

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
 Data File : VU046113.D
 Acq On : 06 Dec 2021 15:42
 Operator : SY/MD
 Sample : M4942-03DL 100X
 Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BGKQ1DL

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\SFAMULM112921WMA.M
 Title : VOC Analysis

Signal : TIC: VU046113.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.601	73	81	96	rBV	99552	110009	12.58%	1.001%
2	1.916	171	179	191	rVB	70212	87624	10.02%	0.797%
3	2.514	358	365	370	rVB2	486	533	0.06%	0.005%
4	2.572	370	383	406	rBV	172893	289478	33.11%	2.634%
5	2.707	422	425	429	rBV2	208	195	0.02%	0.002%
6	2.790	444	451	463	rVB3	376	826	0.09%	0.008%
7	2.961	501	504	505	rBV	149	106	0.01%	0.001%
8	3.047	525	531	539	rBV3	1449	2129	0.24%	0.019%
9	3.504	669	673	675	rBV	165	122	0.01%	0.001%
10	3.684	726	729	732	rBV	160	165	0.02%	0.002%
11	4.035	834	838	841	rVB2	166	131	0.01%	0.001%
12	4.076	849	851	856	rVB	159	107	0.01%	0.001%
13	4.154	870	875	878	rVB2	212	184	0.02%	0.002%
14	4.633	1007	1024	1053	rBV	76679	196975	22.53%	1.793%
15	4.832	1084	1086	1092	rVB	294	217	0.02%	0.002%
16	4.864	1092	1096	1098	rVV	317	255	0.03%	0.002%
17	4.896	1102	1106	1112	rVB2	293	220	0.03%	0.002%
18	5.067	1142	1159	1181	rBV	138035	313508	35.86%	2.853%
19	5.231	1209	1210	1216	rVB	398	252	0.03%	0.002%
20	5.263	1216	1220	1224	rBV	183	252	0.03%	0.002%
21	5.295	1224	1230	1235	rVB2	389	510	0.06%	0.005%
22	5.337	1238	1243	1249	rVB	251	332	0.04%	0.003%
23	5.385	1254	1258	1260	rBV	166	177	0.02%	0.002%
24	5.472	1282	1285	1290	rVB	177	132	0.02%	0.001%
25	5.729	1343	1365	1387	rBV2	299933	804791	92.06%	7.324%
26	6.028	1454	1458	1459	rBV	195	152	0.02%	0.001%
27	6.250	1512	1527	1546	rBV	230023	431990	49.41%	3.931%
28	6.452	1586	1590	1593	rBV2	180	144	0.02%	0.001%
29	6.530	1601	1614	1630	rVB5	4263	9019	1.03%	0.082%
30	6.694	1651	1665	1682	rBV	180854	352466	40.32%	3.208%
31	6.803	1698	1699	1700	rBV	407	107	0.01%	0.001%
32	6.867	1715	1719	1723	rBV	235	203	0.02%	0.002%
33	7.047	1767	1775	1782	rBV3	527	904	0.10%	0.008%
34	7.176	1812	1815	1818	rVV2	248	252	0.03%	0.002%
35	7.234	1829	1833	1836	rVV2	174	145	0.02%	0.001%
36	7.385	1876	1880	1883	rBV2	194	155	0.02%	0.001%

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Integrator: RTE

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Peak Location: TOP

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

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
 Title : VOC Analysis

37	7.459	1901	1903	1906	rVB	241	144	0.02%	0.001%
38	7.501	1911	1916	1923	rVB3	615	679	0.08%	0.006%
39	7.568	1924	1937	1951	rBV	107547	185476	21.22%	1.688%
40	7.661	1960	1966	1969	rBV3	820	809	0.09%	0.007%
41	7.771	1996	2000	2002	rBV3	203	128	0.01%	0.001%
42	7.813	2008	2013	2016	rBV	112	127	0.01%	0.001%
43	7.848	2020	2024	2027	rVV3	571	562	0.06%	0.005%
44	7.899	2027	2040	2054	rVV	387071	670313	76.67%	6.100%
45	7.970	2057	2062	2079	rVB	19118	31190	3.57%	0.284%
46	8.063	2088	2091	2093	rBV2	297	171	0.02%	0.002%
47	8.134	2104	2113	2117	rBV4	1337	1822	0.21%	0.017%
48	8.182	2117	2128	2140	rVV	70231	116074	13.28%	1.056%
49	8.372	2183	2187	2194	rBV4	622	624	0.07%	0.006%
50	8.468	2214	2217	2221	rBV2	465	348	0.04%	0.003%
51	8.539	2233	2239	2243	rBV2	654	574	0.07%	0.005%
52	8.632	2256	2268	2294	rBV	240246	421025	48.16%	3.831%
53	8.867	2334	2341	2351	rVB3	2731	3957	0.45%	0.036%
54	8.925	2351	2359	2366	rBV5	3389	5631	0.64%	0.051%
55	9.038	2385	2394	2395	rBV3	833	984	0.11%	0.009%
56	9.128	2412	2422	2426	rBV6	2824	4834	0.55%	0.044%
57	9.166	2430	2434	2435	rBV2	2443	1777	0.20%	0.016%
58	9.214	2442	2449	2460	rVB4	11666	19676	2.25%	0.179%
59	9.263	2460	2464	2470	rVB4	536	586	0.07%	0.005%
60	9.301	2472	2476	2477	rBV	172	112	0.01%	0.001%
61	9.337	2477	2487	2499	rBV3	11729	19988	2.29%	0.182%
62	9.417	2499	2512	2530	rBV	325670	546305	62.49%	4.972%
63	9.501	2536	2538	2540	rBV2	404	226	0.03%	0.002%
64	9.571	2549	2560	2573	rBV	75631	119557	13.68%	1.088%
65	9.694	2586	2598	2608	rBV	217644	375133	42.91%	3.414%
66	9.954	2672	2679	2688	rBV6	2056	3872	0.44%	0.035%
67	10.021	2689	2700	2713	rBV5	18505	44402	5.08%	0.404%
68	10.102	2716	2725	2740	rVB	71996	130883	14.97%	1.191%
69	10.253	2754	2772	2785	rBV4	38547	80388	9.20%	0.732%
70	10.314	2786	2791	2793	rBV4	4591	4572	0.52%	0.042%
71	10.356	2796	2804	2811	rBV5	26055	43646	4.99%	0.397%
72	10.484	2831	2844	2852	rVB2	19631	38633	4.42%	0.352%
73	10.558	2852	2867	2875	rBV4	28168	60776	6.95%	0.553%
74	10.758	2918	2929	2939	rBV	295281	471131	53.89%	4.287%
75	10.906	2967	2975	2984	rBV	30815	46219	5.29%	0.421%
76	10.999	2996	3004	3019	rBV	83327	193306	22.11%	1.759%

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
 Title : VOC Analysis

77	11.115	3035	3040	3055	rVB	246326	370506	42.38%	3.372%
78	11.247	3075	3081	3086	rBV3	6177	9051	1.04%	0.082%
79	11.298	3087	3097	3113	rVV3	50940	132050	15.10%	1.202%
80	11.375	3116	3121	3127	rVV3	27880	40740	4.66%	0.371%
81	11.417	3127	3134	3141	rVV	89240	129340	14.79%	1.177%
82	11.468	3141	3150	3174	rVB	268967	490976	56.16%	4.468%
83	11.574	3175	3183	3189	rBV2	31327	52014	5.95%	0.473%
84	11.706	3215	3224	3231	rBV5	54265	101960	11.66%	0.928%
85	11.812	3247	3257	3270	rBV	384923	652223	74.60%	5.935%
86	11.906	3271	3286	3294	rBV3	78806	207946	23.79%	1.892%
87	12.099	3337	3346	3349	rBV2	41623	63469	7.26%	0.578%
88	12.192	3364	3375	3390	rVB	493420	874239	100.00%	7.956%
89	12.327	3408	3417	3427	rBV	309845	432774	49.50%	3.938%
90	12.468	3455	3461	3464	rBV	32230	41252	4.72%	0.375%
91	13.353	3730	3736	3745	rBV3	9443	14130	1.62%	0.129%
92	13.433	3753	3761	3781	rVB4	108707	218592	25.00%	1.989%
93	13.735	3848	3855	3864	rBV10	4839	9813	1.12%	0.089%
94	13.838	3882	3887	3903	rVB6	10836	19374	2.22%	0.176%
95	14.086	3955	3964	3984	rVB	256736	422574	48.34%	3.846%
96	14.192	3987	3997	4012	rVB4	22609	47926	5.48%	0.436%
97	14.452	4070	4078	4086	rBV	62714	85681	9.80%	0.780%
98	15.224	4305	4318	4328	rBV	88446	152403	17.43%	1.387%
99	15.404	4366	4374	4385	rBV2	82893	128034	14.65%	1.165%
100	16.262	4634	4641	4646	rBV2	25464	40053	4.58%	0.364%

