

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU013020\
 Data File : VU036639.D
 Acq On : 31 Jan 2020 00:11
 Operator : JC/MD
 Sample : L1324-08
 Misc : 5.00g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 GAT65

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM012920WMA.M
 Title : VOC Analysis

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.488	34	46	54	rBV2	16474	29385	0.11%	0.027%
2	1.607	76	83	99	rVB	87544	112278	0.41%	0.103%
3	1.887	165	170	180	rVB	26133	32057	0.12%	0.029%
4	2.539	361	373	394	rVB	161883	270675	0.99%	0.247%
5	2.639	394	404	411	rBV5	2249	4447	0.02%	0.004%
6	2.993	505	514	517	rBV	958	1376	0.01%	0.001%
7	3.060	527	535	546	rVB3	1431	2648	0.01%	0.002%
8	3.218	566	584	603	rBV2	14755	39015	0.14%	0.036%
9	3.366	626	630	637	rVB2	484	559	0.00%	0.001%
10	3.436	650	652	656	rBV	287	226	0.00%	0.000%
11	3.610	700	706	710	rBV	180	288	0.00%	0.000%
12	3.700	729	734	735	rBV	862	712	0.00%	0.001%
13	4.009	827	830	836	rVB	172	211	0.00%	0.000%
14	4.047	837	842	846	rBV	191	273	0.00%	0.000%
15	4.147	871	873	877	rVB	249	154	0.00%	0.000%
16	4.597	1009	1013	1017	rVB	287	256	0.00%	0.000%
17	4.642	1017	1027	1029	rBV2	455	690	0.00%	0.001%
18	4.706	1029	1047	1090	rVV4	47286	203154	0.75%	0.186%
19	5.105	1154	1171	1195	rVB	138429	334046	1.23%	0.305%
20	5.276	1222	1224	1227	rVB2	397	221	0.00%	0.000%
21	5.301	1227	1232	1235	rBV	293	236	0.00%	0.000%
22	5.324	1235	1239	1248	rBV3	396	692	0.00%	0.001%
23	5.404	1258	1264	1268	rBV2	528	737	0.00%	0.001%
24	5.755	1353	1373	1401	rVV2	371713	919652	3.38%	0.840%
25	6.028	1450	1458	1465	rVB3	584	601	0.00%	0.001%
26	6.282	1520	1537	1566	rVV	313758	630902	2.32%	0.576%
27	6.517	1607	1610	1613	rBV	142	151	0.00%	0.000%
28	6.559	1613	1623	1634	rVV2	9418	17170	0.06%	0.016%
29	6.629	1639	1645	1646	rBV2	591	400	0.00%	0.000%
30	6.645	1646	1650	1654	rVV3	494	423	0.00%	0.000%
31	6.726	1660	1675	1697	rVB	198410	404669	1.49%	0.370%
32	6.838	1708	1710	1719	rVB2	461	460	0.00%	0.000%
33	7.047	1771	1775	1776	rBV	255	159	0.00%	0.000%
34	7.173	1811	1814	1816	rBV2	237	141	0.00%	0.000%

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Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM012920WMA.M
 Title : VOC Analysis

35	7.221	1820	1829	1840	rVV5	3080	5926	0.02%	0.005%
36	7.343	1862	1867	1870	rBV	198	242	0.00%	0.000%
37	7.520	1918	1922	1926	rBV2	186	217	0.00%	0.000%
38	7.597	1933	1946	1969	rBV2	110978	209172	0.77%	0.191%
39	7.716	1977	1983	1991	rBV4	1068	1721	0.01%	0.002%
40	7.825	2008	2017	2021	rBV4	945	1694	0.01%	0.002%
41	7.928	2035	2049	2063	rBV	445202	793948	2.92%	0.725%
42	7.999	2063	2071	2094	rVB	20187	36310	0.13%	0.033%
43	8.102	2100	2103	2106	rBV2	212	134	0.00%	0.000%
44	8.153	2111	2119	2125	rBV3	2189	2781	0.01%	0.003%
45	8.211	2125	2137	2156	rVB	74375	141532	0.52%	0.129%
46	8.385	2187	2191	2192	rBV	249	164	0.00%	0.000%
47	8.504	2223	2228	2231	rBV	318	340	0.00%	0.000%
48	8.578	2238	2251	2268	rBV	283122	475994	1.75%	0.435%
49	8.690	2268	2286	2304	rBV	240150	534545	1.96%	0.488%
50	8.812	2321	2324	2326	rBV2	708	506	0.00%	0.000%
51	8.890	2339	2348	2356	rVB3	2278	3412	0.01%	0.003%
52	8.947	2356	2366	2377	rBV3	4847	8060	0.03%	0.007%
53	8.999	2379	2382	2388	rVB4	750	662	0.00%	0.001%
54	9.092	2400	2411	2418	rVB5	1532	2850	0.01%	0.003%
55	9.140	2418	2426	2433	rBV3	2596	3557	0.01%	0.003%
56	9.189	2433	2441	2447	rVV5	4372	7187	0.03%	0.007%
57	9.237	2449	2456	2466	rVB4	7265	11828	0.04%	0.011%
58	9.359	2484	2494	2504	rBV	9748	15045	0.06%	0.014%
59	9.449	2506	2522	2544	rBV	419880	791489	2.91%	0.723%
60	9.600	2555	2569	2582	rBV	110040	199897	0.73%	0.183%
61	9.726	2597	2608	2615	rBV	97518	178086	0.65%	0.163%
62	9.896	2652	2661	2675	rVB10	5140	10665	0.04%	0.010%
63	9.976	2678	2686	2695	rBV6	2888	4627	0.02%	0.004%
64	10.057	2695	2711	2721	rBV6	14558	40414	0.15%	0.037%
65	10.131	2723	2734	2748	rVV	170625	313390	1.15%	0.286%
66	10.230	2761	2765	2766	rBV3	1287	900	0.00%	0.001%
67	10.272	2766	2778	2793	rVV2	51288	111949	0.41%	0.102%
68	10.382	2805	2812	2816	rBV3	28536	41701	0.15%	0.038%
69	10.510	2839	2852	2860	rBV3	21294	50082	0.18%	0.046%
70	10.645	2888	2894	2899	rBV	38608	49554	0.18%	0.045%
71	10.719	2911	2917	2926	rBV7	12645	22818	0.08%	0.021%

