

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU010522\
 Data File : VU046673.D
 Acq On : 05 Jan 2022 19:29
 Operator : SY/MD
 Sample : N1025-04
 Misc : 6.80g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BGLG8

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\SFAMULM010522WMA.M
 Title : VOC Analysis

Signal : TIC: VU046673.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.600	73	81	112	rVB	100506	207834	15.11%	1.669%
2	1.883	165	169	170	rBV	1487	1017	0.07%	0.008%
3	1.919	172	180	219	rVB	51312	168203	12.23%	1.351%
4	2.555	365	378	419	rBV	241297	430640	31.30%	3.459%
5	2.771	441	445	450	rBV2	316	401	0.03%	0.003%
6	2.813	455	458	461	rBV2	266	155	0.01%	0.001%
7	3.047	521	531	541	rVV2	1602	3194	0.23%	0.026%
8	3.089	541	544	547	rVB2	272	169	0.01%	0.001%
9	3.137	556	559	561	rBV	203	158	0.01%	0.001%
10	3.186	561	574	585	rVV2	4369	9591	0.70%	0.077%
11	3.520	676	678	684	rBB	156	175	0.01%	0.001%
12	3.681	725	728	731	rBV2	386	284	0.02%	0.002%
13	4.661	1017	1033	1091	rBV	53757	329604	23.96%	2.648%
14	5.067	1142	1159	1186	rVB	229324	539857	39.24%	4.336%
15	5.369	1248	1253	1258	rVB5	717	877	0.06%	0.007%
16	5.584	1316	1320	1325	rBB2	804	796	0.06%	0.006%
17	5.723	1342	1363	1389	rVV2	534134	1375751	100.00%	11.051%
18	5.886	1411	1414	1415	rBV2	304	147	0.01%	0.001%
19	5.983	1433	1444	1457	rBV6	2784	6886	0.50%	0.055%
20	6.041	1460	1462	1467	rVV2	674	530	0.04%	0.004%
21	6.128	1482	1489	1497	rBB3	1260	1979	0.14%	0.016%
22	6.250	1511	1527	1561	rVV	273862	552877	40.19%	4.441%
23	6.369	1561	1564	1567	rVV	335	310	0.02%	0.002%
24	6.385	1567	1569	1574	rVV2	306	300	0.02%	0.002%
25	6.526	1602	1613	1621	rVB7	1917	3489	0.25%	0.028%
26	6.604	1627	1637	1638	rBV	1127	1284	0.09%	0.010%
27	6.694	1649	1665	1692	rBV	287500	608156	44.21%	4.885%
28	6.806	1695	1700	1708	rVB4	1141	1367	0.10%	0.011%
29	6.835	1708	1709	1714	rVB2	230	177	0.01%	0.001%
30	6.857	1714	1716	1719	rBB2	235	149	0.01%	0.001%
31	7.060	1775	1779	1783	rBV	305	323	0.02%	0.003%
32	7.105	1789	1793	1796	rBB2	234	195	0.01%	0.002%
33	7.179	1813	1816	1817	rBV	239	166	0.01%	0.001%
34	7.192	1817	1820	1825	rBV4	510	473	0.03%	0.004%
35	7.234	1829	1833	1834	rBV2	575	378	0.03%	0.003%
36	7.301	1851	1854	1858	rBB	164	153	0.01%	0.001%

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Integrator: RTE

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Peak Location: TOP

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM010522WMA.M
 Title : VOC Analysis

37	7.340	1863	1866	1870	rBV	159	190	0.01%	0.002%
38	7.411	1883	1888	1889	rBV	145	149	0.01%	0.001%
39	7.497	1908	1915	1919	rBB3	626	696	0.05%	0.006%
40	7.568	1923	1937	1972	rBV2	146952	323224	23.49%	2.596%
41	7.693	1972	1976	1979	rBV2	405	420	0.03%	0.003%
42	7.780	2000	2003	2004	rBV2	454	253	0.02%	0.002%
43	7.899	2025	2040	2056	rVV	612852	1166981	84.83%	9.374%
44	7.970	2058	2062	2083	rVB	25772	43943	3.19%	0.353%
45	8.185	2116	2129	2162	rVV	80535	223250	16.23%	1.793%
46	8.365	2179	2185	2191	rBB3	705	799	0.06%	0.006%
47	8.462	2212	2215	2219	rBV	383	325	0.02%	0.003%
48	8.536	2233	2238	2241	rVB2	313	255	0.02%	0.002%
49	8.584	2250	2253	2256	rBV	177	162	0.01%	0.001%
50	8.681	2261	2283	2314	rBV	280710	796423	57.89%	6.397%
51	8.864	2338	2340	2347	rVB2	1378	1331	0.10%	0.011%
52	8.925	2351	2359	2363	rBV4	1001	1337	0.10%	0.011%
53	9.044	2392	2396	2399	rBV2	698	753	0.05%	0.006%
54	9.124	2414	2421	2428	rVV6	1801	2624	0.19%	0.021%
55	9.169	2429	2435	2440	rVV4	2027	3092	0.22%	0.025%
56	9.214	2442	2449	2458	rVB4	6231	10794	0.78%	0.087%
57	9.279	2464	2469	2472	rBV	131	185	0.01%	0.001%
58	9.340	2477	2488	2497	rVV3	6955	11521	0.84%	0.093%
59	9.430	2501	2516	2541	rBV	302704	714403	51.93%	5.738%
60	9.574	2550	2561	2577	rVV	15891	32021	2.33%	0.257%
61	9.639	2577	2581	2582	rVV3	616	490	0.04%	0.004%
62	9.655	2583	2586	2587	rVV2	926	626	0.05%	0.005%
63	9.703	2588	2601	2614	rVV	40829	87599	6.37%	0.704%
64	9.764	2618	2620	2631	rVB6	1939	2443	0.18%	0.020%
65	9.915	2664	2667	2673	rBV	128	201	0.01%	0.002%
66	9.957	2673	2680	2685	rBV3	919	1107	0.08%	0.009%
67	10.018	2689	2699	2712	rBV7	8302	16088	1.17%	0.129%
68	10.115	2718	2729	2740	rBV3	20085	37618	2.73%	0.302%
69	10.169	2740	2746	2754	rVB5	4023	6303	0.46%	0.051%
70	10.218	2754	2761	2762	rBV2	1382	1426	0.10%	0.011%
71	10.250	2762	2771	2784	rVB3	13975	25011	1.82%	0.201%
72	10.317	2784	2792	2794	rBV2	2639	3570	0.26%	0.029%
73	10.353	2797	2803	2812	rVV9	7594	13562	0.99%	0.109%
74	10.394	2813	2816	2828	rVB3	7663	11237	0.82%	0.090%
75	10.491	2828	2846	2855	rBV3	7931	20165	1.47%	0.162%
76	10.558	2855	2867	2874	rBV5	9976	20147	1.46%	0.162%