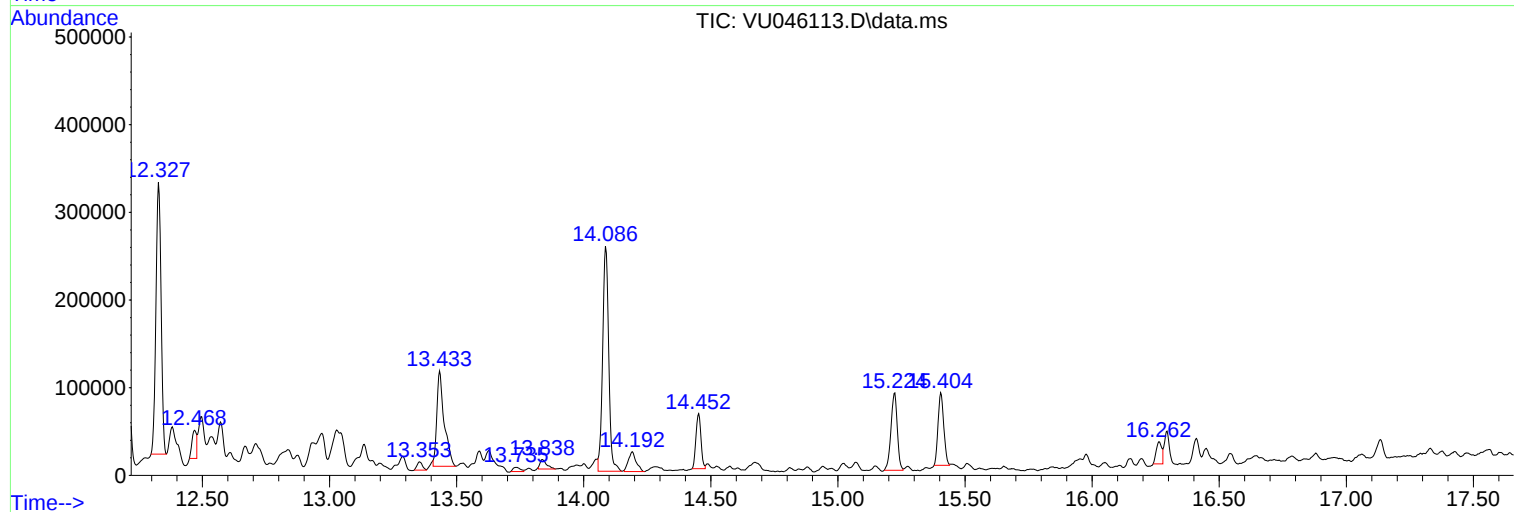
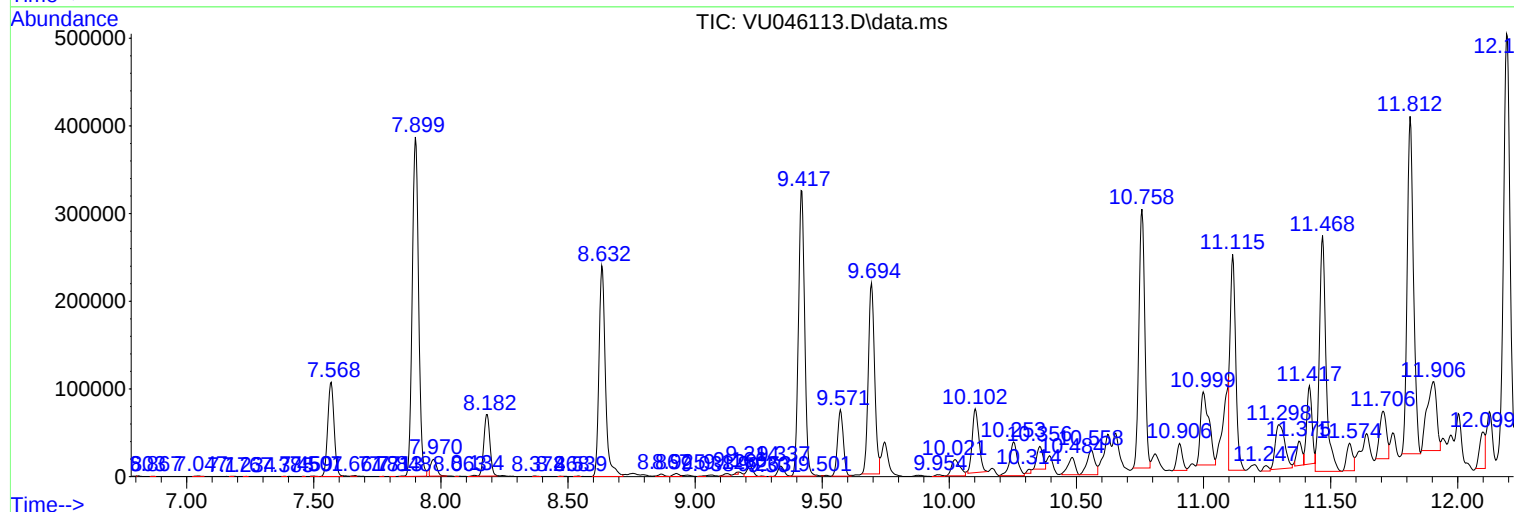
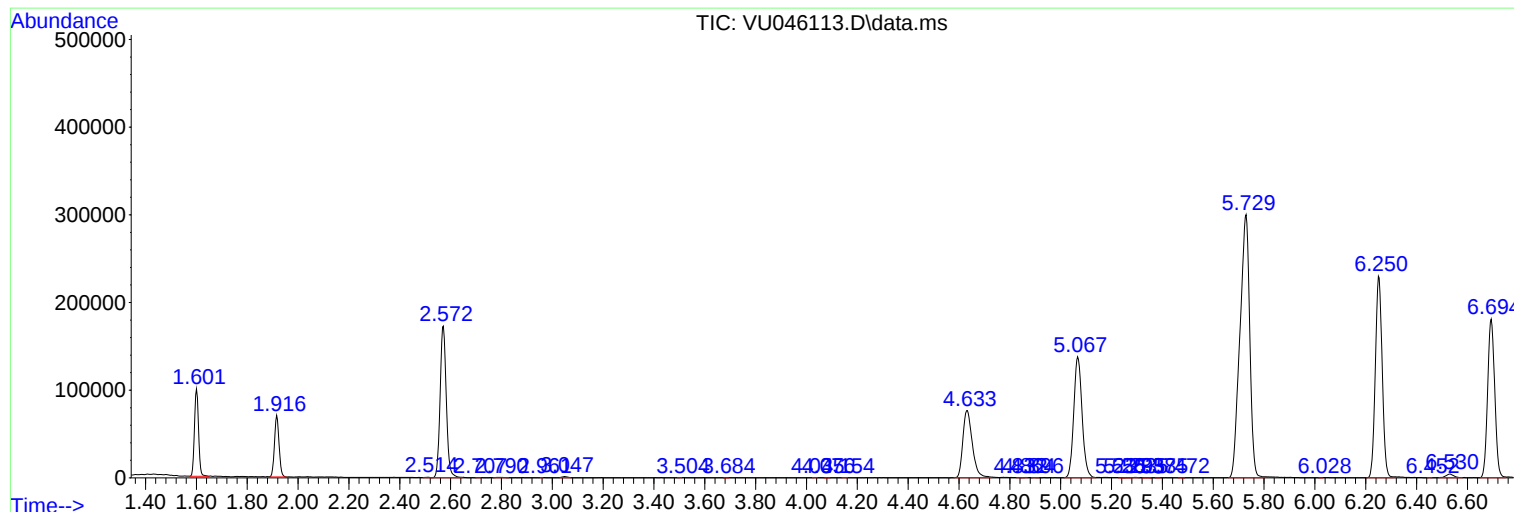
Sum of corrected areas: 10988547

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ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKQ1DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.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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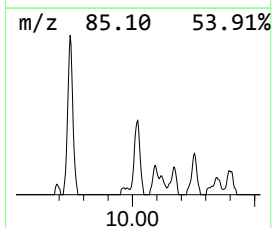
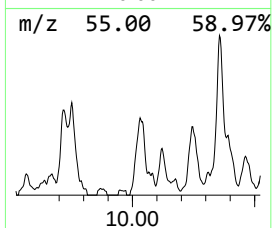
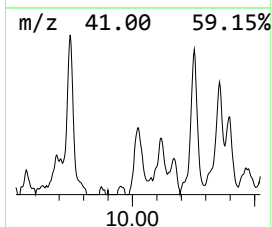
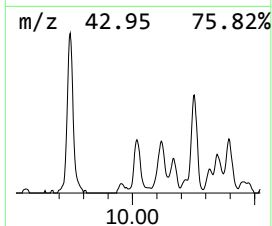
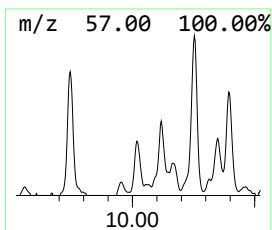
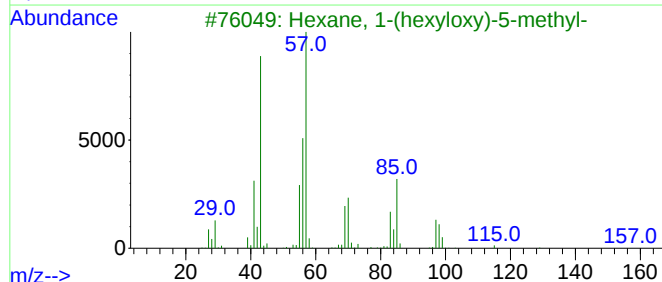
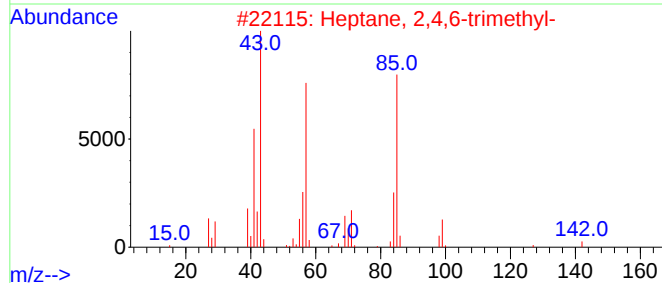
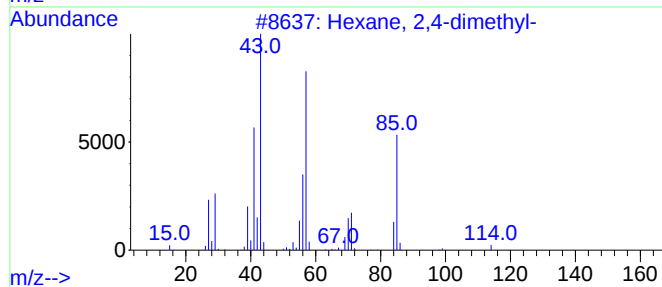
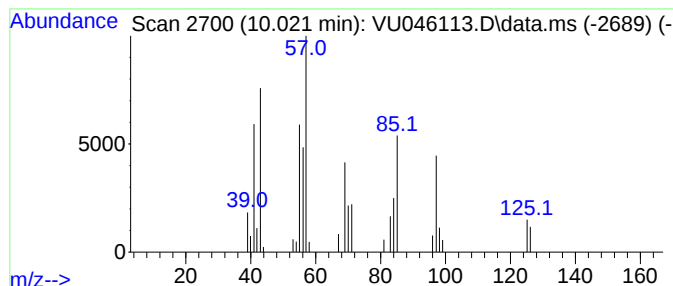
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.022	4.06 ug/L	44402	Chlorobenzene-d5	9.420

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	52
2			Heptane, 2,4,6-trimethyl-	142	C10H22	002613-61-8	43
3			Hexane, 1-(hexyloxy)-5-methyl-	200	C13H28O	074421-19-5	43
4			1-Heptanol, 6-methyl-	130	C8H18O	001653-40-3	27
5			Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	27



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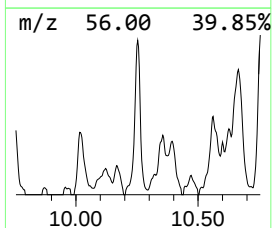
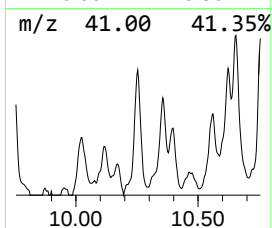
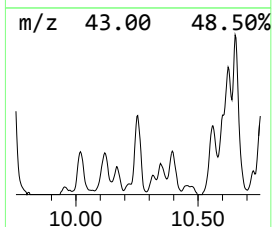
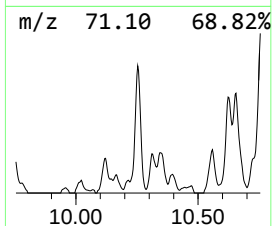
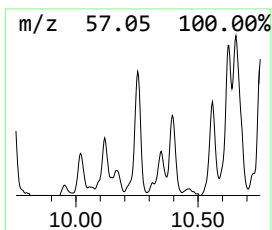
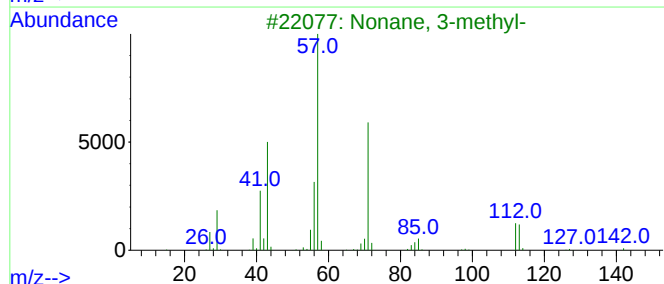
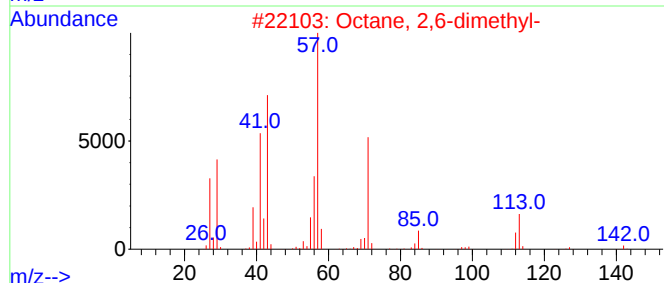
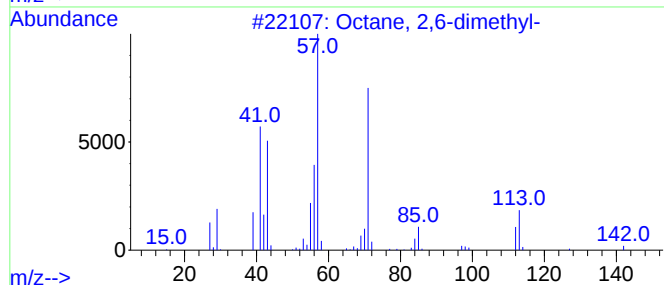
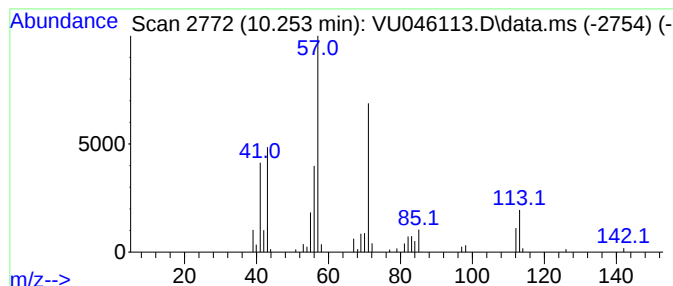
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.253	7.36 ug/L	80388	Chlorobenzene-d5	9.420

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-			142	C10H22	002051-30-1	90
2	Octane, 2,6-dimethyl-			142	C10H22	002051-30-1	90
3	Nonane, 3-methyl-			142	C10H22	005911-04-6	87
4	Nonane, 3-methyl-			142	C10H22	005911-04-6	83
5	Octane, 2,6-dimethyl-			142	C10H22	002051-30-1	81