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72	10.787	2927	2938	2947	rBV	331694	589759	2.17%	0.539%
73	11.050	3005	3020	3029	rBV	777054	1430386	5.26%	1.307%
74	11.118	3033	3041	3061	rVB2	672873	1709856	6.28%	1.562%
75	11.320	3094	3104	3119	rVB2	221246	537506	1.98%	0.491%
76	11.436	3133	3140	3147	rBV	149828	195177	0.72%	0.178%
77	11.494	3147	3158	3169	rVV	2056087	3200523	11.76%	2.924%
78	11.562	3169	3179	3193	rVB	1483745	2254576	8.29%	2.060%
79	11.732	3226	3232	3238	rBV2	145800	215501	0.79%	0.197%
80	11.838	3255	3265	3276	rVB2	559131	972951	3.58%	0.889%
81	11.918	3283	3290	3300	rVB	950141	1492939	5.49%	1.364%
82	11.970	3301	3306	3314	rVB	168199	220148	0.81%	0.201%
83	12.214	3372	3382	3395	rVB3	1188005	2188836	8.05%	2.000%
84	12.340	3410	3421	3433	rVB2	3556436	5943979	21.85%	5.431%
85	12.767	3549	3554	3564	rVB	840553	1206993	4.44%	1.103%
86	12.947	3602	3610	3618	rBV3	776570	1393991	5.12%	1.274%
87	13.050	3636	3642	3657	rVB3	936237	2291110	8.42%	2.093%
88	13.539	3788	3794	3803	rVB2	1471058	2098483	7.71%	1.917%
89	13.648	3820	3828	3851	rVB	3221492	6293992	23.13%	5.751%
90	14.108	3962	3971	3990	rVB	10251120	17242477	63.38%	15.755%
91	15.246	4313	4325	4335	rVB	16945359	27206370	100.00%	24.860%
92	15.430	4372	4382	4393	rBV	9180686	14584177	53.61%	13.326%
93	16.159	4601	4609	4618	rBV	1521249	2259907	8.31%	2.065%
94	16.275	4636	4645	4667	rVB	1632598	2783641	10.23%	2.544%
95	16.423	4681	4691	4698	rBV	1281285	2042235	7.51%	1.866%
96	16.892	4830	4837	4848	rVB2	198322	305047	1.12%	0.279%
97	17.111	4897	4905	4922	rVB2	346630	825023	3.03%	0.754%
98	17.388	4985	4991	5010	rVB2	192427	392102	1.44%	0.358%
99	17.548	5035	5041	5053	rVB2	153571	252944	0.93%	0.231%
100	17.616	5055	5062	5071	rBV3	141079	227081	0.83%	0.207%

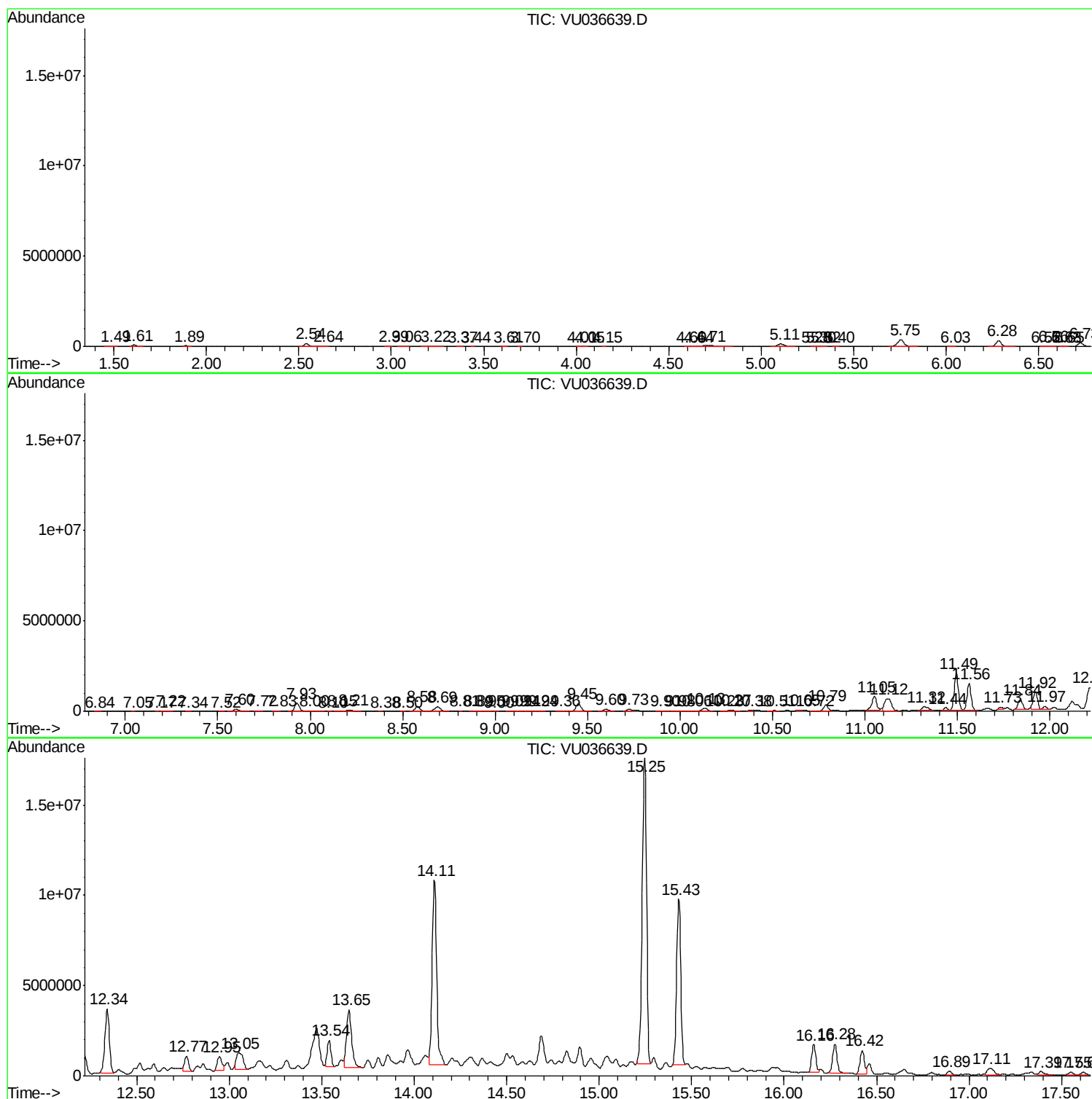
Sum of corrected areas: 109440173

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ALS Vial : 35 Sample Multiplier: 1

Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM012920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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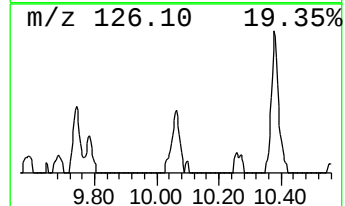
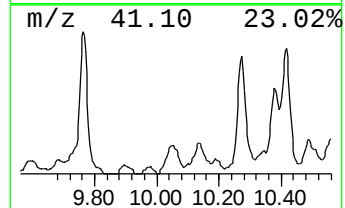
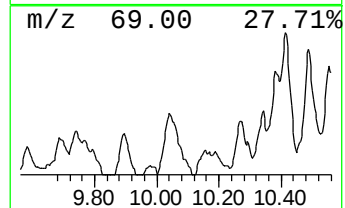
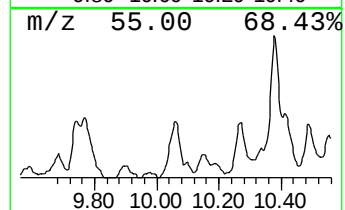
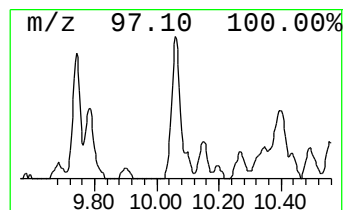
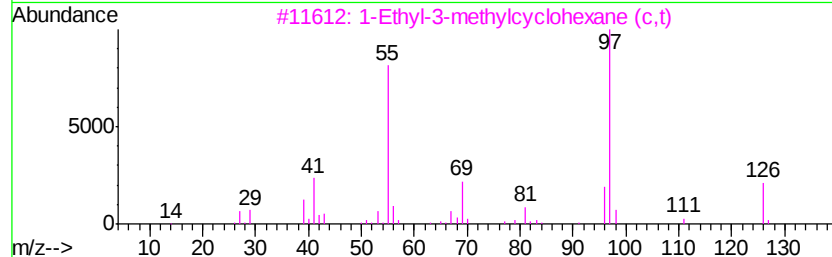
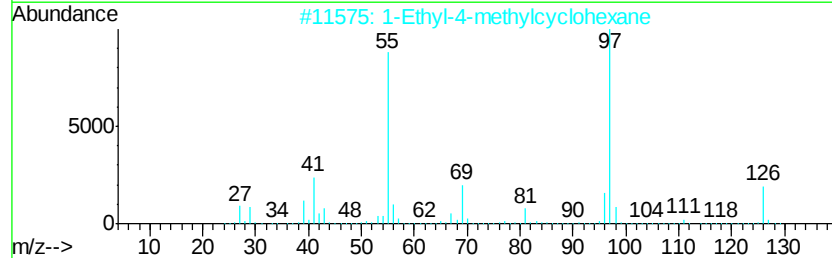
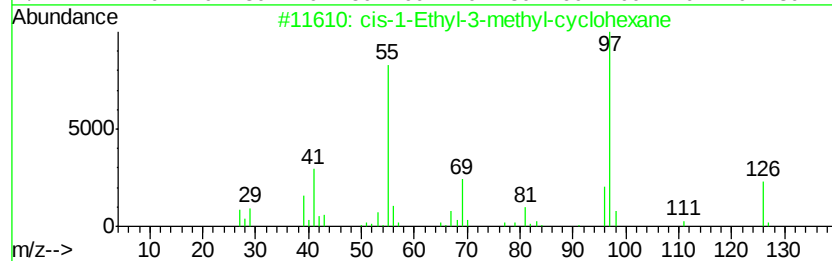
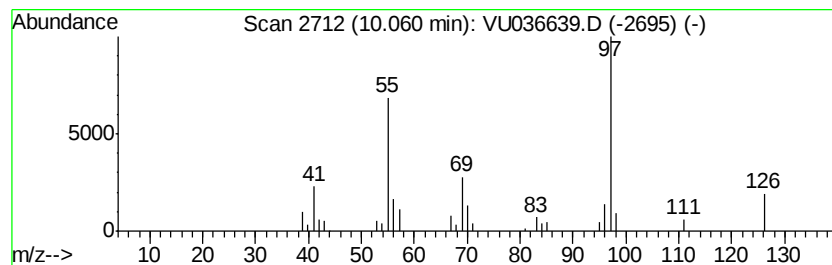
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM012920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Cyclic10.06 Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.06	2.55 ug/L	40414	Chlorobenzene-d5	9.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	64
2		1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	64
3		1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	64
4		Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	59
5		Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	59



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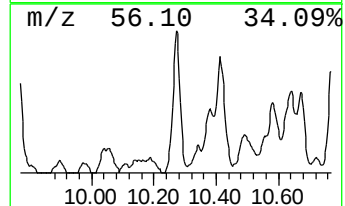
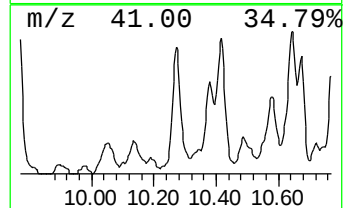
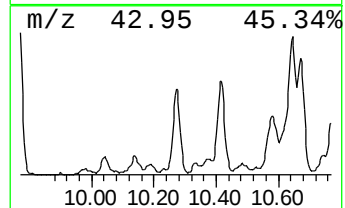
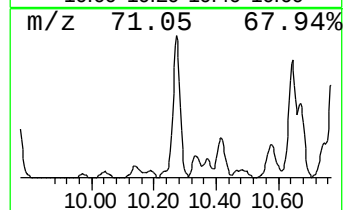
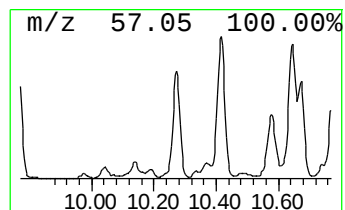
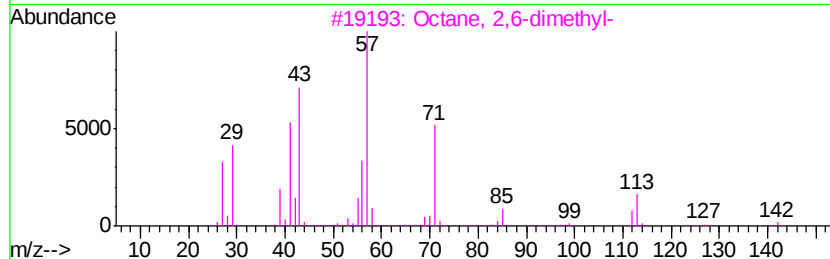
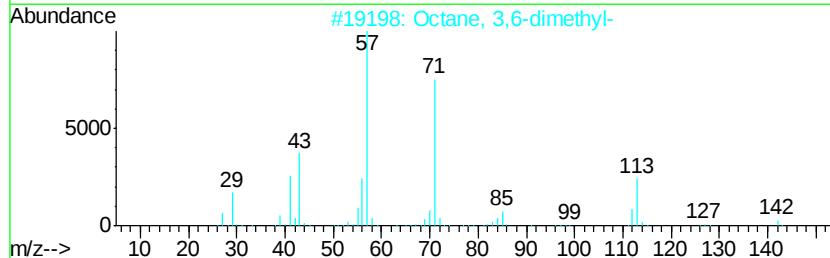
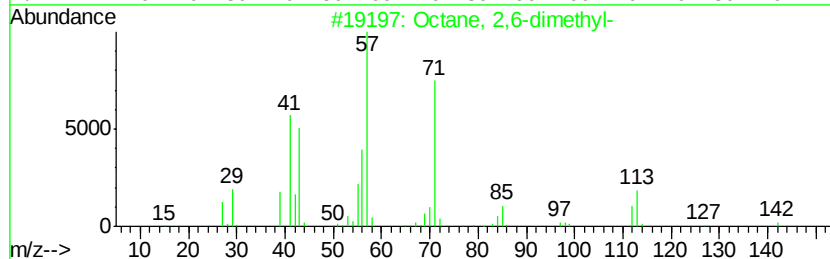
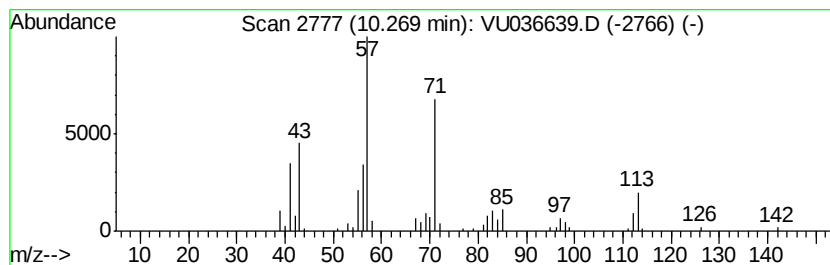
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.27	7.07 ug/L	111949	Chlorobenzene-d5	9.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	91
2		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	90
3		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	90
4		Nonane, 3-methyl-	142	C10H22	005911-04-6	87
5		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	81