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 Title : VOC Analysis

77	10.771	2921	2933	2959	rVB	464934	795175	57.80%	6.387%
78	10.912	2970	2977	2986	rVB	13276	19913	1.45%	0.160%
79	11.028	2999	3013	3019	rBV	16900	42670	3.10%	0.343%
80	11.246	3077	3081	3088	rBV6	2404	2330	0.17%	0.019%
81	11.307	3090	3100	3113	rVV5	18407	42449	3.09%	0.341%
82	11.378	3115	3122	3127	rVV3	12194	18849	1.37%	0.151%
83	11.417	3128	3134	3141	rVV2	34070	49981	3.63%	0.401%
84	11.475	3142	3152	3173	rVB	119860	226564	16.47%	1.820%
85	11.574	3174	3183	3191	rBV2	15825	28034	2.04%	0.225%
86	11.706	3215	3224	3232	rBV6	21928	41635	3.03%	0.334%
87	11.819	3248	3259	3271	rBV	482733	770154	55.98%	6.186%
88	11.873	3272	3276	3282	rBV3	14153	17325	1.26%	0.139%
89	12.105	3339	3348	3360	rBV5	13853	30826	2.24%	0.248%
90	12.201	3364	3378	3392	rVB	714907	1187276	86.30%	9.537%
91	12.475	3456	3463	3465	rBV4	12897	16287	1.18%	0.131%
92	12.668	3518	3523	3531	rVB2	11273	14814	1.08%	0.119%
93	13.838	3882	3887	3902	rVB7	10741	17936	1.30%	0.144%
94	14.089	3954	3965	3985	rVB	489112	784403	57.02%	6.301%
95	14.192	3988	3997	4012	rVB5	20833	44680	3.25%	0.359%
96	14.449	4071	4077	4084	rBV2	12686	16460	1.20%	0.132%
97	15.224	4306	4318	4329	rBV	104097	174543	12.69%	1.402%
98	15.407	4365	4375	4398	rVB2	127709	223993	16.28%	1.799%
99	16.262	4634	4641	4644	rBV3	14802	19684	1.43%	0.158%
100	16.410	4680	4687	4694	rBV	23433	32626	2.37%	0.262%