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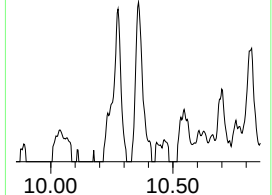
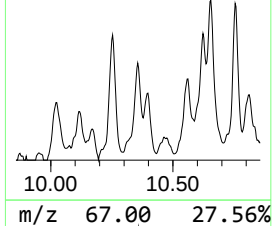
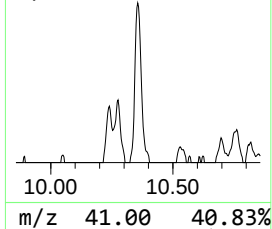
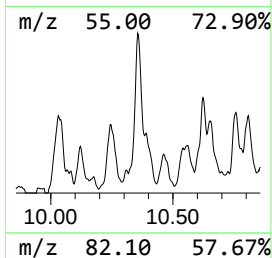
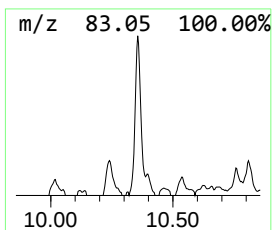
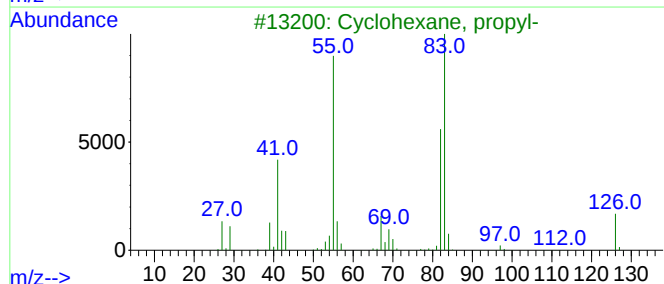
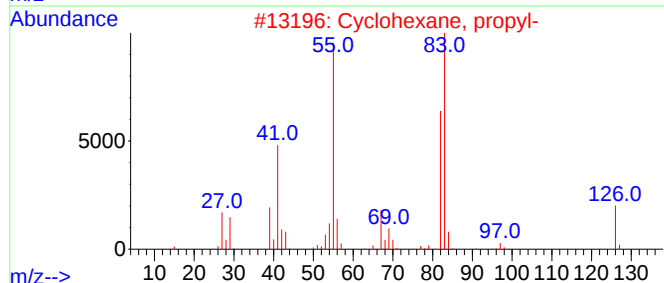
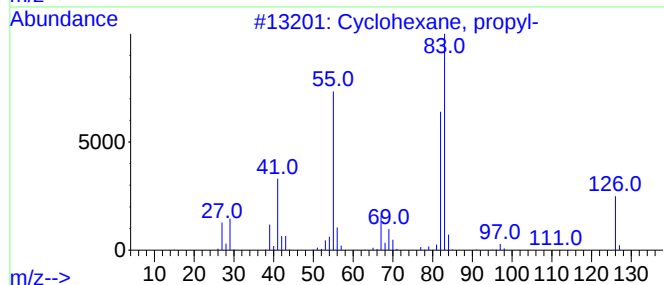
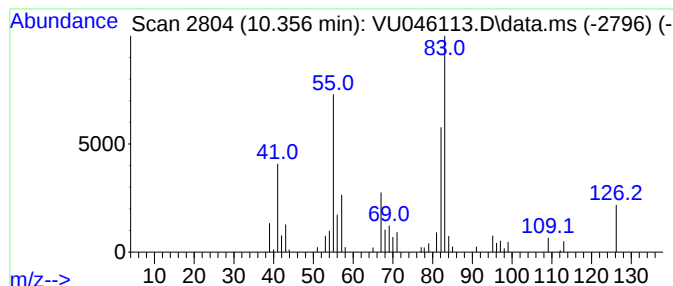
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Quant Title : VOC Analysis

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

Peak Number 4 (DEL) Alkane: Cyclic10.356 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.356	3.99 ug/L	43646	Chlorobenzene-d5	9.420

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, propyl-	126	C9H18	001678-92-8	81
2			Cyclohexane, propyl-	126	C9H18	001678-92-8	81
3			Cyclohexane, propyl-	126	C9H18	001678-92-8	74
4			Cyclohexane, propyl-	126	C9H18	001678-92-8	72
5			Cyclohexane, decyl-	224	C16H32	001795-16-0	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKQ1DL

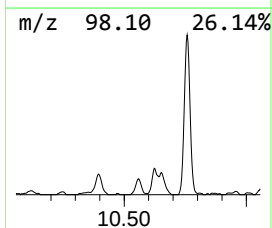
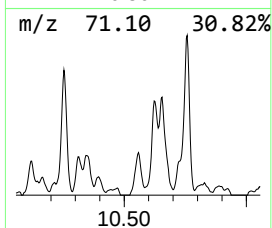
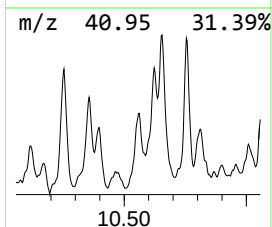
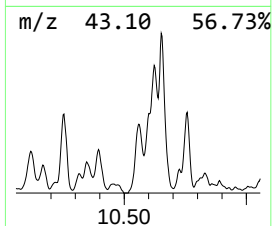
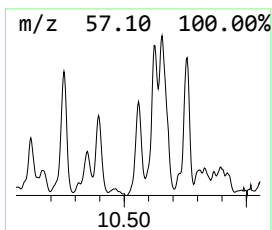
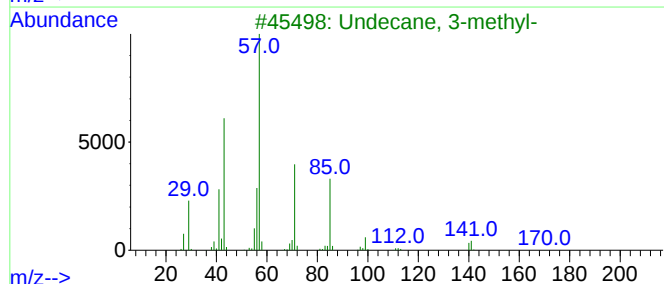
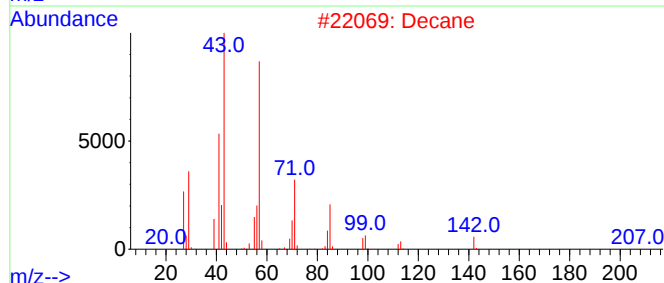
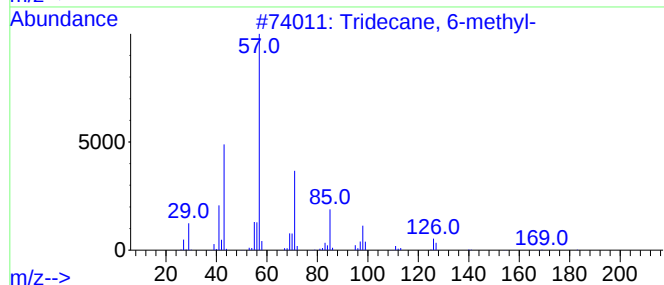
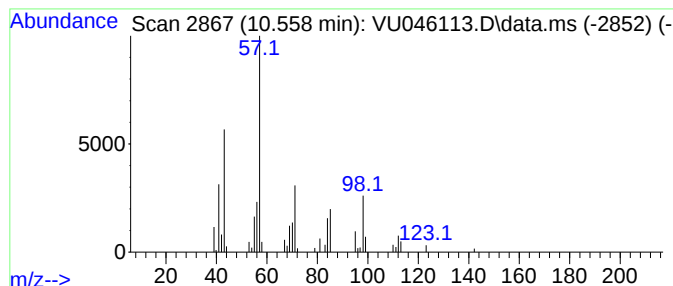
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Quant Title : VOC Analysis

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.558	5.56 ug/L	60776	Chlorobenzene-d5	9.420

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 6-methyl-	198	C14H30	013287-21-3	47
2			Decane	142	C10H22	000124-18-5	47
3			Undecane, 3-methyl-	170	C12H26	001002-43-3	47
4			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	47
5			Decane, 2,5,6-trimethyl-	184	C13H28	062108-23-0	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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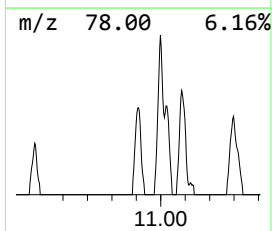
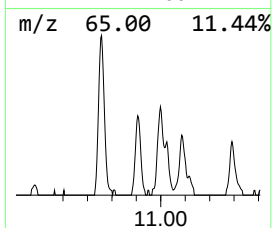
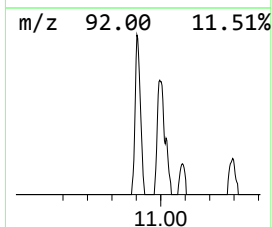
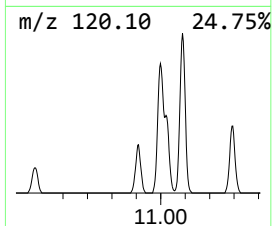
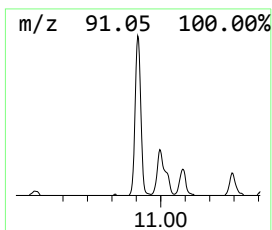
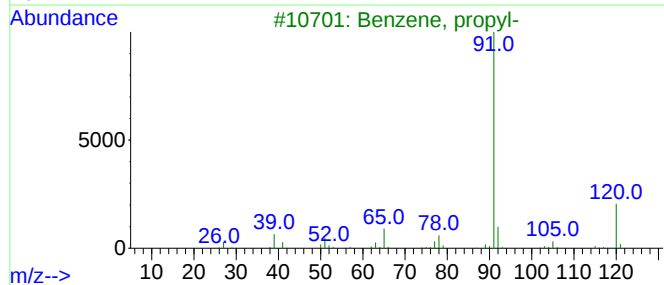
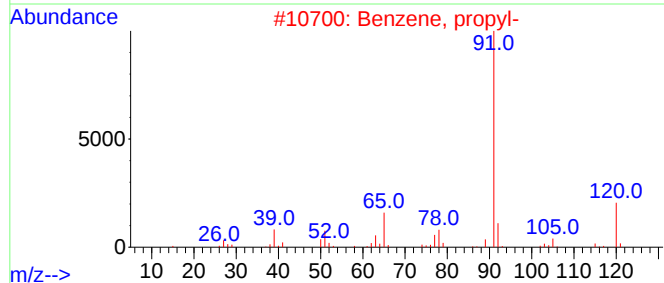
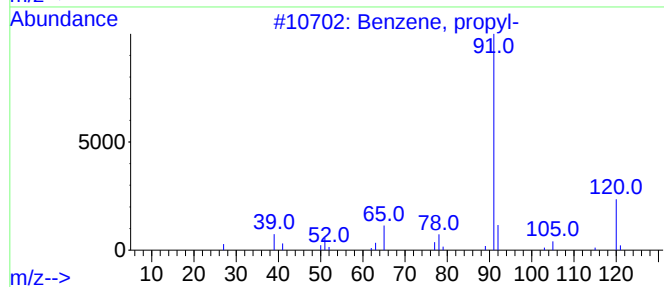
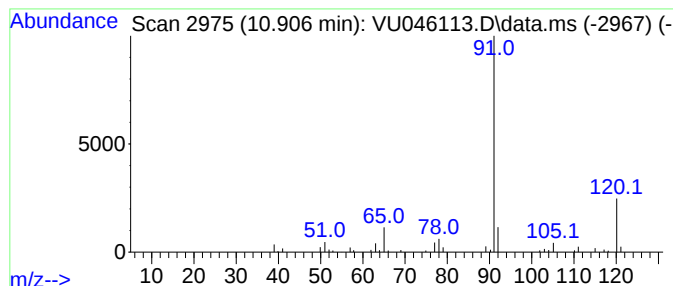
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Quant Title : VOC Analysis