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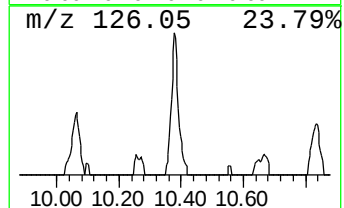
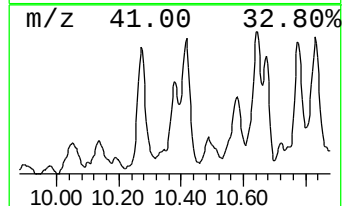
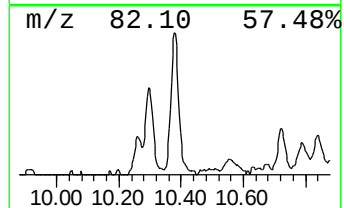
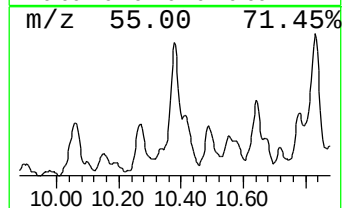
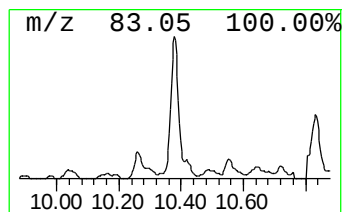
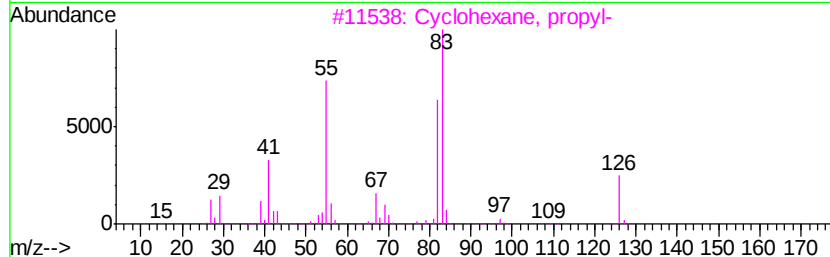
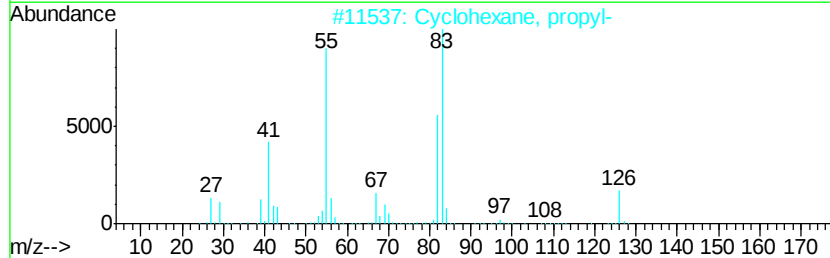
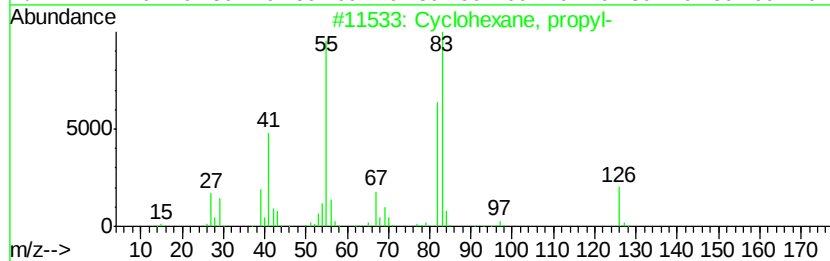
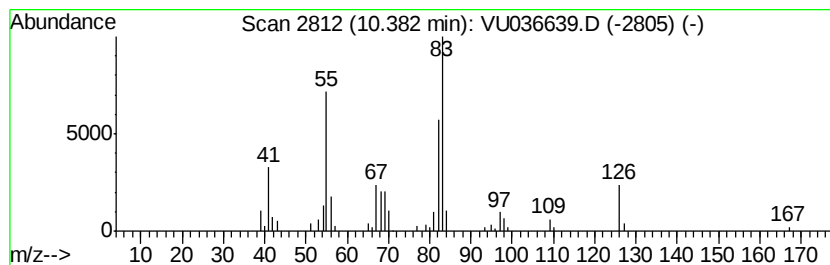
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM012920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Cyclic10.38 Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.38	2.63 ug/L	41701	Chlorobenzene-d5	9.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, propyl-	126	C9H18	001678-92-8	80
2		Cyclohexane, propyl-	126	C9H18	001678-92-8	72
3		Cyclohexane, propyl-	126	C9H18	001678-92-8	64
4		Cyclohexane, 1,1'-(1,4-butanedi...	222	C16H30	006165-44-2	59
5		Cyclohexane, propyl-	126	C9H18	001678-92-8	53



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00uL/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleID :
GAT65

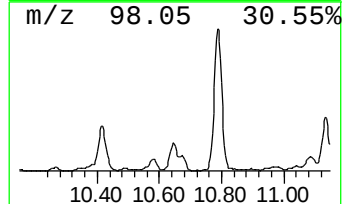
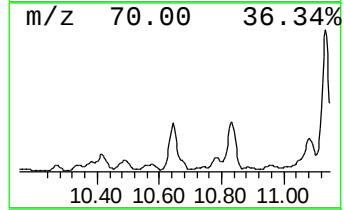
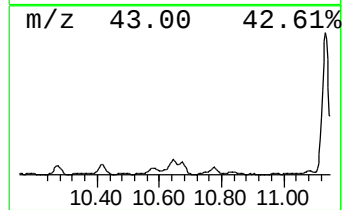
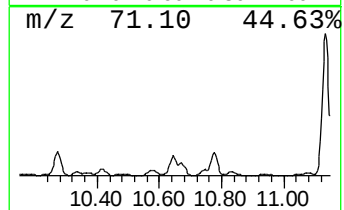
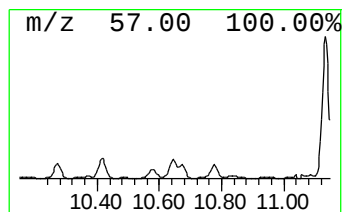
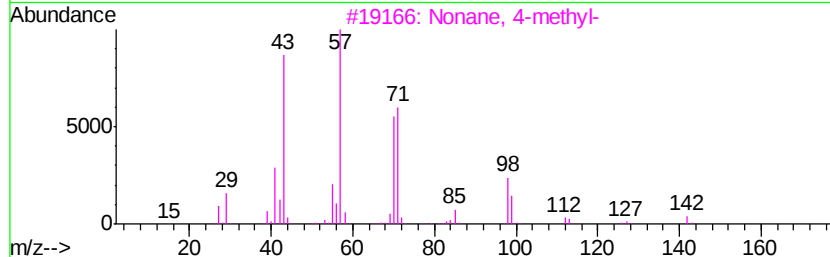
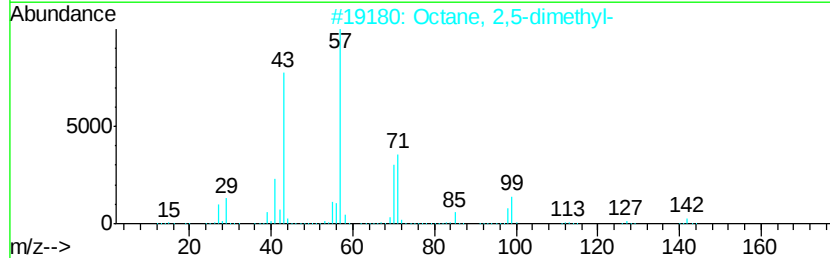
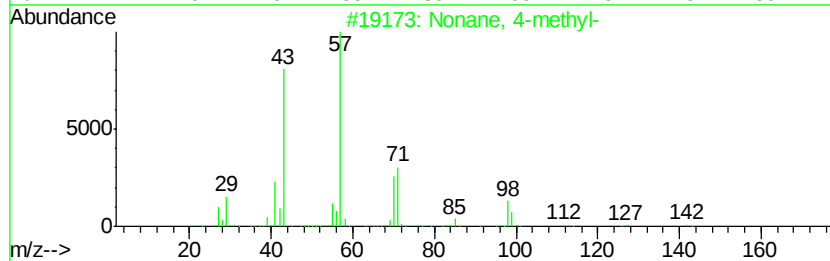
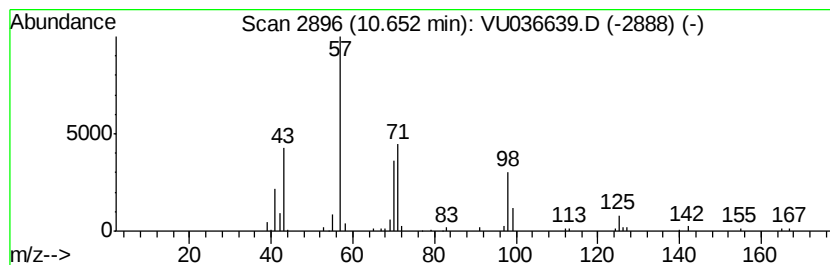
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	2.55 ug/L	49554	1,4-Dichlorobenzene-d4	11.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonane, 4-methyl-	142	C10H22	017301-94-9	87
2		Octane, 2,5-dimethyl-	142	C10H22	015869-89-3	72
3		Nonane, 4-methyl-	142	C10H22	017301-94-9	72
4		Nonane, 4-methyl-	142	C10H22	017301-94-9	64
5		Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00µL/5.0mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