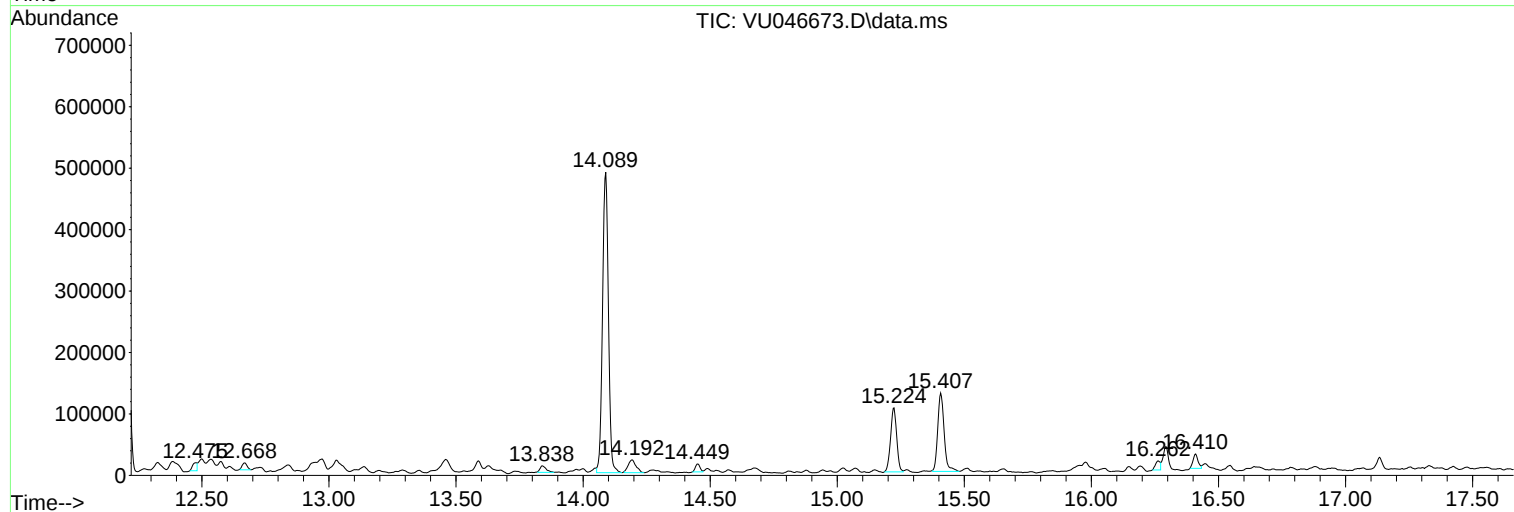
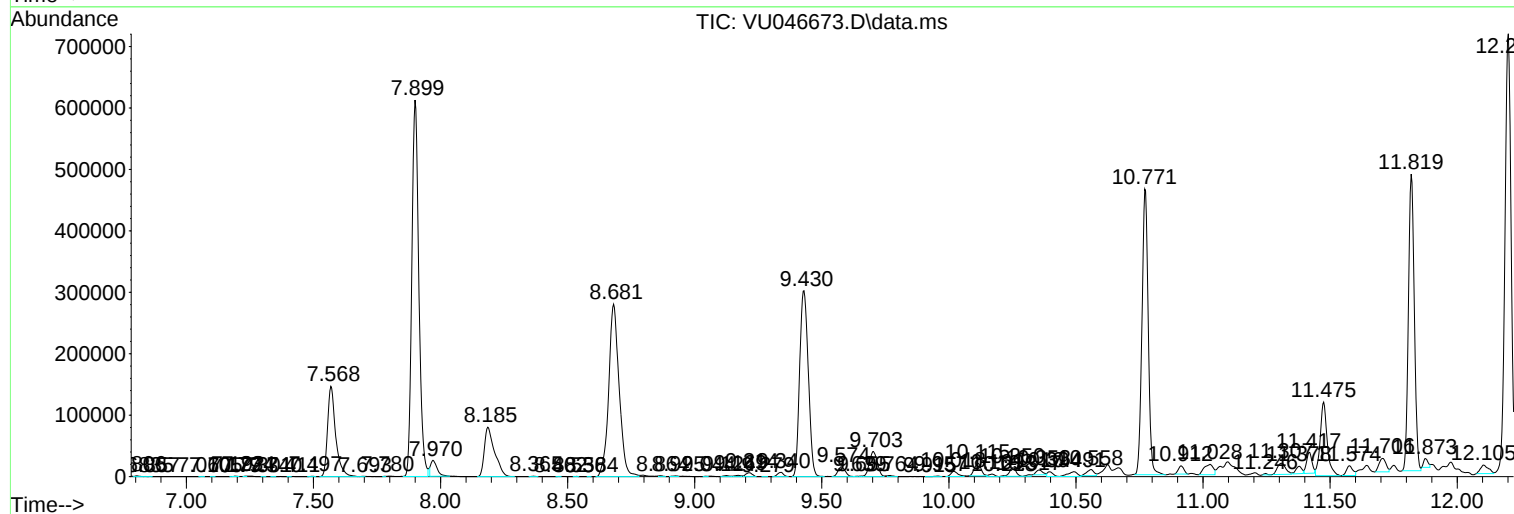
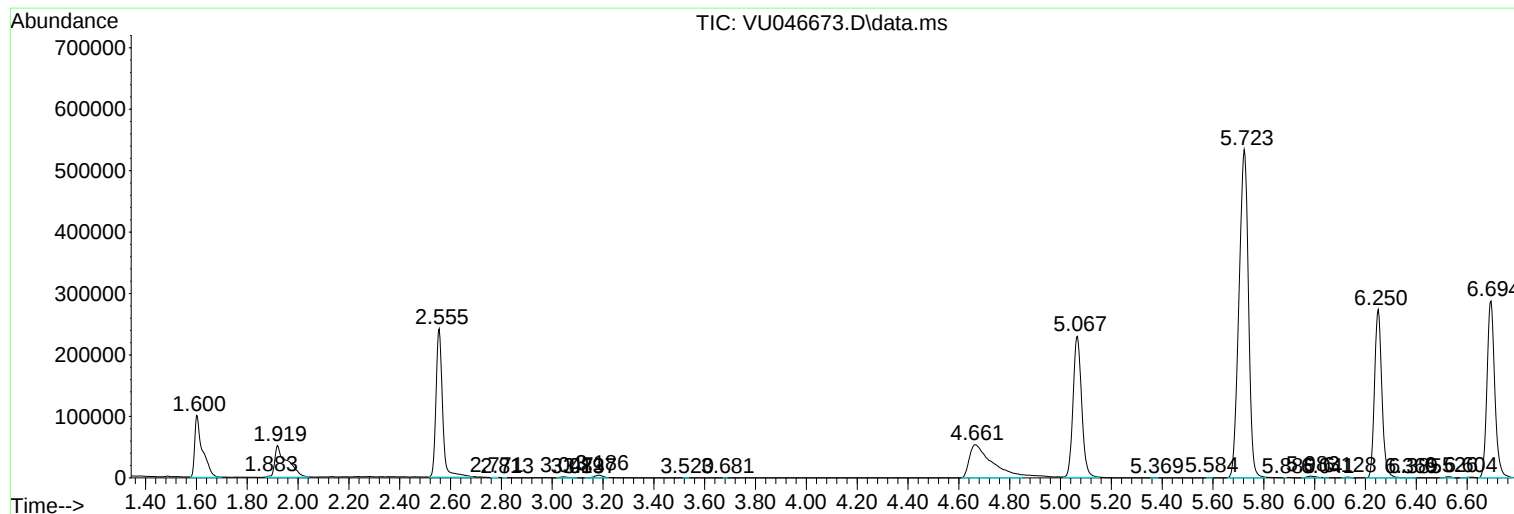
Sum of corrected areas: 12449376

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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM010522WMA.M
Quant Title : VOC Analysis