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

Peak Number 6 Benzene, propyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.906	3.54 ug/L	46219	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	93
2			Benzene, propyl-	120	C9H12	000103-65-1	90
3			Benzene, propyl-	120	C9H12	000103-65-1	87
4			Benzene, propyl-	120	C9H12	000103-65-1	83
5			N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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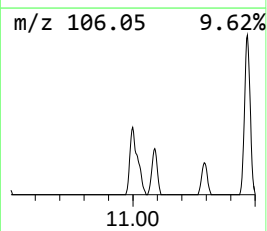
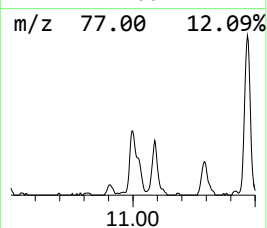
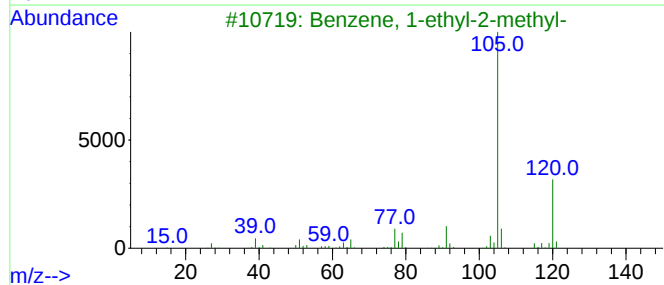
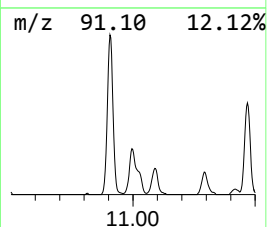
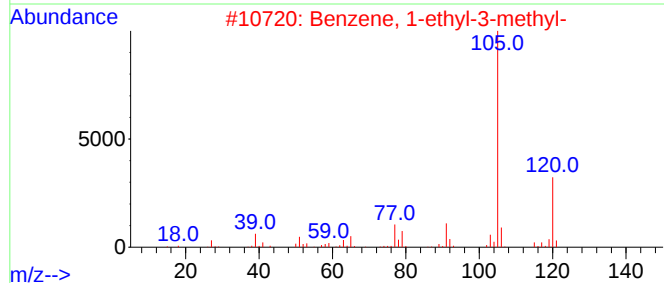
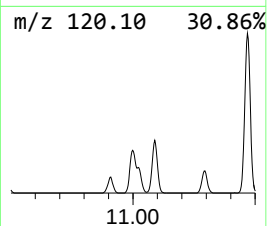
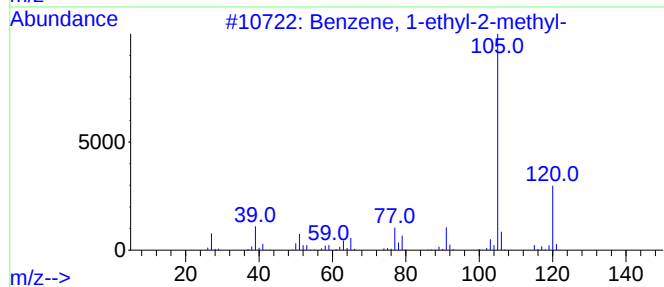
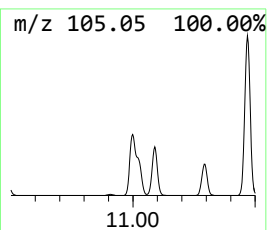
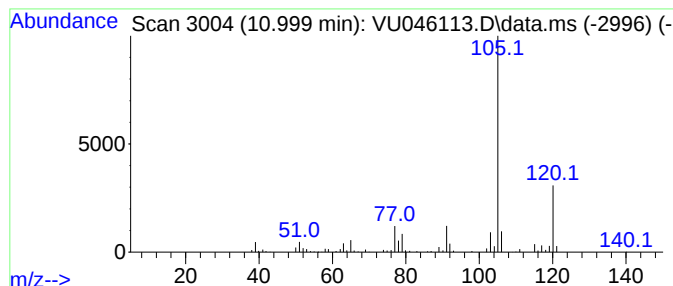
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Quant Title : VOC Analysis

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.999	14.82 ug/L	193306	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
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Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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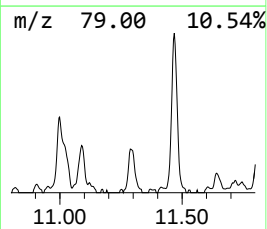
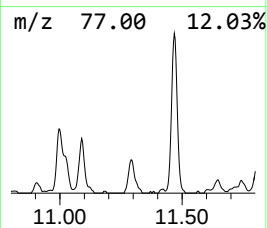
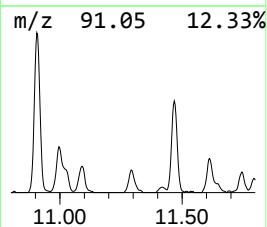
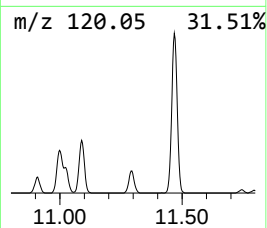
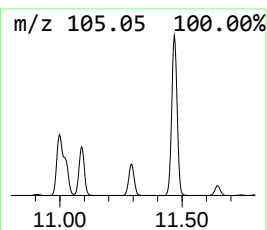
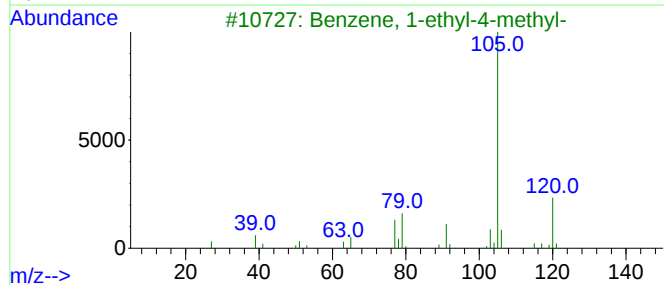
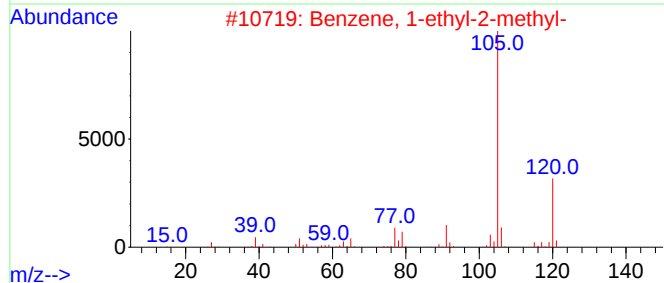
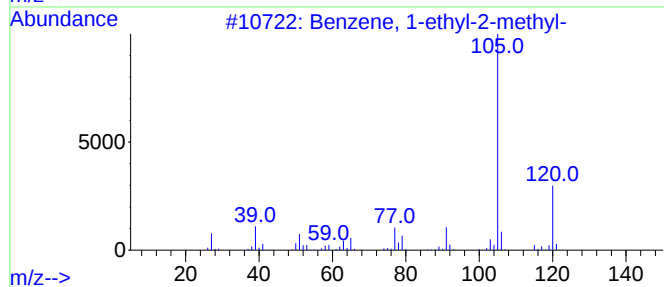
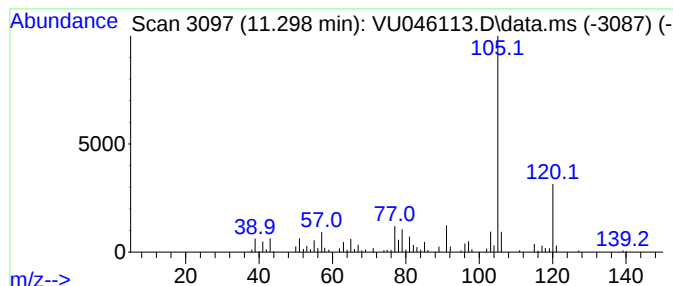
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Quant Title : VOC Analysis

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

Peak Number 8 Benzene, 1-ethyl-4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.298	10.12 ug/L	132050	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	89
4			Mesitylene	120	C9H12	000108-67-8	87
5			Mesitylene	120	C9H12	000108-67-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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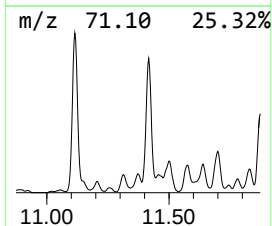
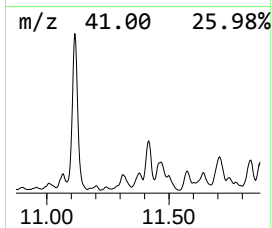
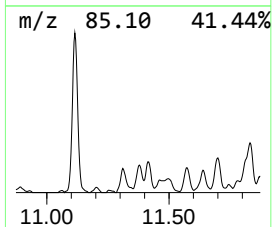
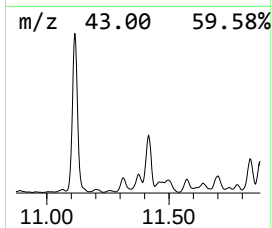
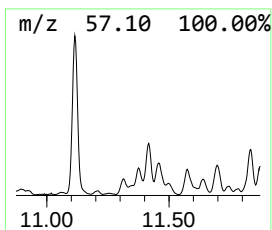
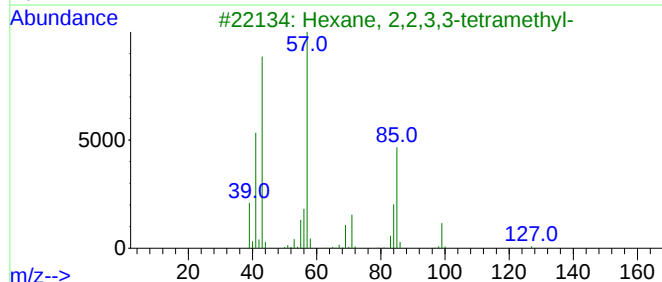
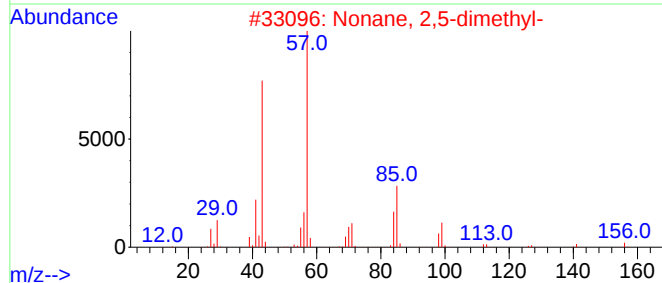
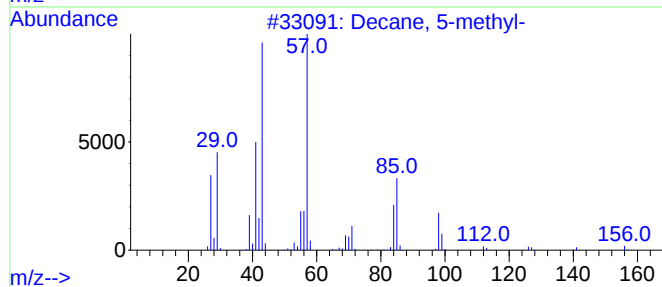
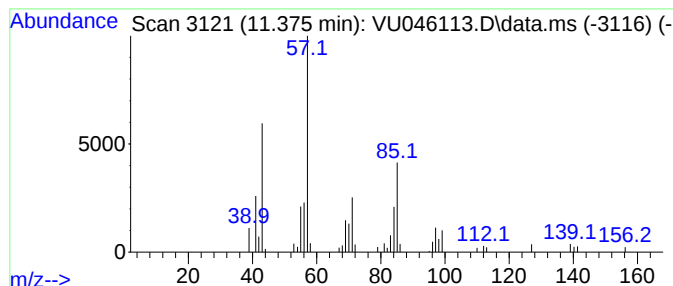
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Quant Title : VOC Analysis

