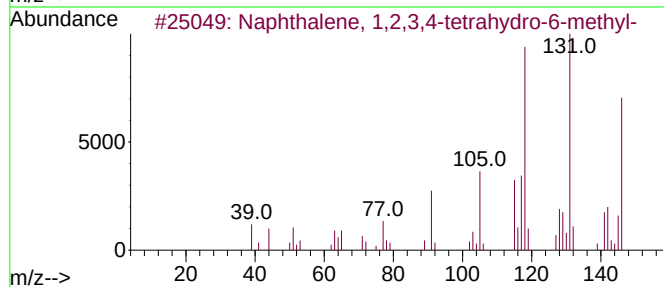
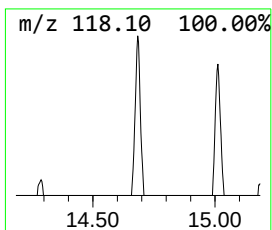
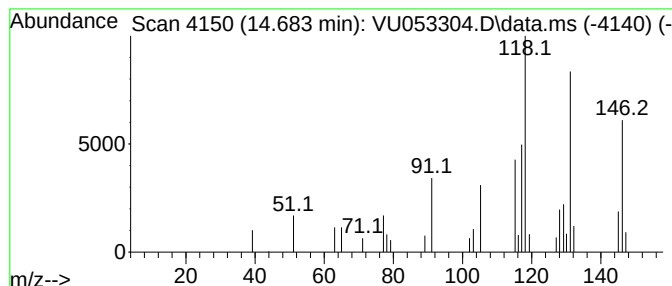
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	2.74 ug/L	18519	1,4-Dichlorobenzene-d4	11.806

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	93
2			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	83
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	83
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	81
5			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032023\
Data File : VU053304.D
Acq On : 20 Mar 2023 13:50
Operator : JC/MD
Sample : 01956-05DL 20X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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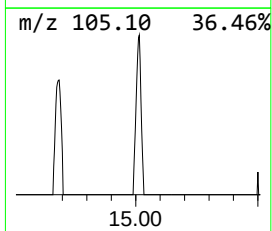
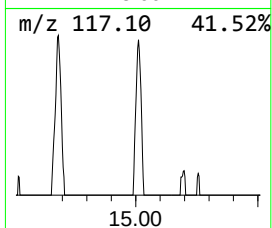
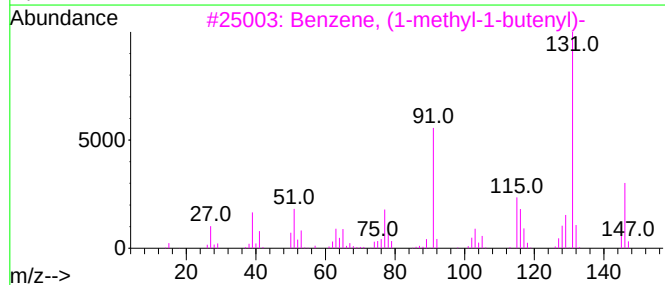
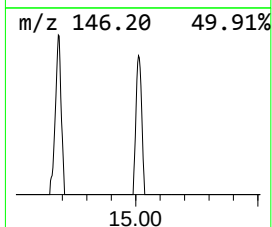
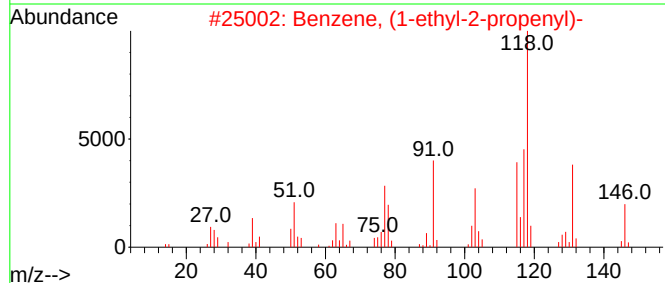
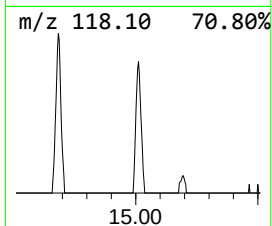
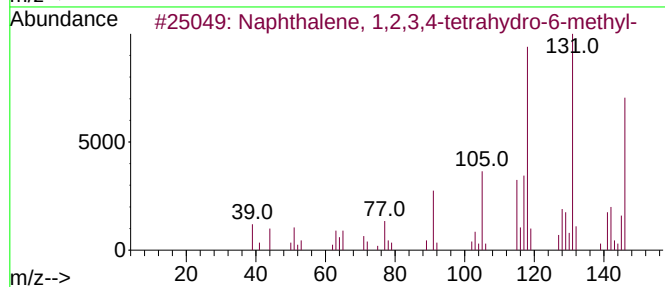
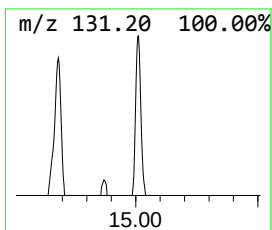
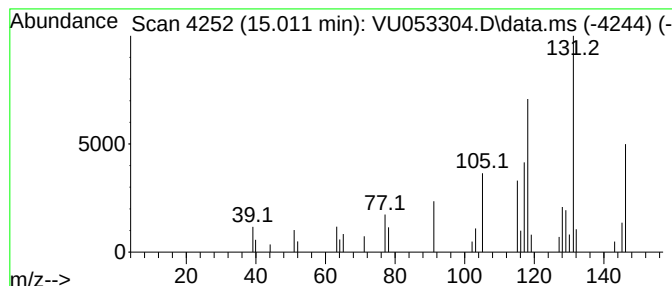