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



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Instrument :
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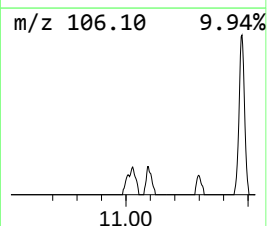
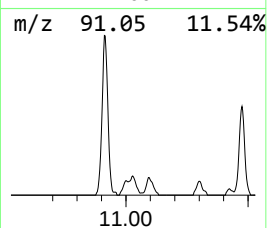
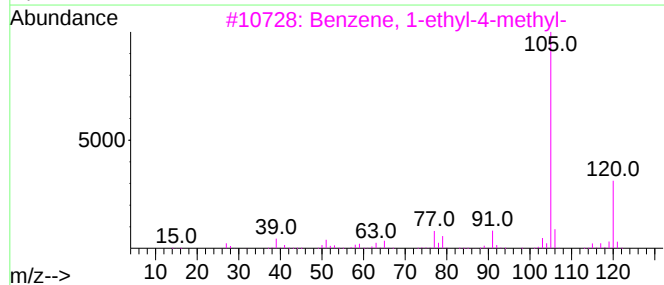
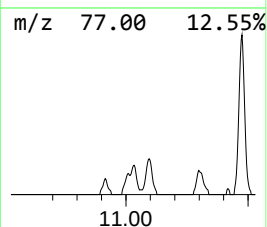
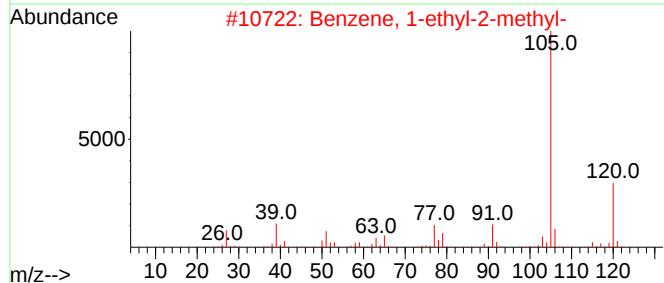
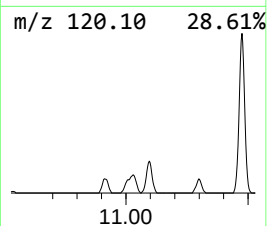
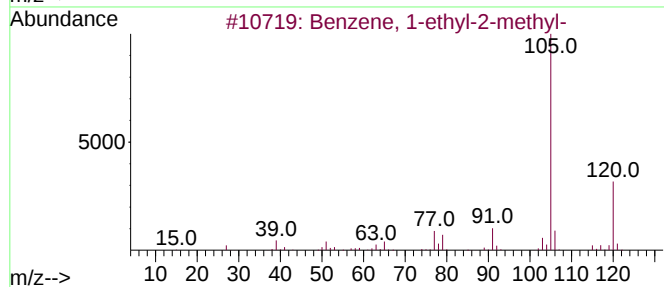
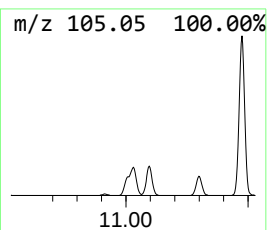
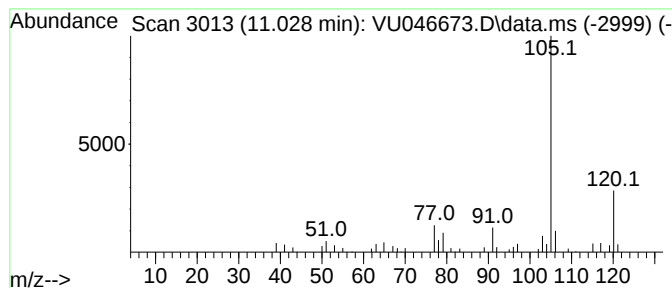
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM010522WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.028	2.77 ug/L	42670	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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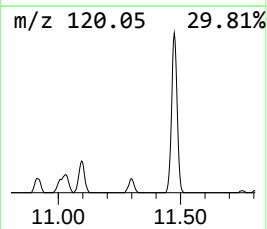
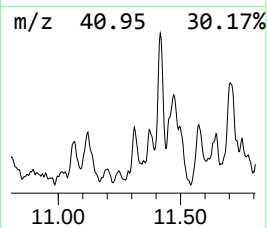
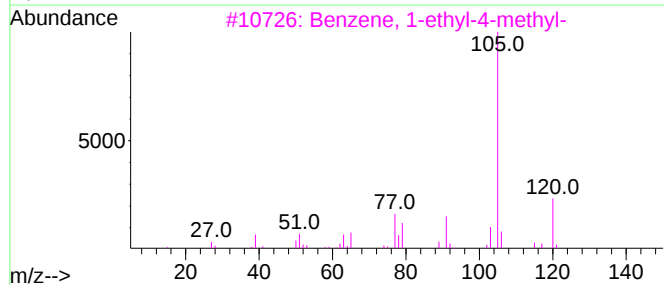
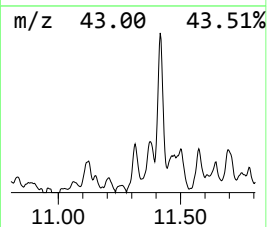
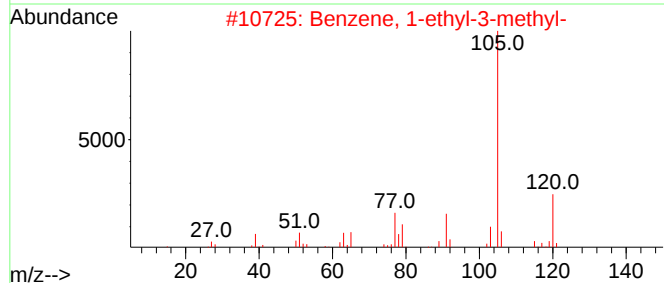
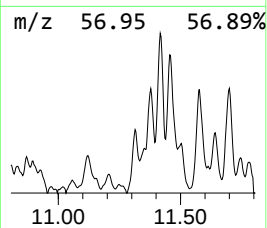
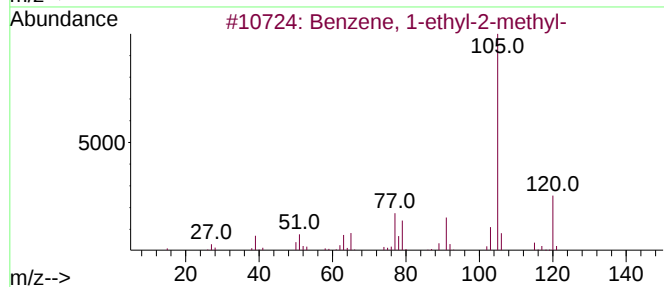
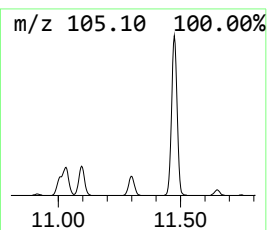
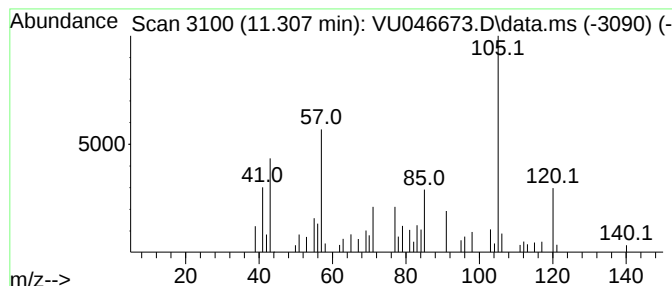
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM010522WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 Benzene, 1-ethyl-3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.307	2.76 ug/L	42449	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	50
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	50
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	50
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	45
5			2,4-Nonadiyne	120	C9H12	063621-15-8	45