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.375	3.12 ug/L	40740	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 5-methyl-	156	C11H24	013151-35-4	72
2			Nonane, 2,5-dimethyl-	156	C11H24	017302-27-1	62
3			Hexane, 2,2,3,3-tetramethyl-	142	C10H22	013475-81-5	59
4			Tritetracontane	605	C43H88	007098-21-7	53
5			Heptane, 2,2,3,3,5,6,6-heptamethyl-	198	C14H30	007225-67-4	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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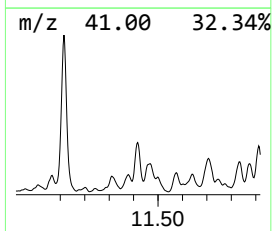
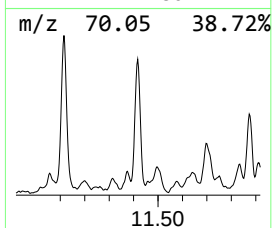
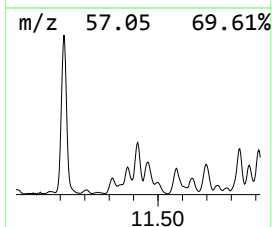
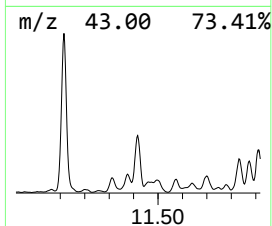
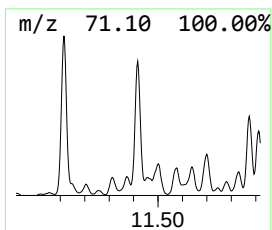
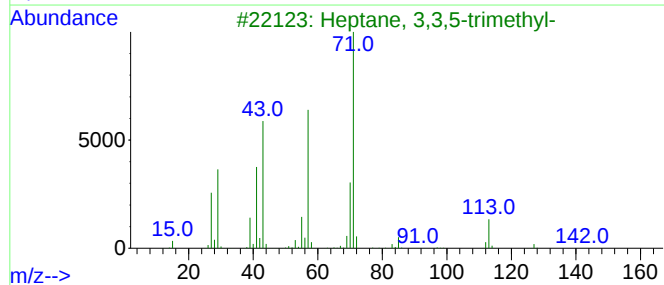
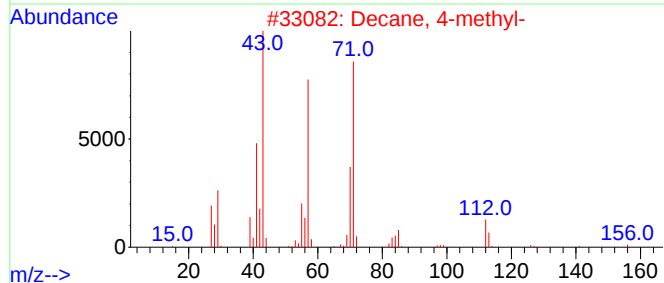
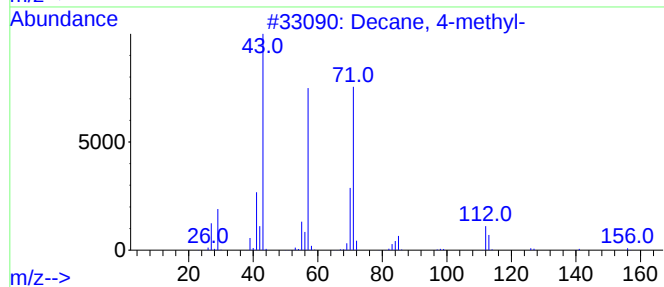
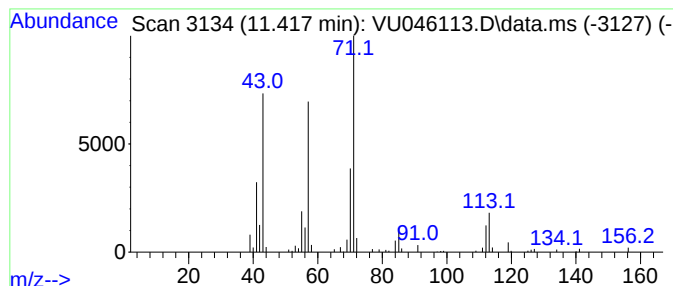
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Quant Title : VOC Analysis

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.417	9.92 ug/L	129340	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 4-methyl-	156	C11H24	002847-72-5	87
2			Decane, 4-methyl-	156	C11H24	002847-72-5	74
3			Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	64
4			Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	64
5			Decane, 2,8,8-trimethyl-	184	C13H28	1000060-81-2	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKQ1DL

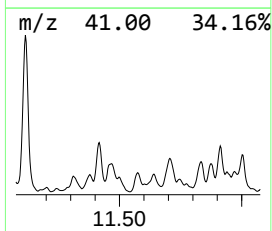
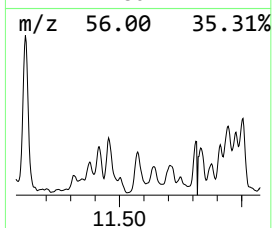
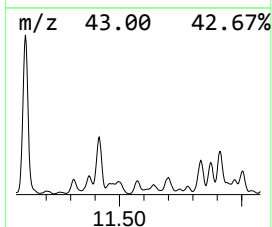
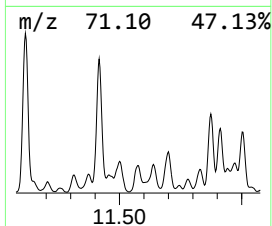
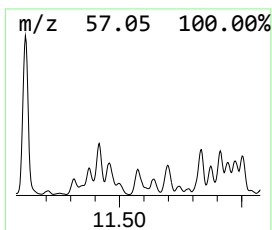
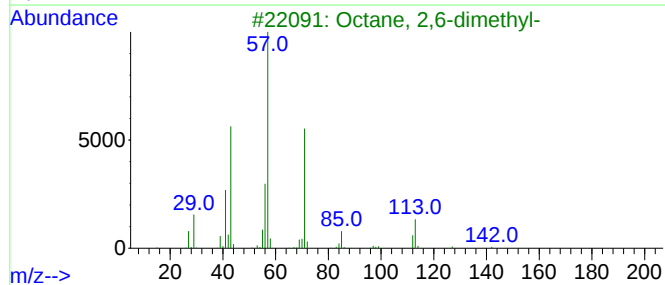
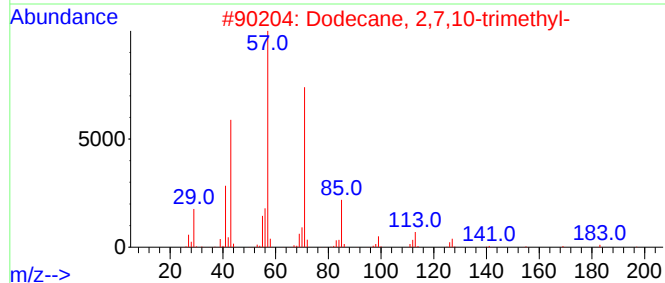
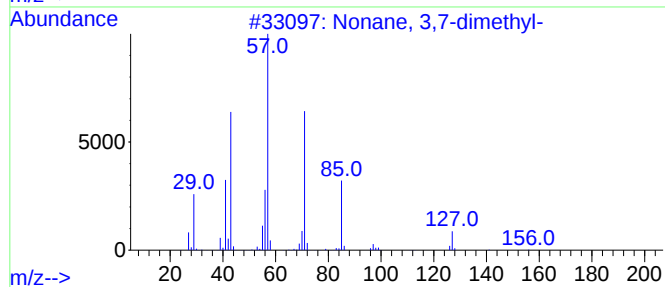
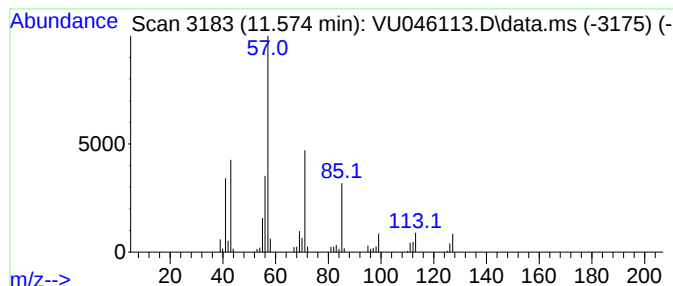
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Quant Title : VOC Analysis

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.575	3.99 ug/L	52014	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	59
2			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	53
3			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	53
4			Sulfurous acid, butyl octyl ester	250	C12H26O3S	1000309-17-5	50
5			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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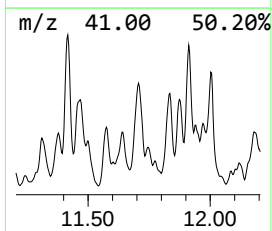
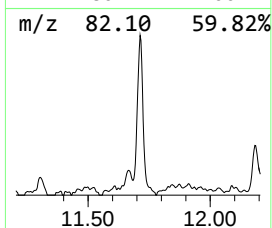
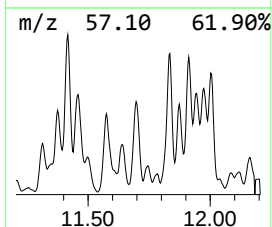
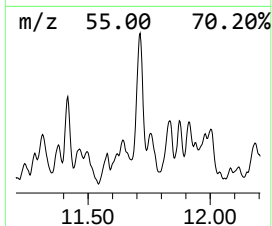
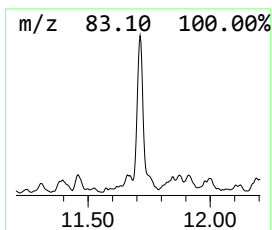
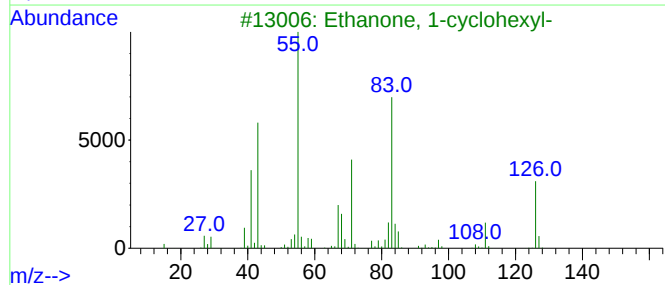
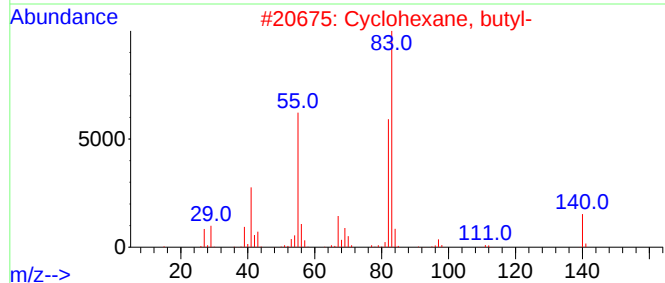
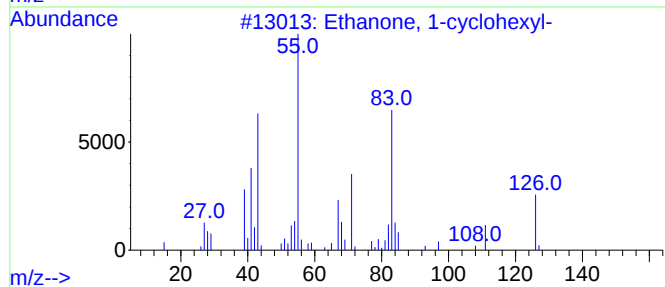
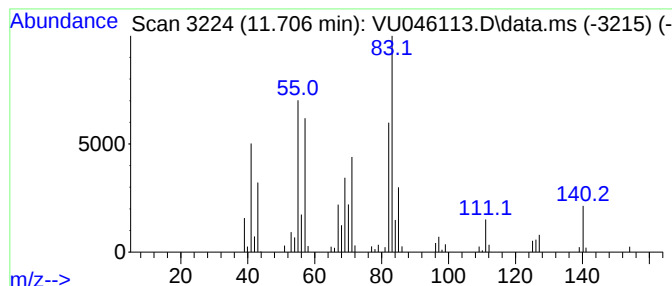
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Quant Title : VOC Analysis