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, (1-ethyl-2-propenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.011	2.68 ug/L	18055	1,4-Dichlorobenzene-d4	11.806

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
2			Benzene, (1-ethyl-2-propenyl)-	146	C11H14	019947-22-9	62
3			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	49
4			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	43
5			Benzene, (1-methylenebutyl)-	146	C11H14	005676-32-4	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032023\
Data File : VU053304.D
Acq On : 20 Mar 2023 13:50
Operator : JC/MD
Sample : 01956-05DL 20X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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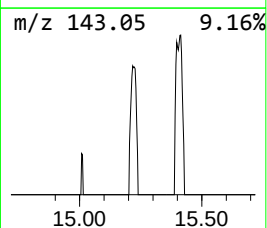
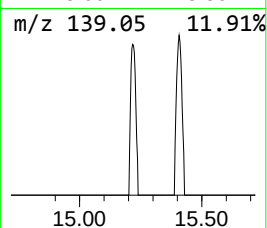
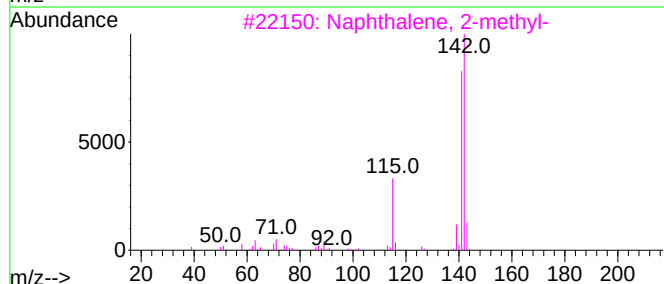
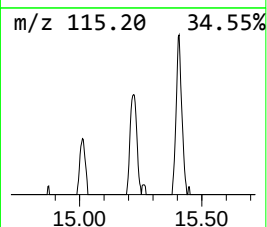
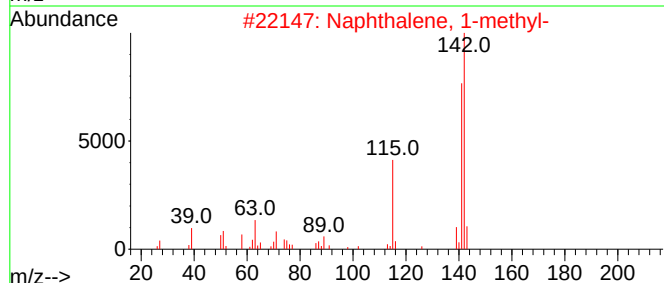
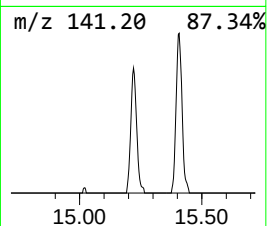
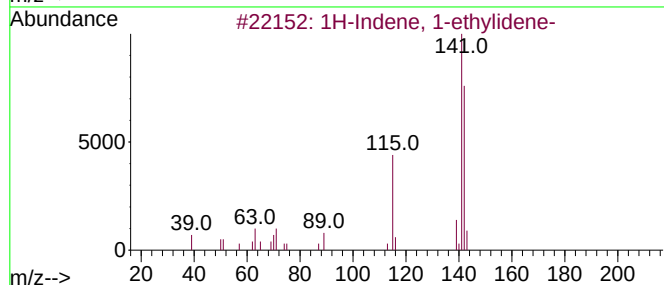
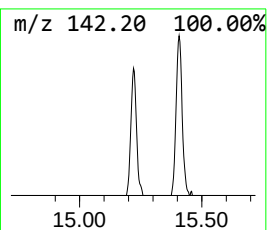
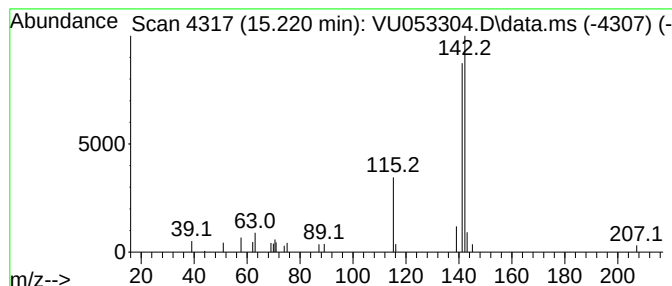