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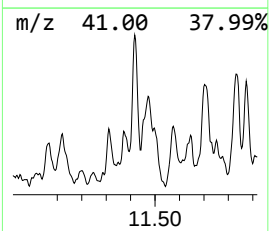
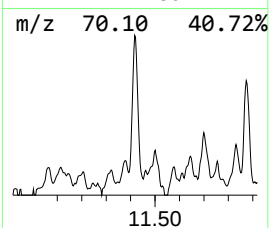
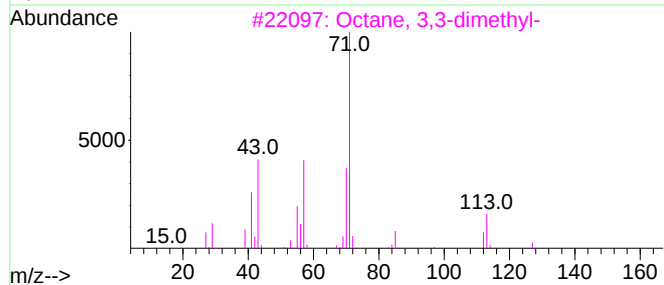
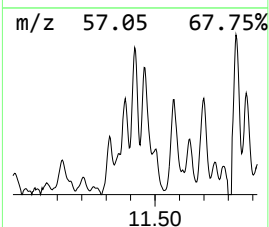
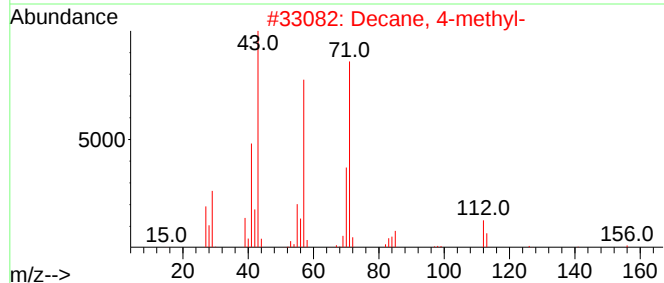
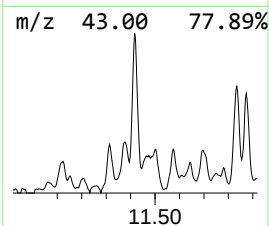
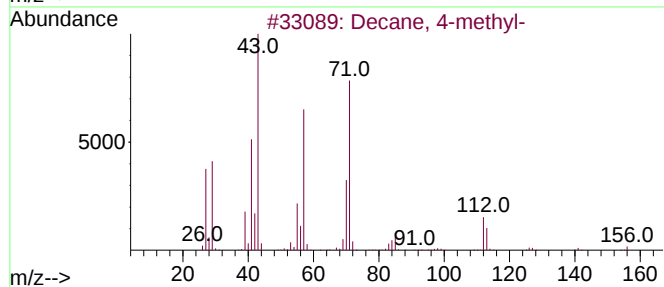
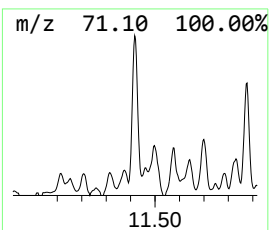
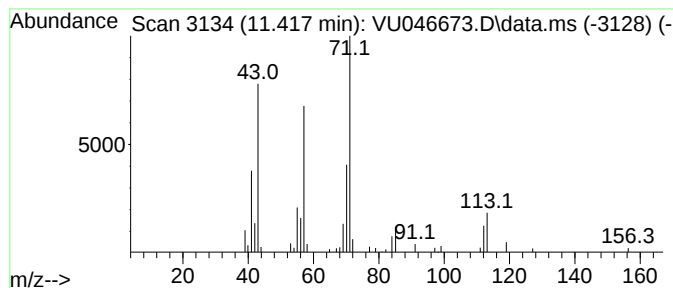
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ClientSampleId :
BGLG8

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Quant Title : VOC Analysis

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.417	3.24 ug/L	49981	1,4-Dichlorobenzene-d4		11.819	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane, 4-methyl-		156	C11H24	002847-72-5	83
2	Decane, 4-methyl-		156	C11H24	002847-72-5	83
3	Octane, 3,3-dimethyl-		142	C10H22	004110-44-5	78
4	Octane, 3,3-dimethyl-		142	C10H22	004110-44-5	72
5	4-Methyl-3-heptanol, trifluoroac...		226	C10H17F3O2	1000365-51-8	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU010522\
Data File : VU046673.D
Acq On : 05 Jan 2022 19:29
Operator : SY/MD
Sample : N1025-04
Misc : 6.80g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGLG8

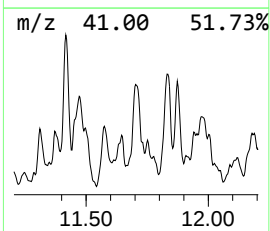
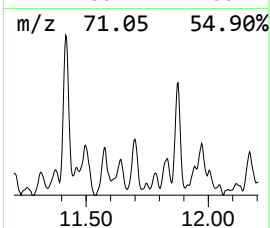
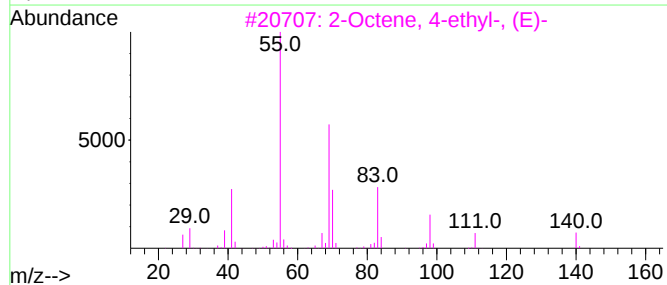
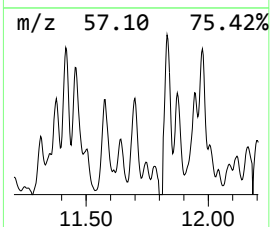
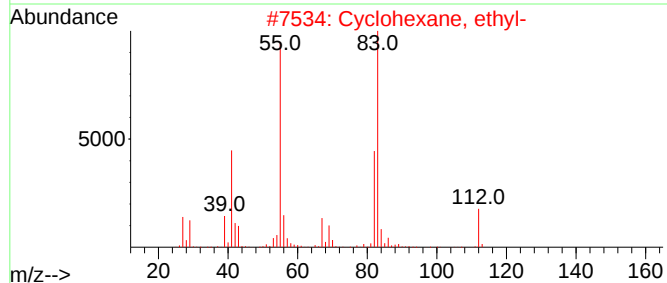
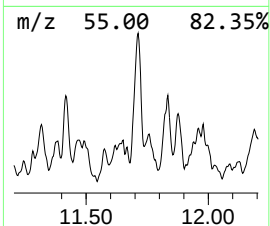
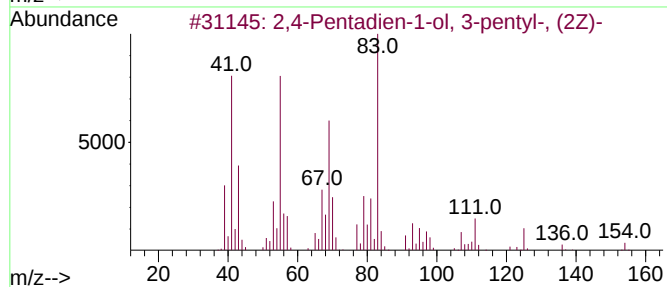
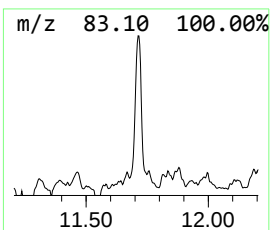
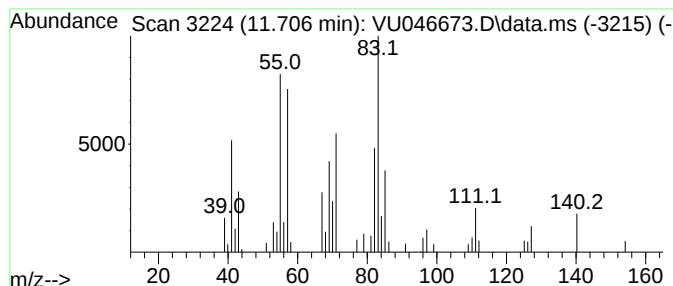
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Quant Title : VOC Analysis