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

Peak Number 12 Ethanone, 1-cyclohexyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.706	7.82 ug/L	101960	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanone, 1-cyclohexyl-	126	C8H14O	000823-76-7	64
2			Cyclohexane, butyl-	140	C10H20	001678-93-9	53
3			Ethanone, 1-cyclohexyl-	126	C8H14O	000823-76-7	52
4			Cyclohexane, pentyl-	154	C11H22	004292-92-6	49
5			Cyclohexane, butyl-	140	C10H20	001678-93-9	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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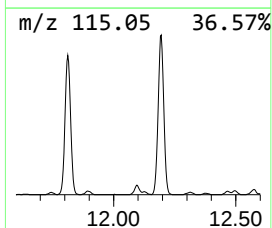
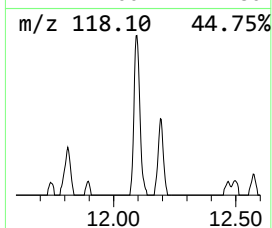
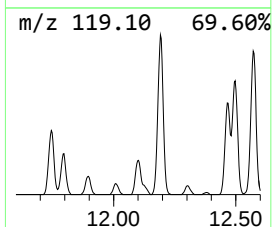
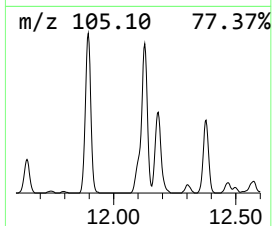
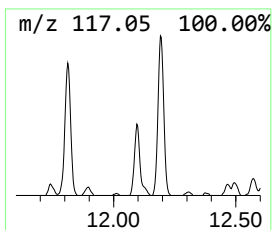
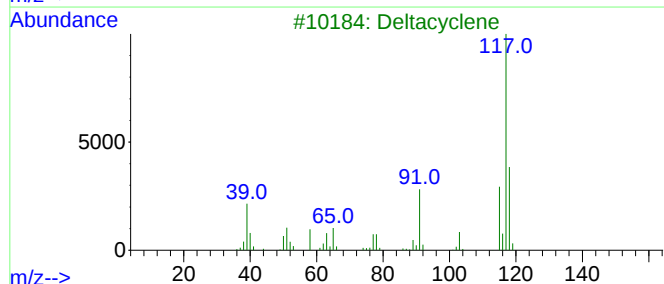
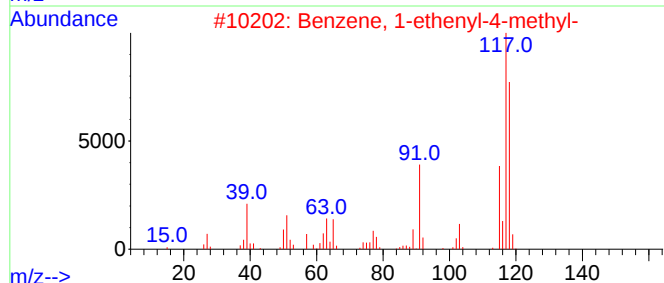
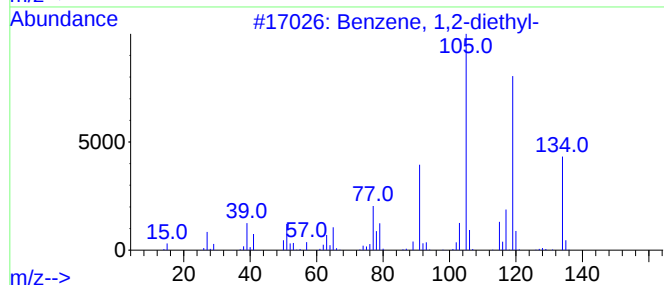
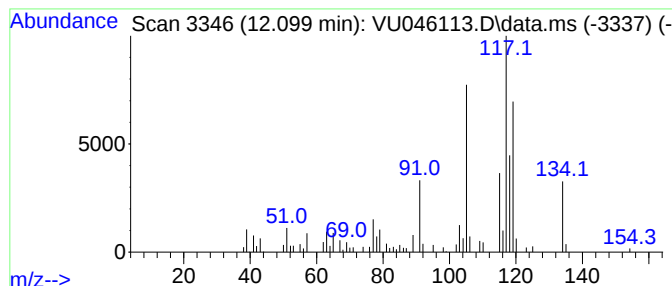
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Quant Title : VOC Analysis

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

Peak Number 14 Benzene, 1,2-diethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.099	4.87 ug/L	63469	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	83
2			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	55
3			Deltacyclene	118	C9H10	007785-10-6	55
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	55
5			Deltacyclene	118	C9H10	007785-10-6	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKQ1DL

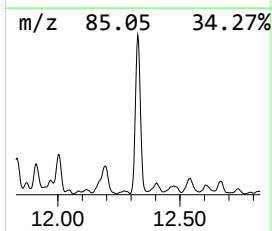
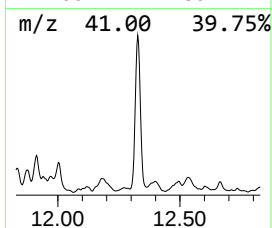
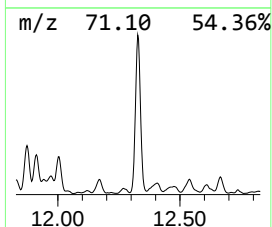
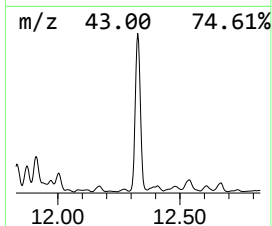
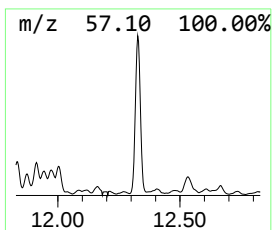
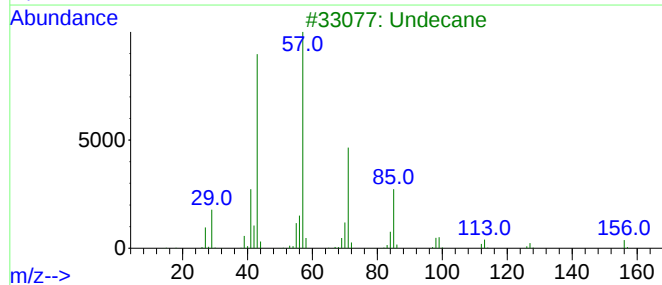
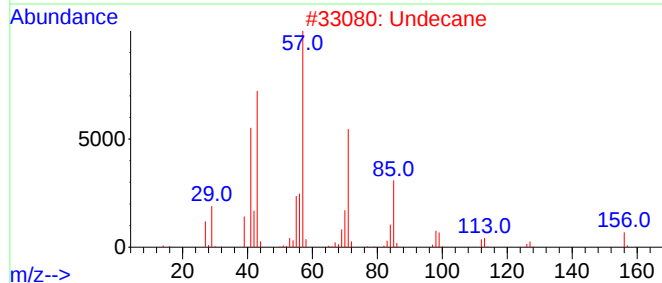
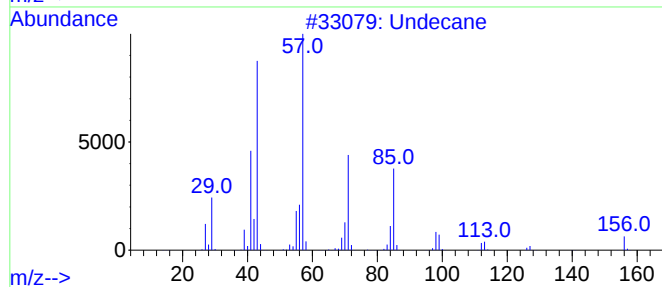
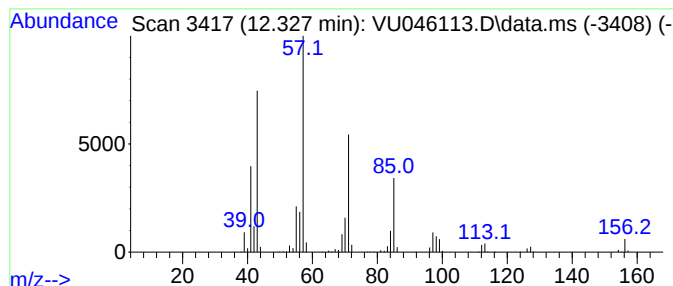
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Quant Title : VOC Analysis

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.327	33.18 ug/L	432774	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	96
2	Undecane			156	C11H24	001120-21-4	94
3	Undecane			156	C11H24	001120-21-4	91
4	Carbonic acid, prop-1-en-2-yl tr...			284	C17H32O3	1000382-90-8	86
5	Hexadecane			226	C16H34	000544-76-3	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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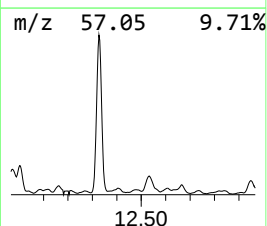
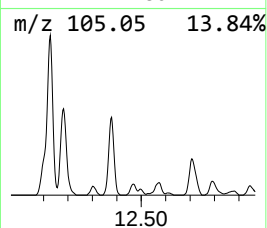
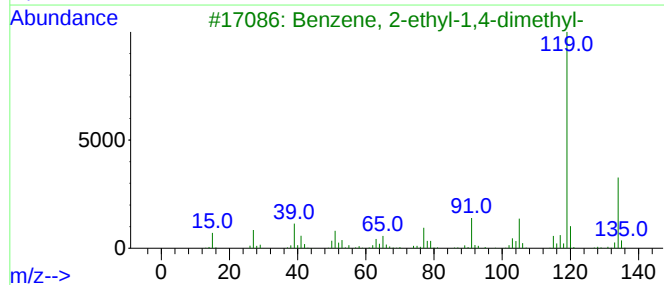
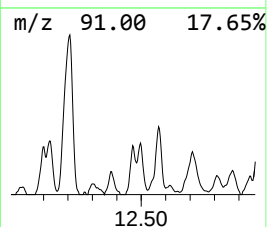
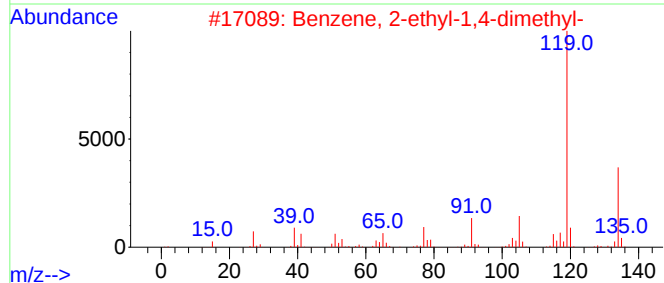
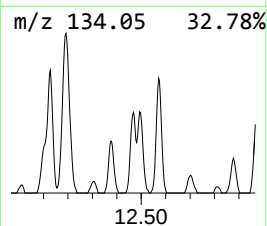
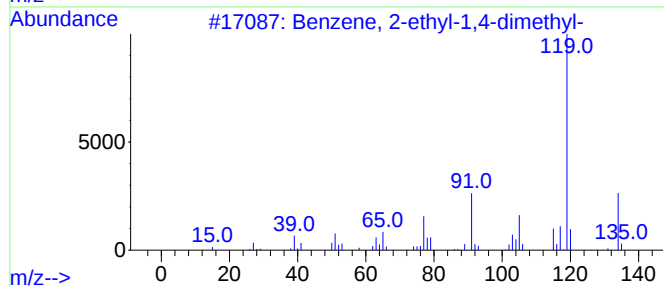
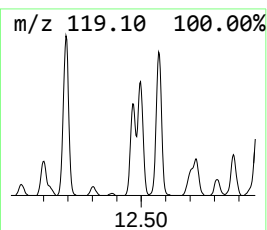
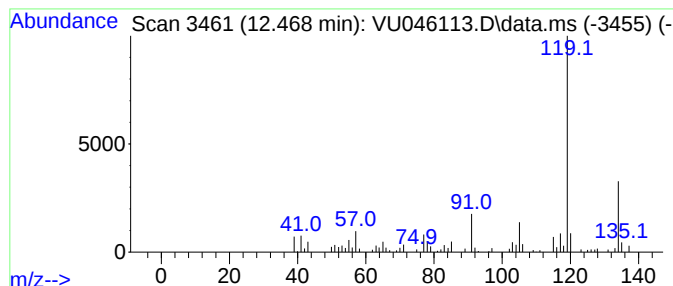
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Quant Title : VOC Analysis