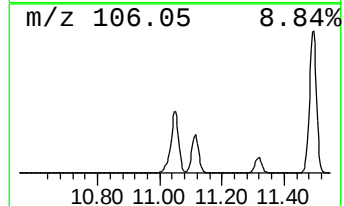
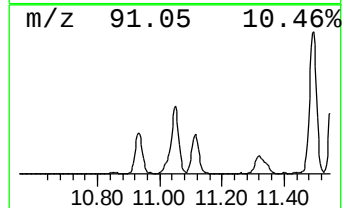
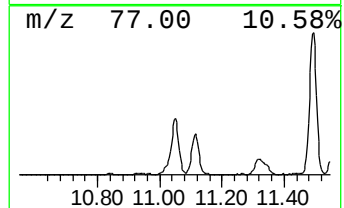
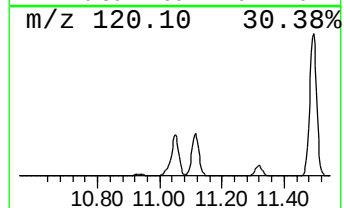
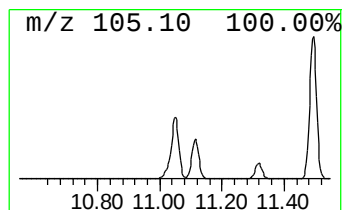
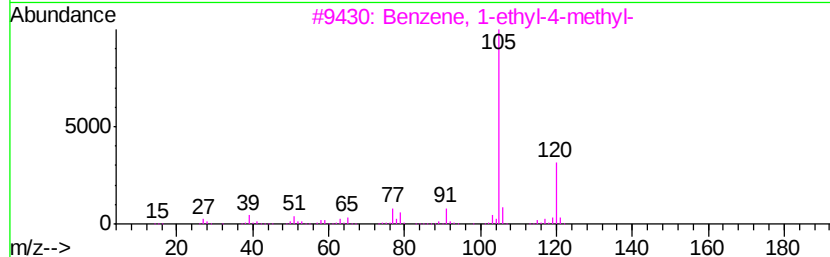
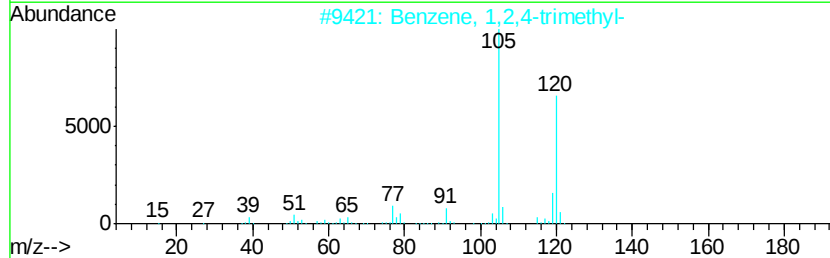
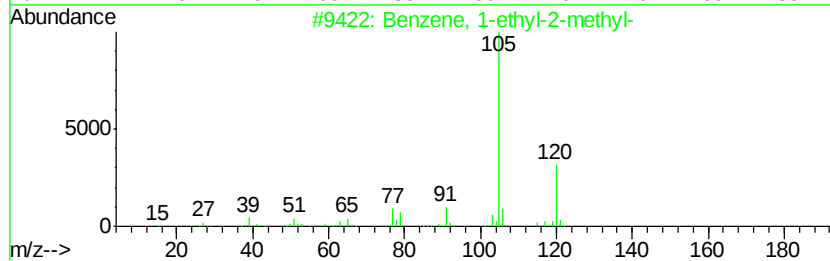
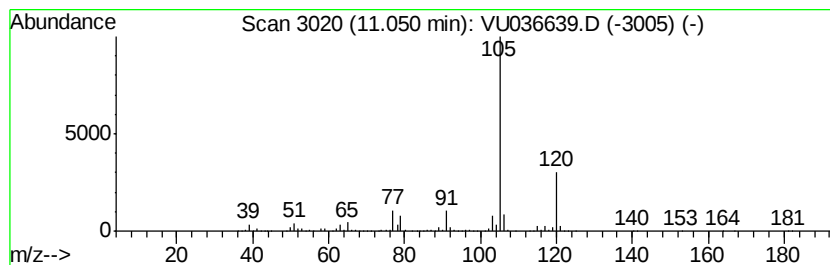
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.05	73.51 ug/L	1430390	1,4-Dichlorobenzene-d4	11.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00µ/5.0mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