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

Peak Number 5 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.706	2.70 ug/L	41635	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4-Pentadien-1-ol, 3-pentyl-, (...	154	C10H18O	1010142-19-7	45
2			Cyclohexane, ethyl-	112	C8H16	001678-91-7	43
3			2-Octene, 4-ethyl-, (E)-	140	C10H20	074630-09-4	43
4			Cyclohexane, butyl-	140	C10H20	001678-93-9	43
5			Cyclohexane, ethyl-	112	C8H16	001678-91-7	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU010522\
Data File : VU046673.D
Acq On : 05 Jan 2022 19:29
Operator : SY/MD
Sample : N1025-04
Misc : 6.80g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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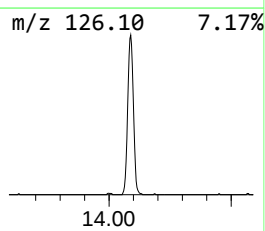
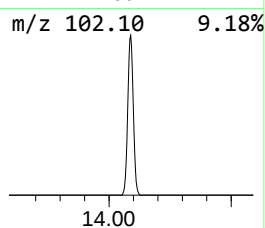
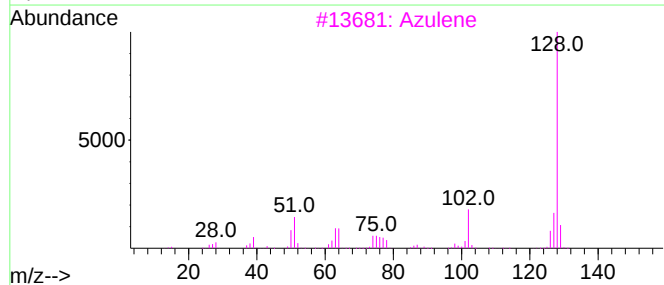
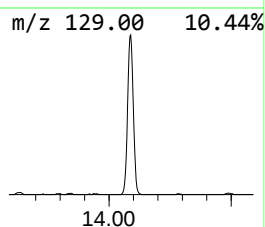
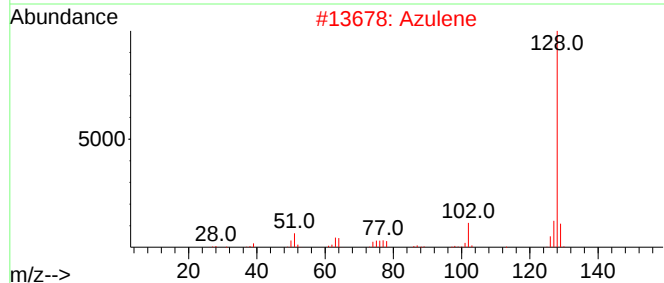
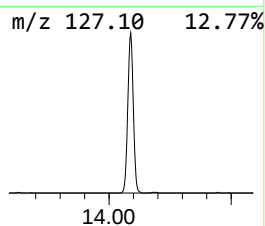
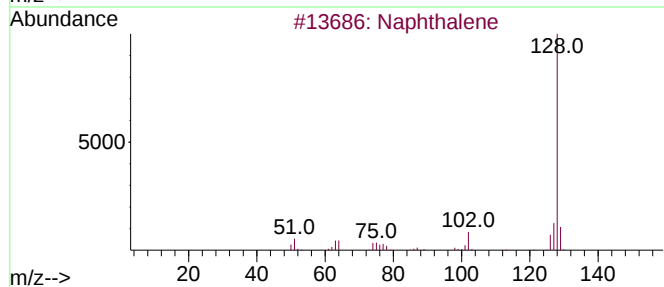
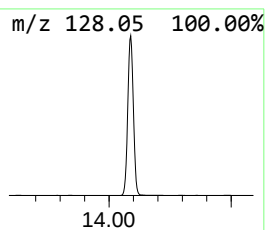
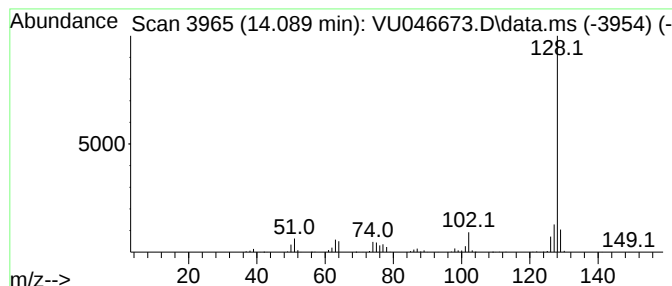
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Quant Title : VOC Analysis

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

Peak Number 6 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.089	50.93 ug/L	784403	1,4-Dichlorobenzene-d4	11.819