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

Peak Number 16 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.468	3.16 ug/L	41252	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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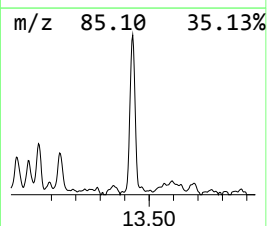
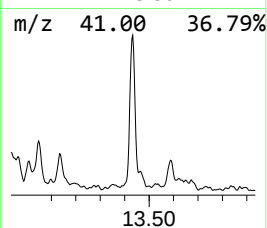
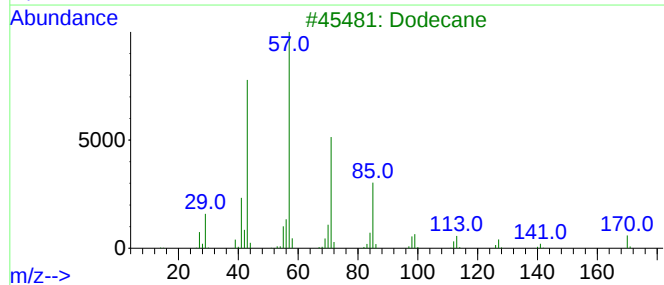
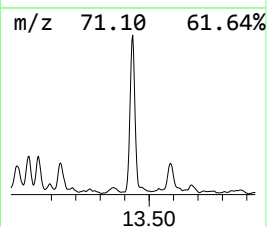
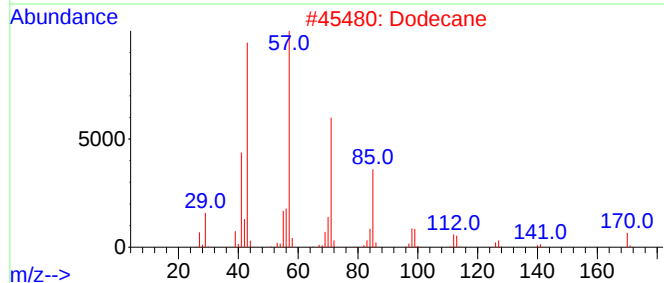
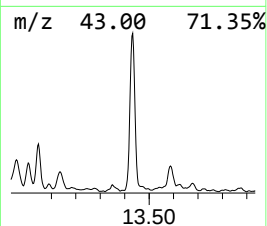
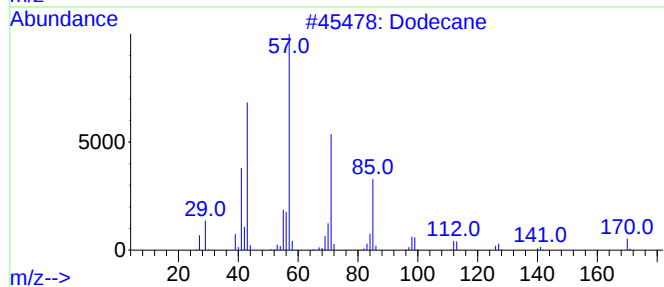
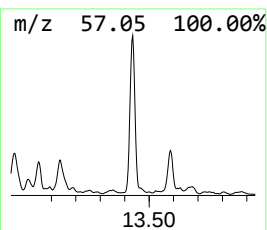
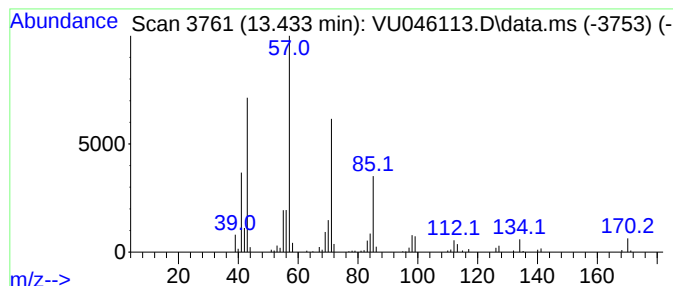
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Quant Title : VOC Analysis

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

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.433	16.76 ug/L	218592	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	96
2			Dodecane	170	C12H26	000112-40-3	94
3			Dodecane	170	C12H26	000112-40-3	87
4			Tetradecane	198	C14H30	000629-59-4	87
5			Dodecane	170	C12H26	000112-40-3	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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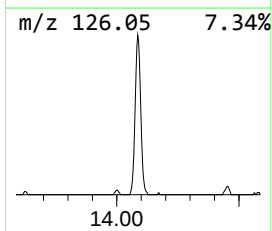
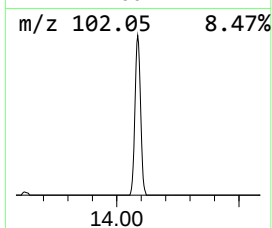
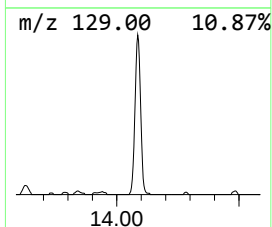
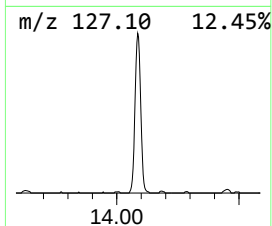
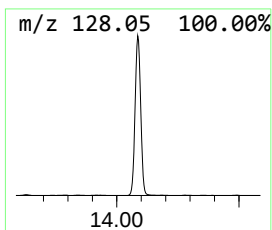
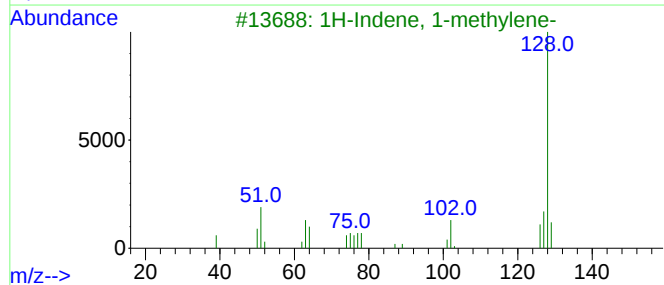
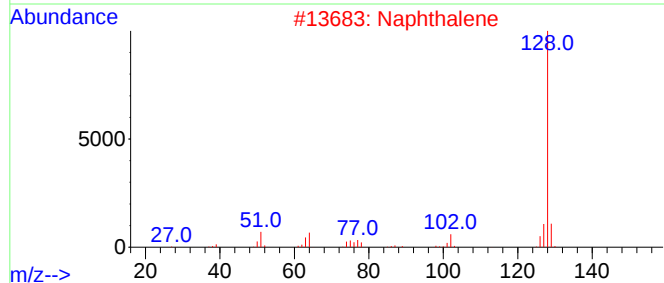
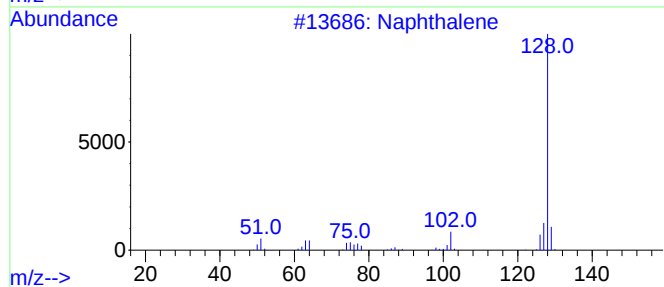
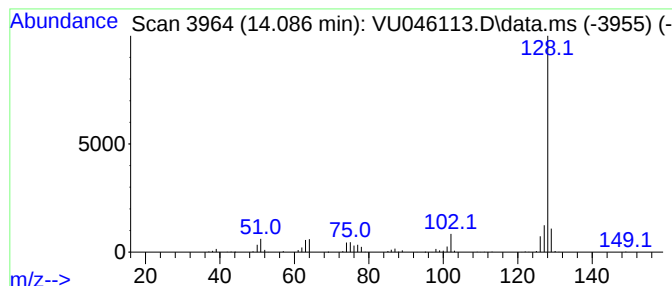
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Quant Title : VOC Analysis

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

Peak Number 18 Azulene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.086	32.39 ug/L	422574	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Naphthalene	128	C10H8	000091-20-3	94
3			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	91
4			Azulene	128	C10H8	000275-51-4	91
5			Azulene	128	C10H8	000275-51-4	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKQ1DL

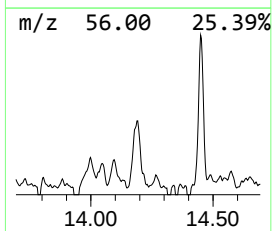
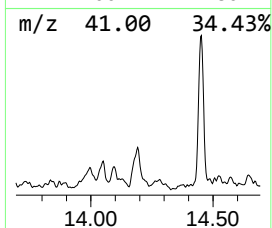
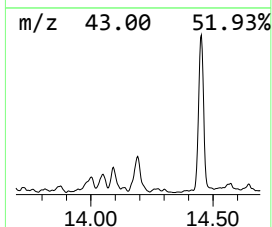
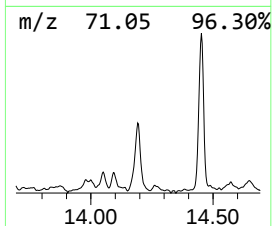
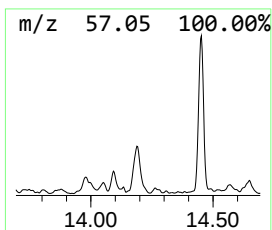
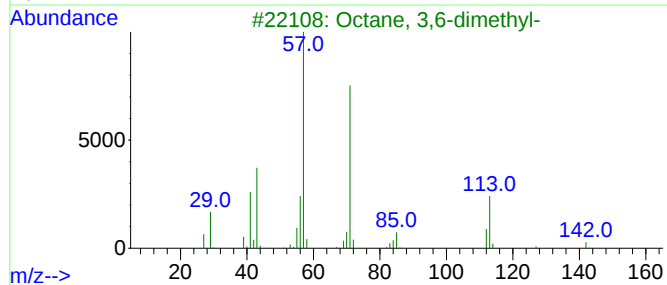
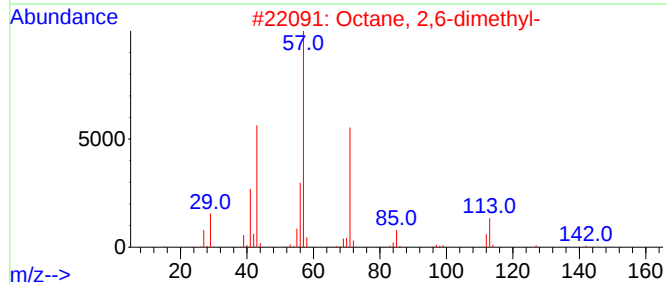
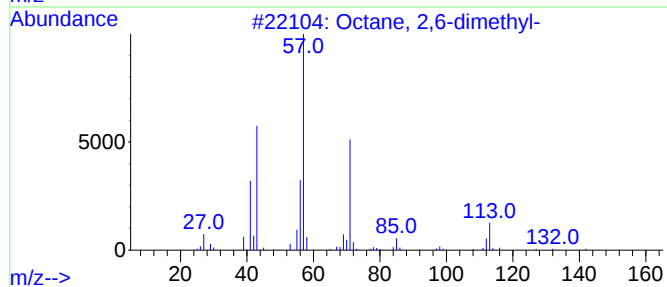
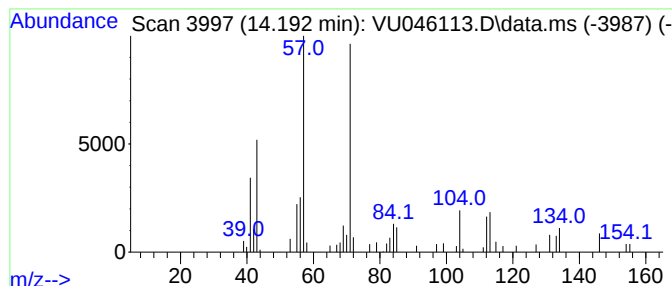
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Quant Title : VOC Analysis

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.192	3.67 ug/L	47926	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
2			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	59
3			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	59
4			Carbonic acid, neopentyl 2-ethyl...	244	C14H28O3	1000357-85-7	50
5			Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKQ1DL