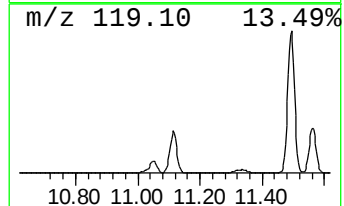
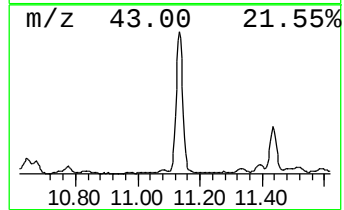
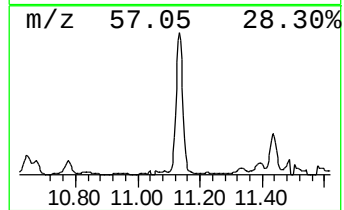
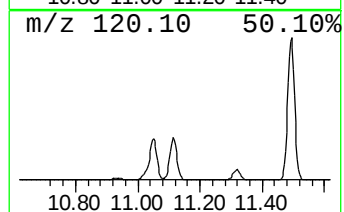
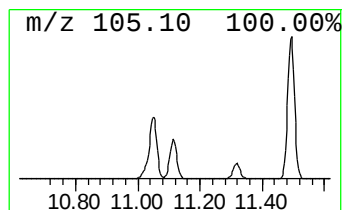
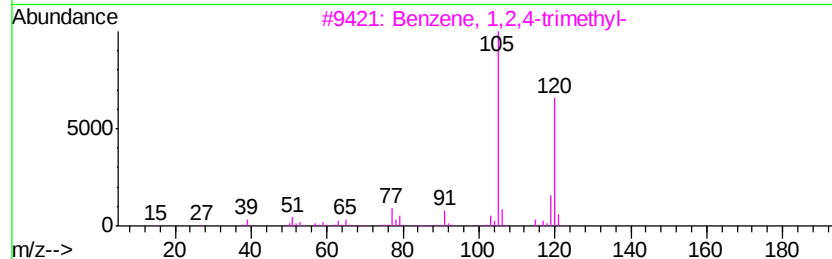
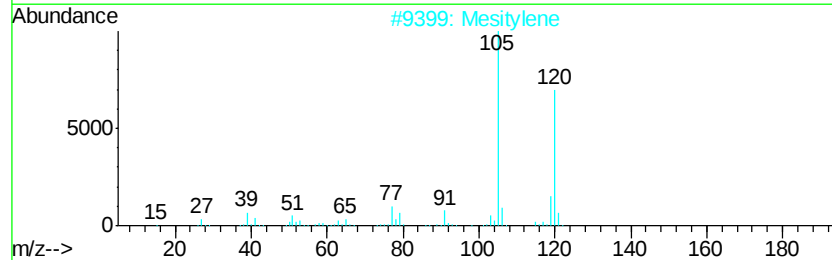
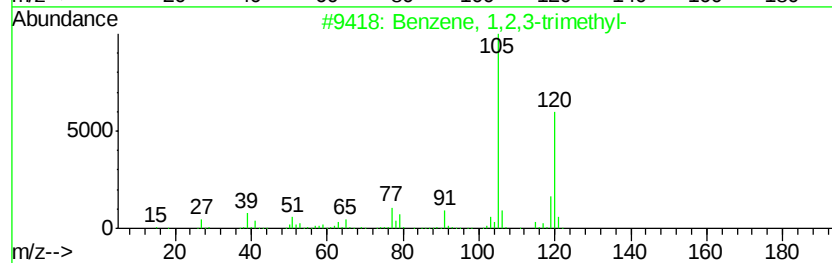
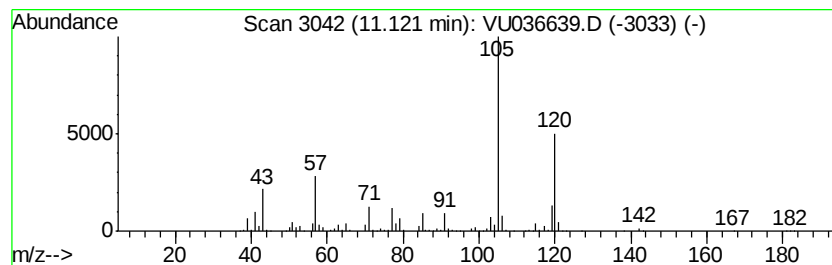
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.12	87.87 ug/L	1709860	1,4-Dichlorobenzene-d4	11.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Mesitylene	120	C9H12	000108-67-8	95
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4		Mesitylene	120	C9H12	000108-67-8	93
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00uL/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampled :
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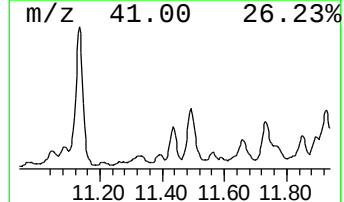
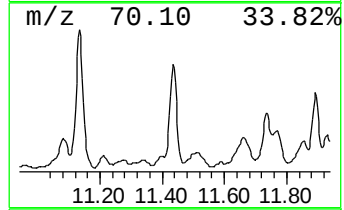
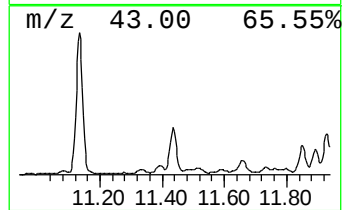
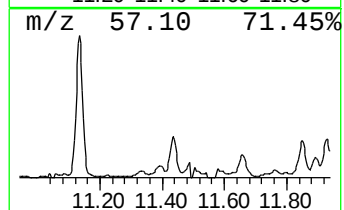
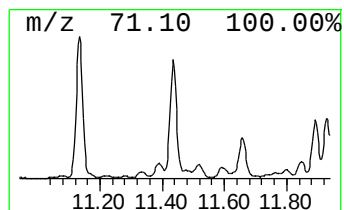
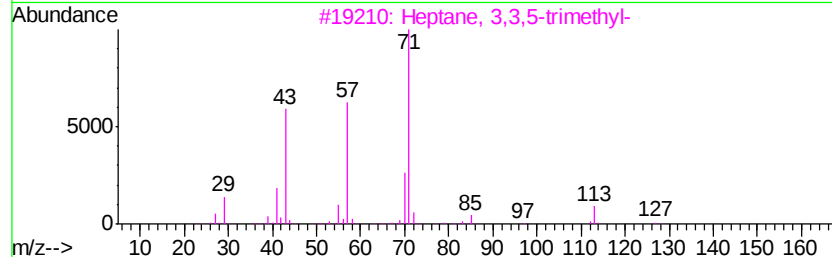
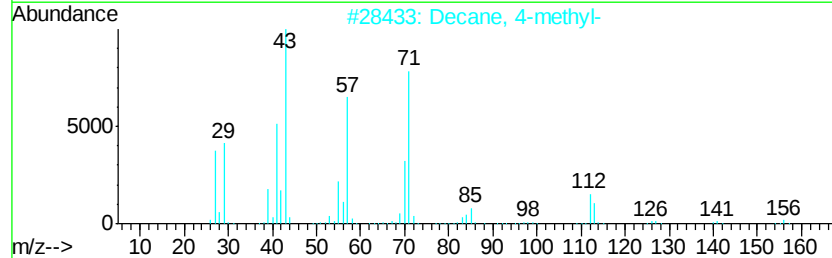
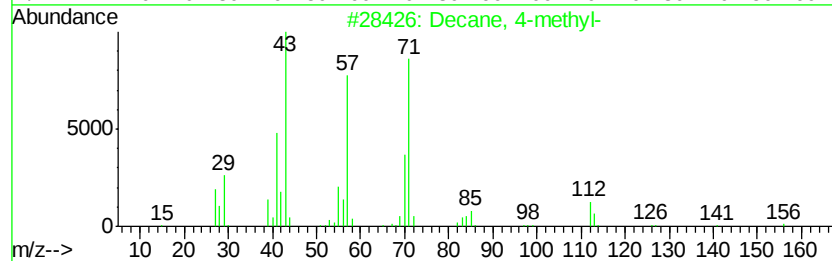
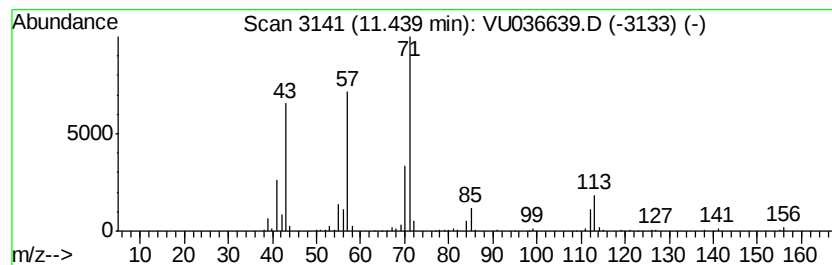
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	10.03 ug/L	195177	1,4-Dichlorobenzene-d4	11.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane, 4-methyl-	156	C11H24	002847-72-5	91
2		Decane, 4-methyl-	156	C11H24	002847-72-5	86
3		Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	72
4		Heptane, 3-ethyl-5-methyl-	142	C10H22	052896-90-9	72
5		4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00uL/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