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU010522\
Data File : VU046673.D
Acq On : 05 Jan 2022 19:29
Operator : SY/MD
Sample : N1025-04
Misc : 6.80g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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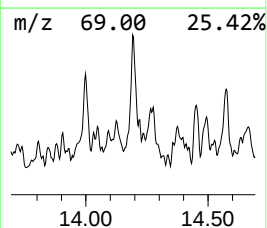
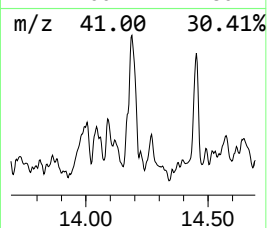
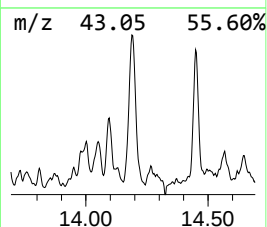
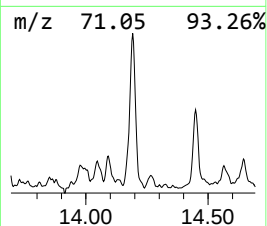
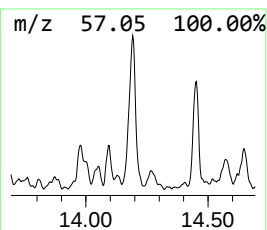
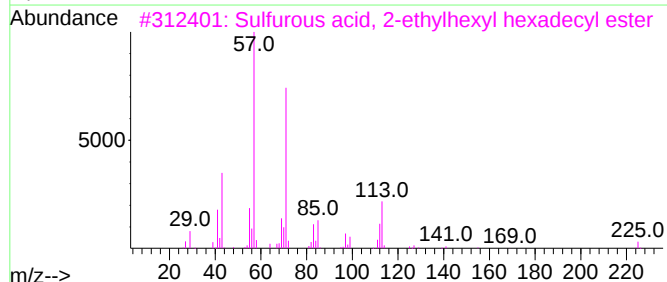
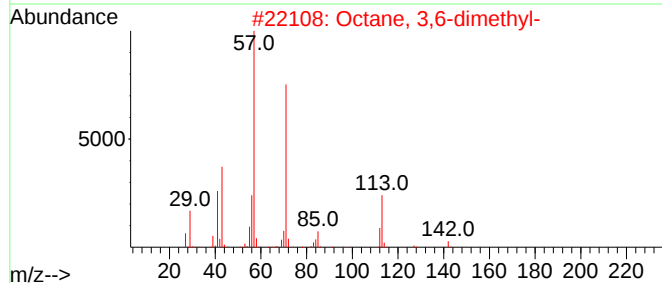
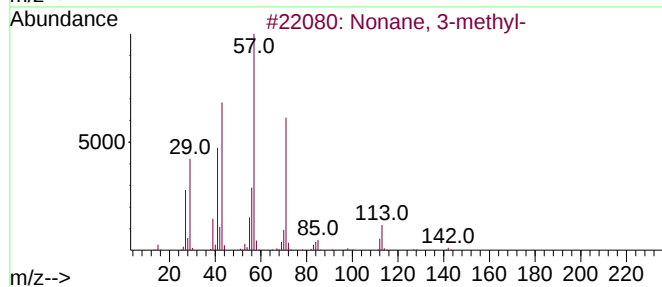
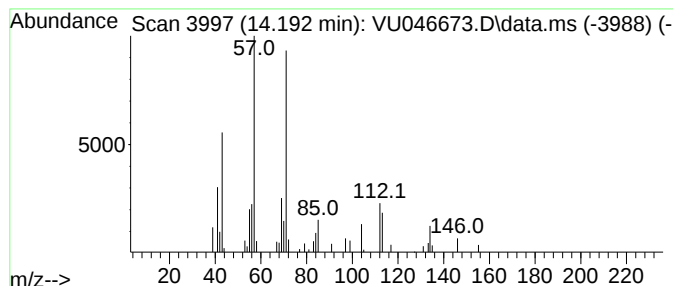
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Quant Title : VOC Analysis

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.192	2.90 ug/L	44680	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3-methyl-	142	C10H22	005911-04-6	59
2			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	59
3			Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	53
4			Heptane, 3-ethyl-5-methyl-	142	C10H22	052896-90-9	53
5			Octane, 2-bromo-	192	C8H17Br	000557-35-7	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU010522\
Data File : VU046673.D
Acq On : 05 Jan 2022 19:29
Operator : SY/MD
Sample : N1025-04
Misc : 6.80g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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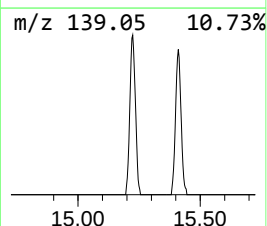
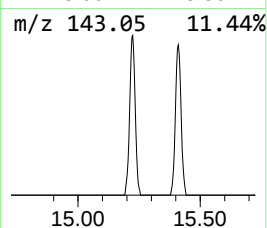
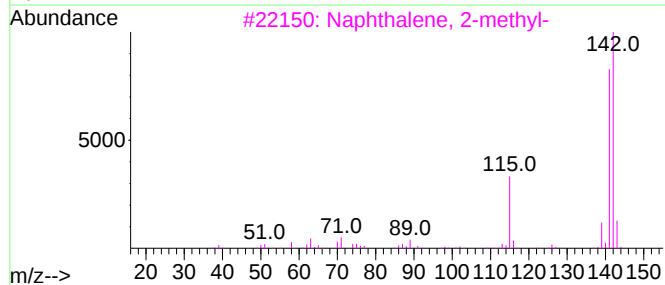
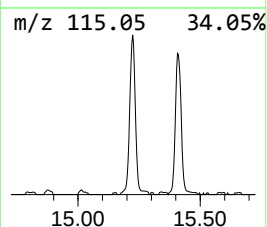
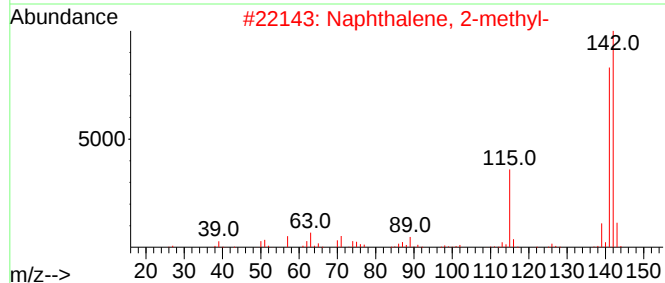
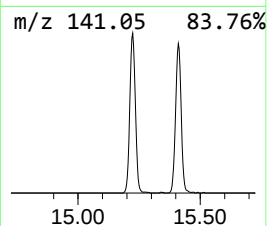
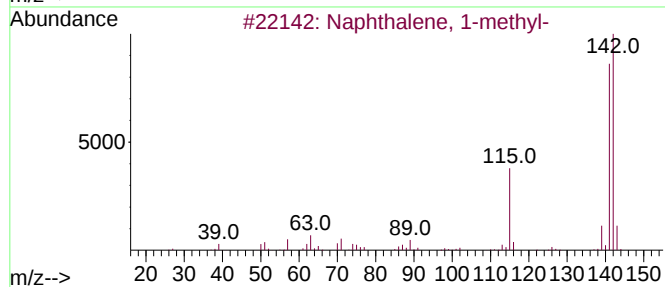
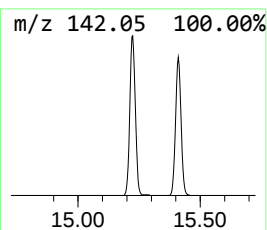
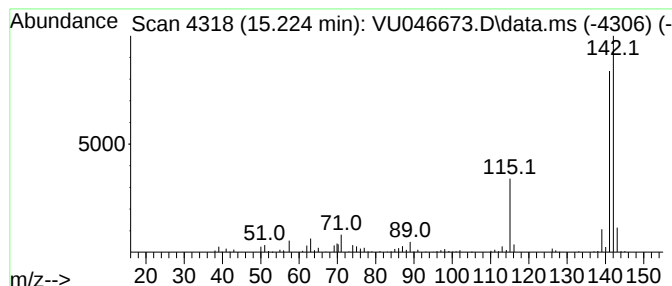
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Quant Title : VOC Analysis