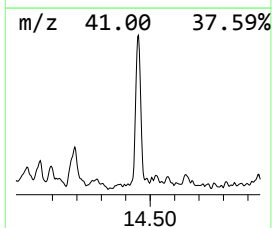
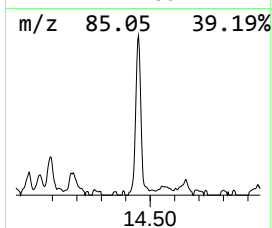
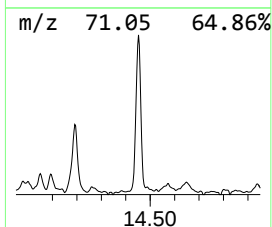
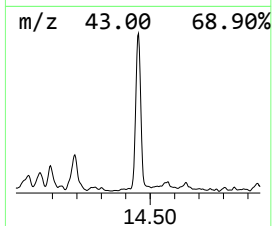
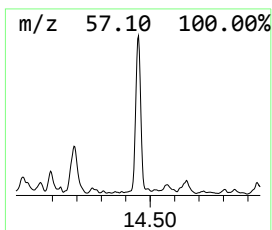
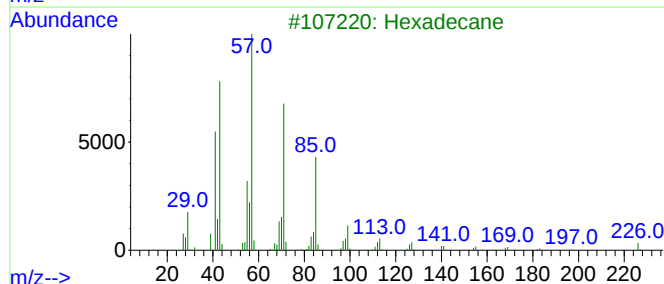
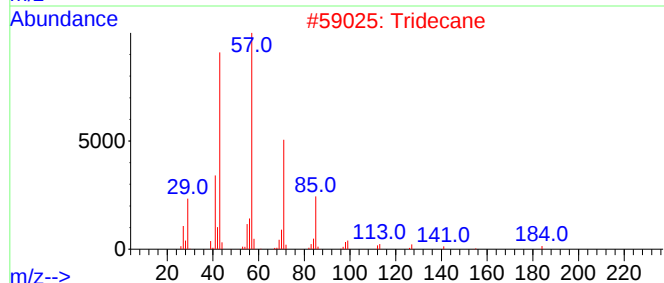
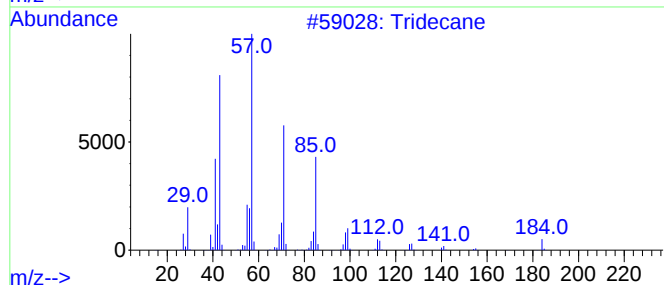
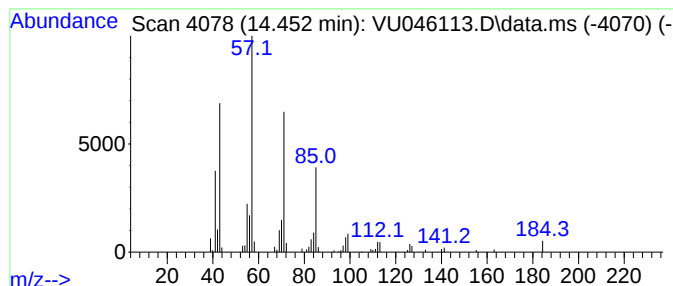
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Quant Title : VOC Analysis

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.452	6.57 ug/L	85681	1,4-Dichlorobenzene-d4	11.812

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane	184	C13H28	000629-50-5	98
2		Tridecane	184	C13H28	000629-50-5	91
3		Hexadecane	226	C16H34	000544-76-3	90
4		Dodecane	170	C12H26	000112-40-3	90
5		Dodecane	170	C12H26	000112-40-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 3.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKQ1DL

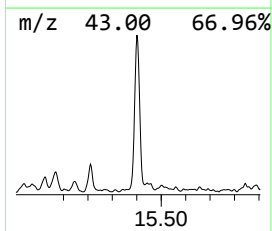
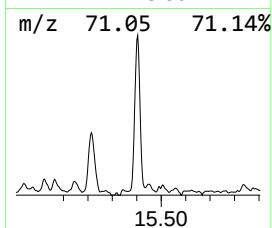
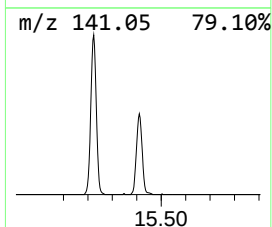
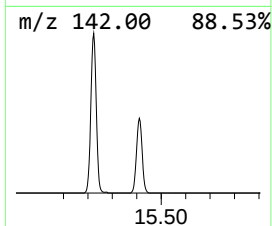
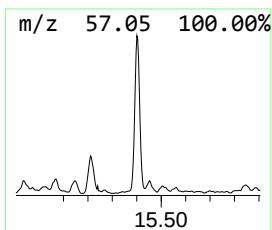
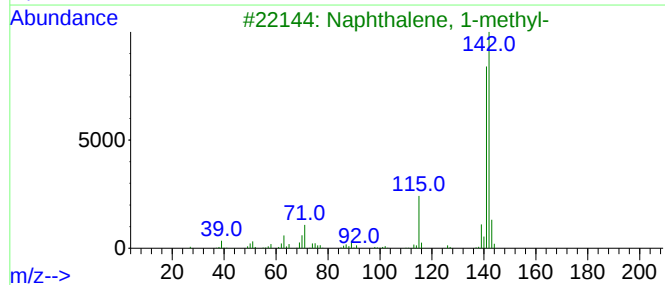
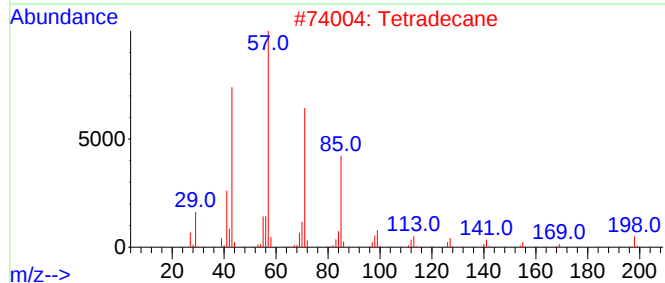
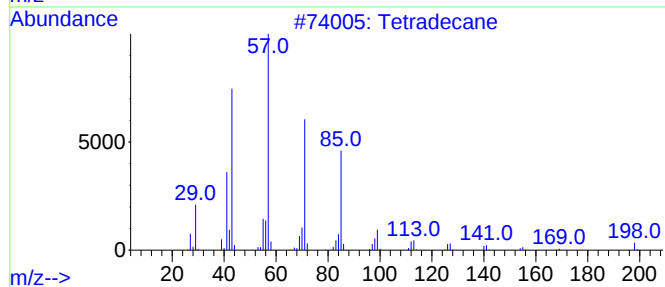
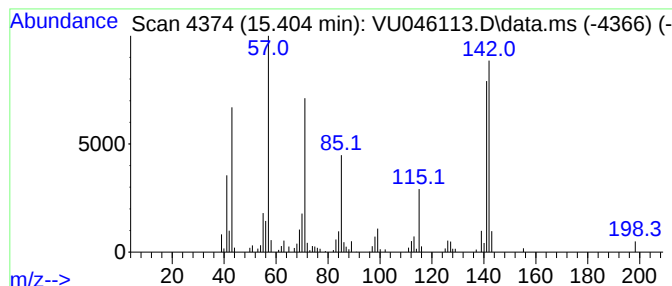
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Quant Title : VOC Analysis

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

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.404	9.82 ug/L	128034	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	78
2			Tetradecane	198	C14H30	000629-59-4	78
3			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	76
4			Tetradecane	198	C14H30	000629-59-4	70
5			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046113.D
Acq On : 06 Dec 2021 15:42
Operator : SY/MD
Sample : M4942-03DL 100X
Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKQ1DL

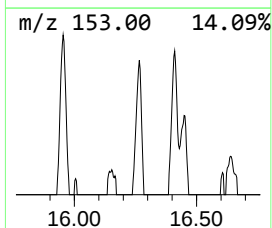
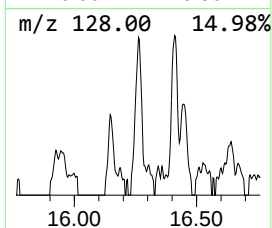
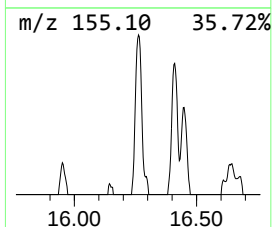
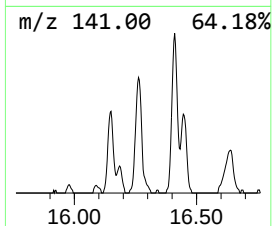
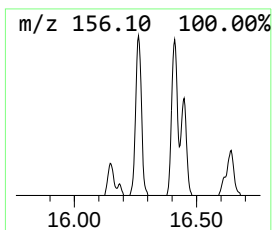
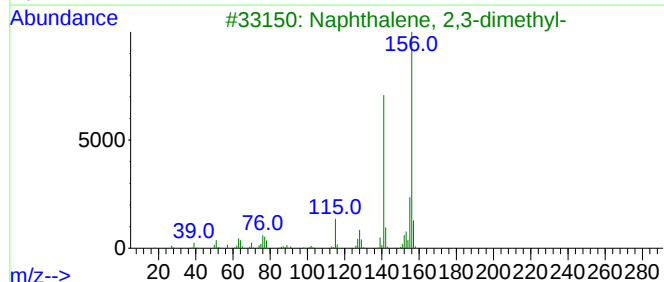
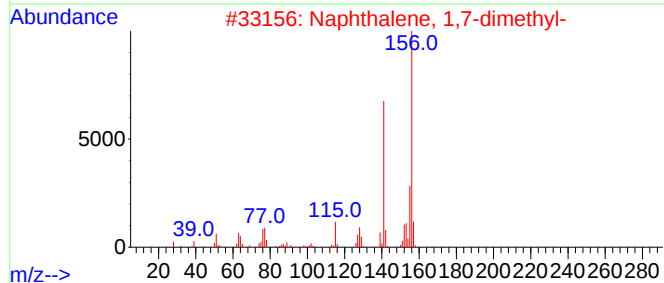
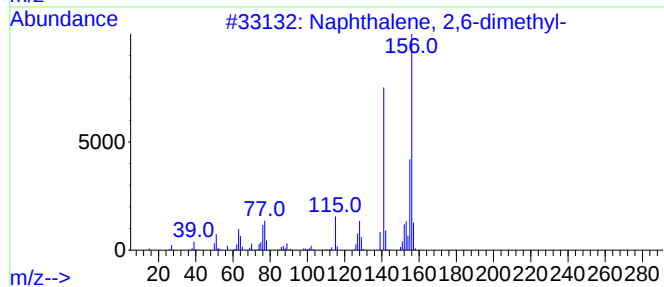
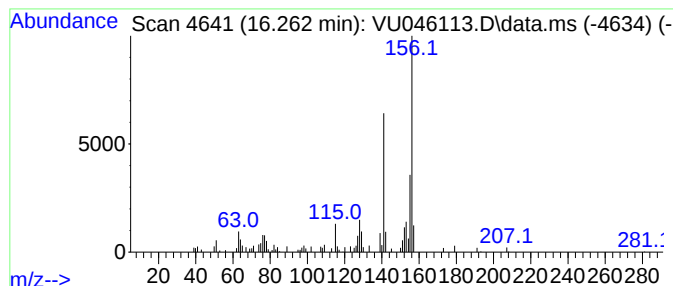
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Quant Title : VOC Analysis

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

Peak Number 23 Naphthalene, 2,6-dimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.262	3.07 ug/L	40053	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
5			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
 Data File : VU046113.D
 Acq On : 06 Dec 2021 15:42
 Operator : SY/MD
 Sample : M4942-03DL 100X
 Misc : 5.89g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BGKQ1DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.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
(DEL) Alkane: S...	10.022	4.1	ug/L	44402	2	9.420	546305	50.0
(DEL) Alkane: S...	10.253	7.4	ug/L	80388	2	9.420	546305	50.0
(DEL) Alkane: C...	10.356	4.0	ug/L	43646	2	9.420	546305	50.0
(DEL) Alkane: S...	10.558	5.6	ug/L	60776	2	9.420	546305	50.0
Benzene, propyl-	10.906	3.5	ug/L	46219	3	11.812	652223	50.0
Benzene, 1-ethy...	10.999	14.8	ug/L	193306	3	11.812	652223	50.0
Benzene, 1-ethy...	11.298	10.1	ug/L	132050	3	11.812	652223	50.0
(DEL) Alkane: S...	11.375	3.1	ug/L	40740	3	11.812	652223	50.0
(DEL) Alkane: S...	11.417	9.9	ug/L	129340	3	11.812	652223	50.0
(DEL) Alkane: S...	11.575	4.0	ug/L	52014	3	11.812	652223	50.0
Ethanone, 1-cyc...	11.706	7.8	ug/L	101960	3	11.812	652223	50.0
Benzene, 1,2-di...	12.099	4.9	ug/L	63469	3	11.812	652223	50.0
(DEL) Alkane: S...	12.327	33.2	ug/L	432774	3	11.812	652223	50.0
Benzene, 2-ethy...	12.468	3.2	ug/L	41252	3	11.812	652223	50.0
(DEL) Alkane: S...	13.433	16.8	ug/L	218592	3	11.812	652223	50.0
Azulene	14.086	32.4	ug/L	422574	3	11.812	652223	50.0
(DEL) Alkane: S...	14.192	3.7	ug/L	47926	3	11.812	652223	50.0
(DEL) Alkane: S...	14.452	6.6	ug/L	85681	3	11.812	652223	50.0
(DEL) Alkane: S...	15.404	9.8	ug/L	128034	3	11.812	652223	50.0
Naphthalene, 2,...	16.262	3.1	ug/L	40053	3	11.812	652223	50.0