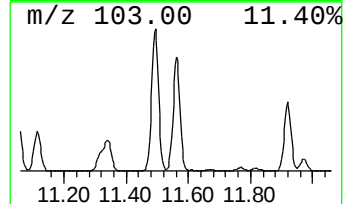
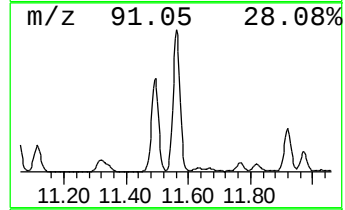
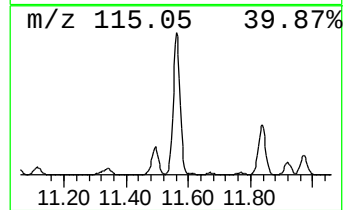
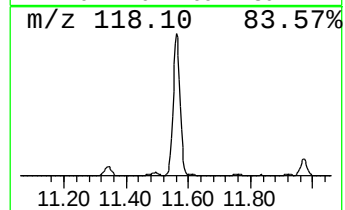
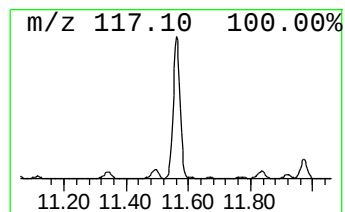
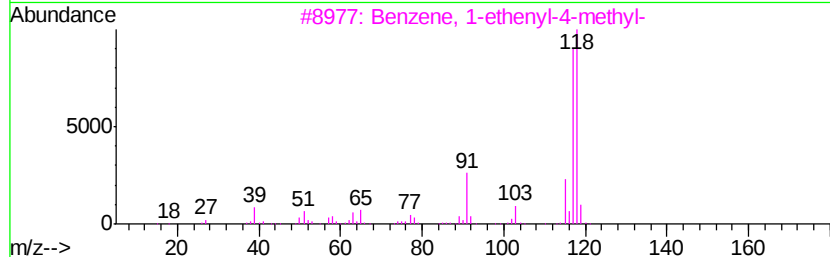
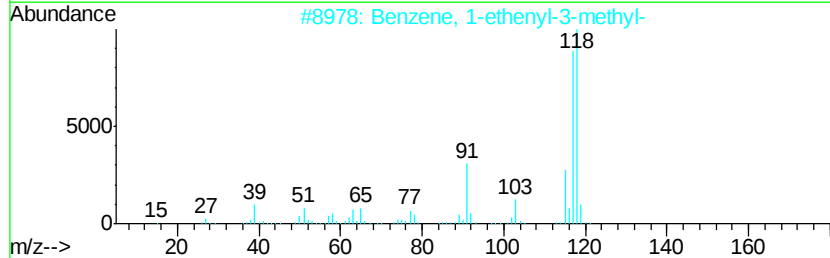
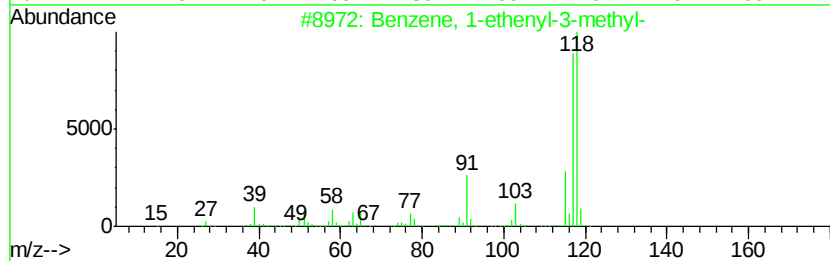
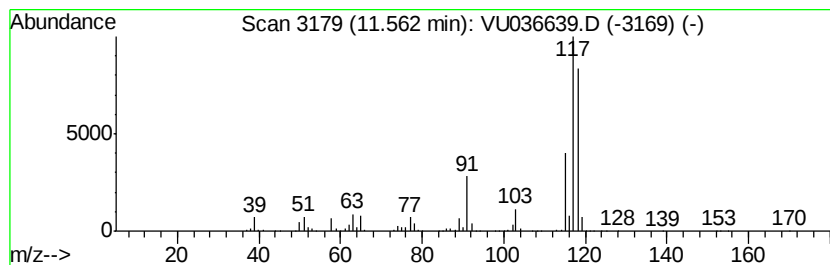
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Benzene, 1-ethenyl-3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.56	115.86 ug/L	2254580	1,4-Dichlorobenzene-d4	11.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	96
2		Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	96
3		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	95
4		Benzene, cyclopropyl-	118	C9H10	000873-49-4	94
5		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00uL/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