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1H-Indene, 1-ethylidene- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.220	2.76 ug/L	18653	1,4-Dichlorobenzene-d4	11.806

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	91
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	90
3			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	87
4			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	87
5			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032023\
Data File : VU053304.D
Acq On : 20 Mar 2023 13:50
Operator : JC/MD
Sample : 01956-05DL 20X
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ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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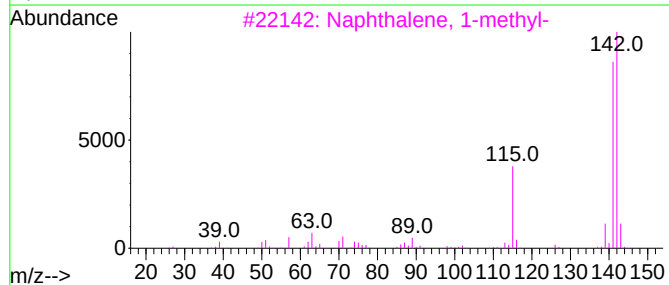
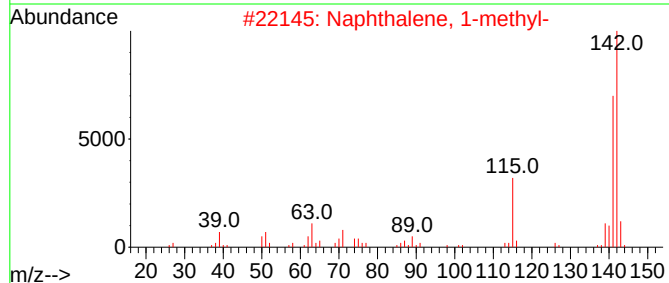
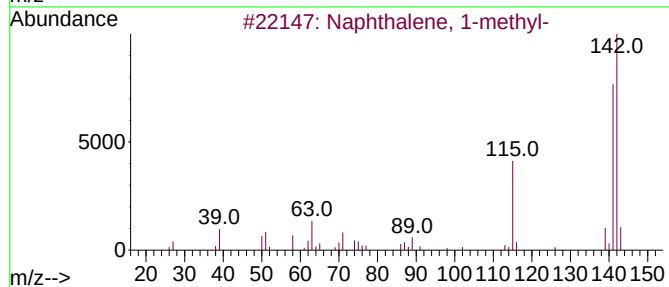
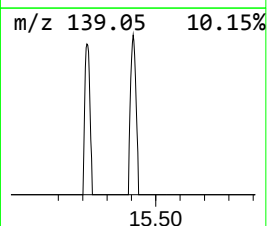
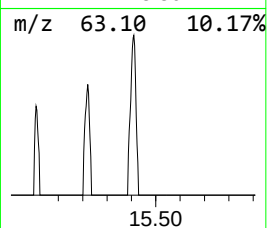
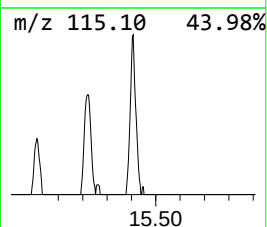
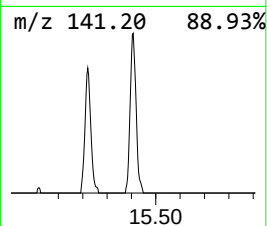
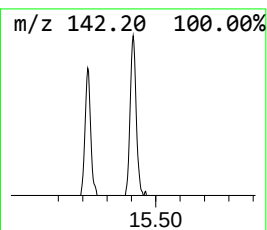
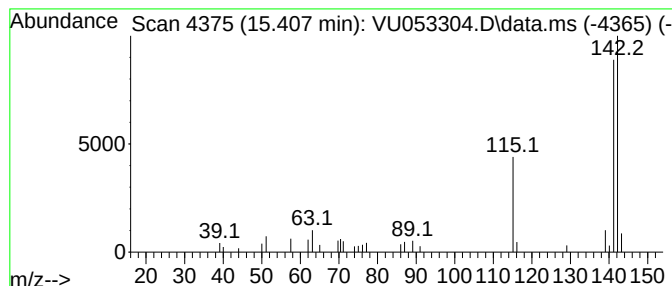
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.407	3.89 ug/L	26228	1,4-Dichlorobenzene-d4	11.806

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
3		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
4		Benzocycloheptatriene	142	C11H10	000264-09-5	90
5		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032023\
Data File : VU053304.D
Acq On : 20 Mar 2023 13:50
Operator : JC/MD
Sample : 01956-05DL 20X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E0288DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM031523WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Furan, tetrahyd...	5.957	5.0	ug/L	27697	1	6.243	277250	50.0
Benzene, 1,2,4,...	13.442	4.5	ug/L	30395	3	11.806	337427	50.0
Naphthalene, 1,...	13.619	10.7	ug/L	72208	3	11.806	337427	50.0
Azulene	14.085	4.5	ug/L	30127	3	11.806	337427	50.0
Naphthalene, 1,...	14.684	2.7	ug/L	18519	3	11.806	337427	50.0
Benzene, (1-eth...	15.011	2.7	ug/L	18055	3	11.806	337427	50.0
1H-Indene, 1-et...	15.220	2.8	ug/L	18653	3	11.806	337427	50.0
Naphthalene, 1-...	15.407	3.9	ug/L	26228	3	11.806	337427	50.0