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

Peak Number 8 Naphthalene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.224	11.33 ug/L	174543	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
5			Benzocycloheptatriene	142	C11H10	000264-09-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU010522\
Data File : VU046673.D
Acq On : 05 Jan 2022 19:29
Operator : SY/MD
Sample : N1025-04
Misc : 6.80g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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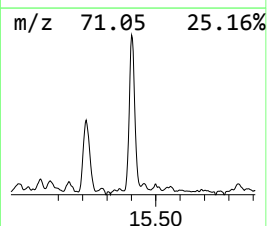
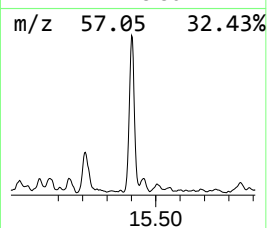
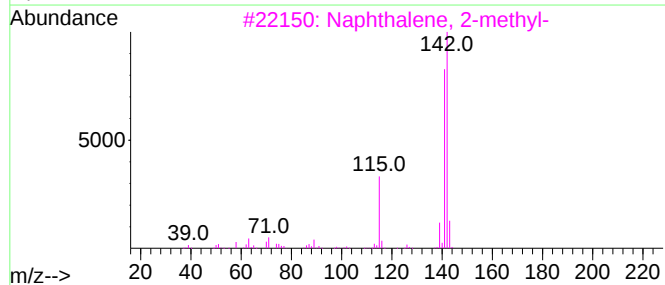
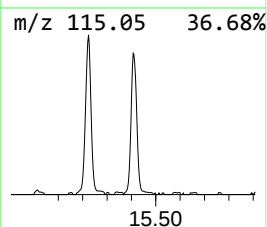
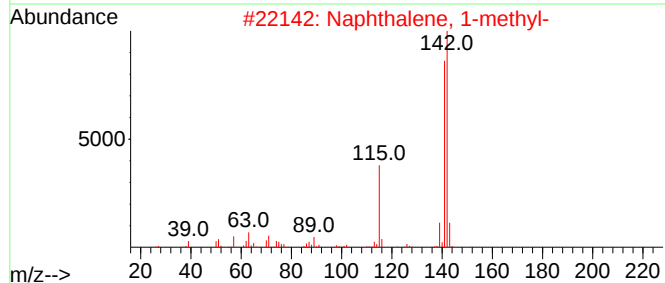
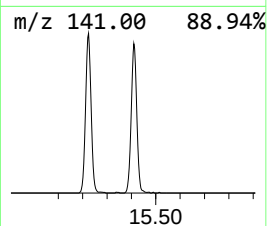
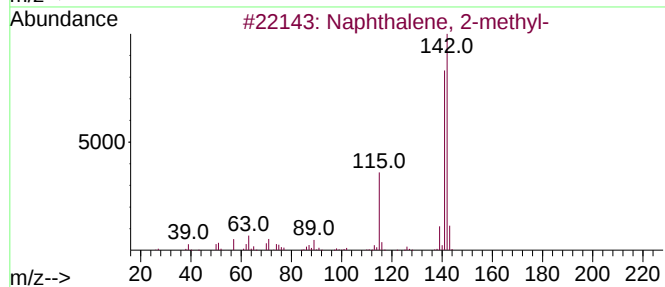
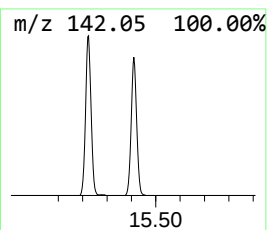
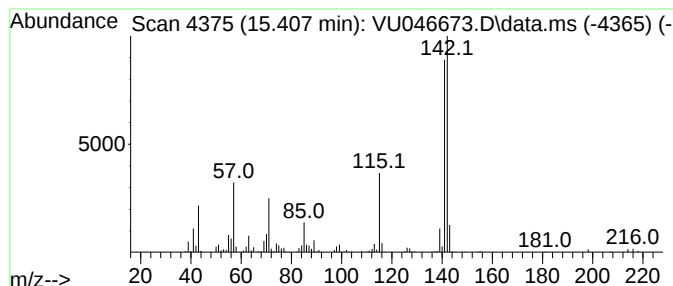
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Quant Title : VOC Analysis

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

Peak Number 9 Naphthalene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.407	14.54 ug/L	223993	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
4			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	93
5			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU010522\
 Data File : VU046673.D
 Acq On : 05 Jan 2022 19:29
 Operator : SY/MD
 Sample : N1025-04
 Misc : 6.80g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BGLG8

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM010522WMA.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
Benzene, 1-ethy...	11.028	2.8	ug/L	42670	3	11.819	770154	50.0
Benzene, 1-ethy...	11.307	2.8	ug/L	42449	3	11.819	770154	50.0
(DEL) Alkane: S...	11.417	3.2	ug/L	49981	3	11.819	770154	50.0
unknown-01	11.706	2.7	ug/L	41635	3	11.819	770154	50.0
Azulene	14.089	50.9	ug/L	784403	3	11.819	770154	50.0
(DEL) Alkane: S...	14.192	2.9	ug/L	44680	3	11.819	770154	50.0
Naphthalene, 1-...	15.224	11.3	ug/L	174543	3	11.819	770154	50.0
Naphthalene, 2-...	15.407	14.5	ug/L	223993	3	11.819	770154	50.0