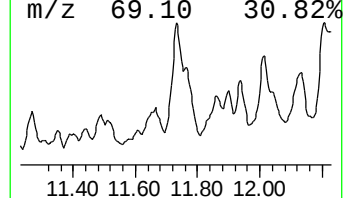
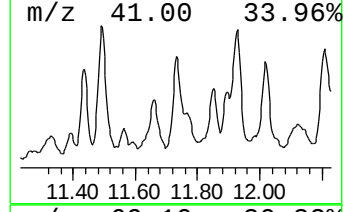
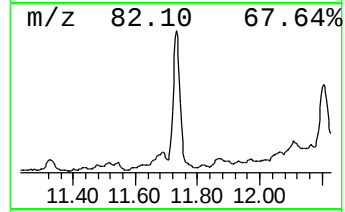
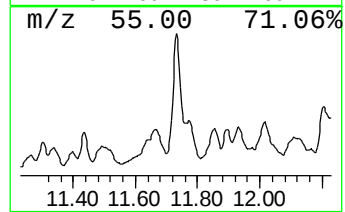
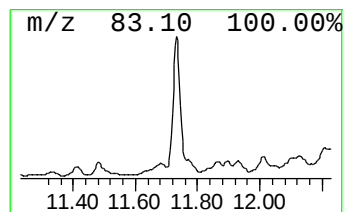
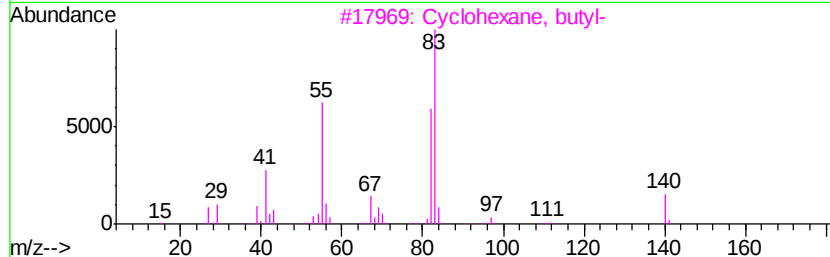
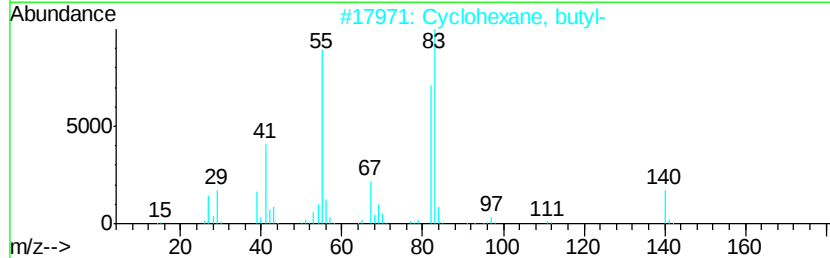
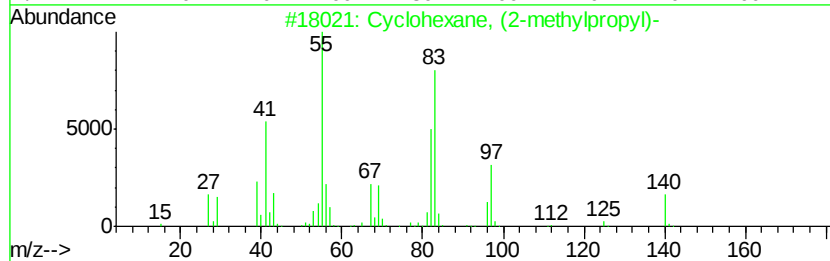
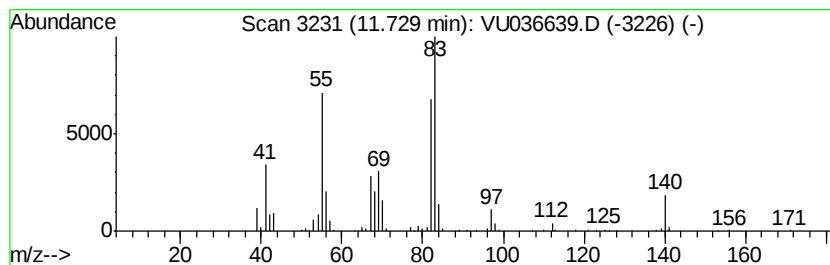
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 (DEL) Alkane: Cyclic11.73 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.73	11.07 ug/L	215501	1,4-Dichlorobenzene-d4	11.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	72
2		Cyclohexane, butyl-	140	C10H20	001678-93-9	64
3		Cyclohexane, butyl-	140	C10H20	001678-93-9	64
4		Cyclohexane, butyl-	140	C10H20	001678-93-9	58
5		Cyclohexane, pentyl-	154	C11H22	004292-92-6	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00µ/5.0mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

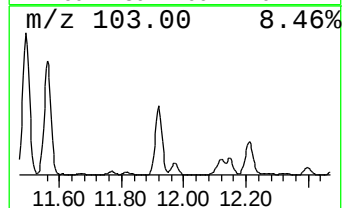
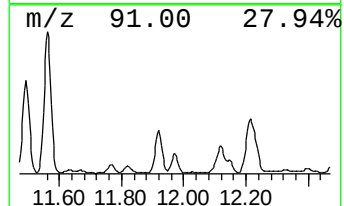
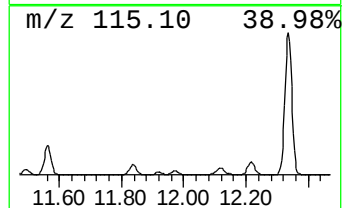
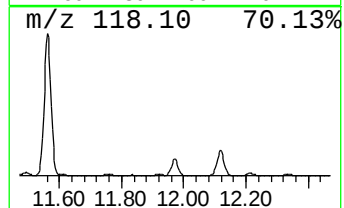
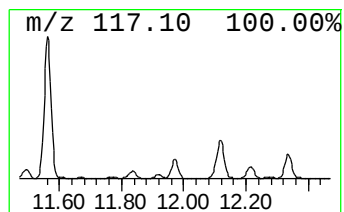
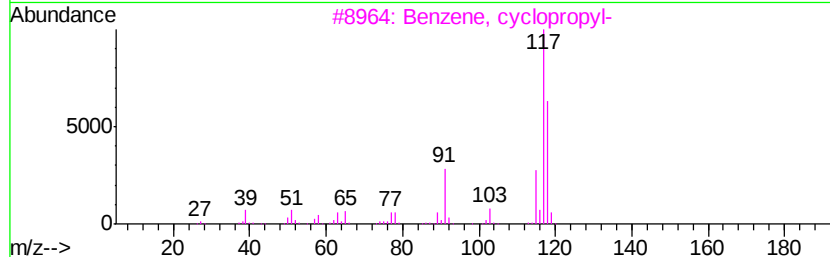
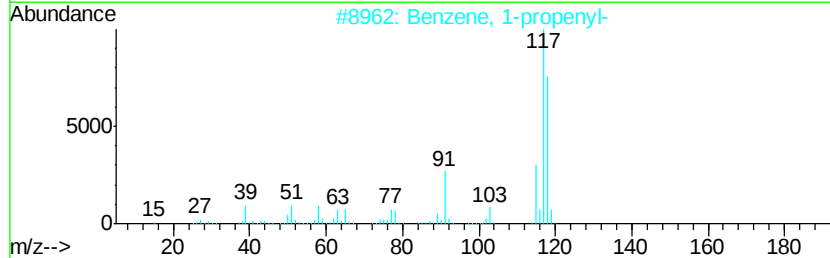
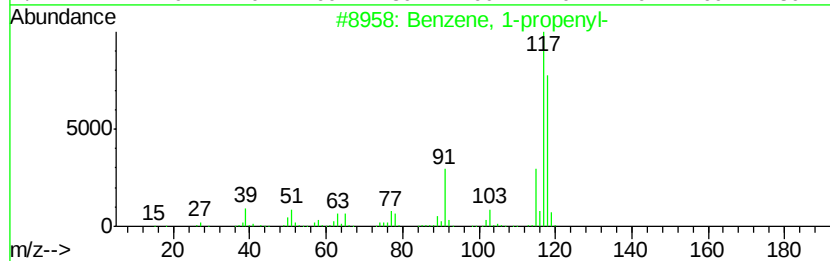
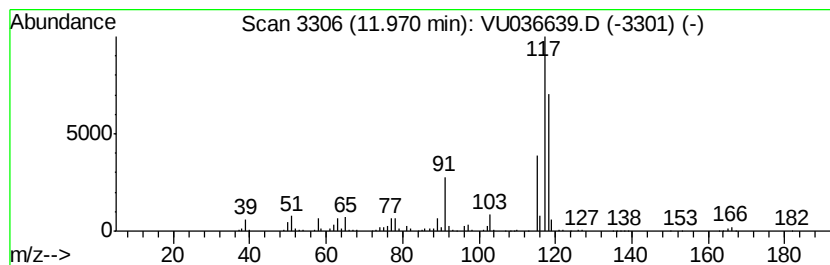
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Benzene, 1-propenyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.97	11.31 ug/L	220148	1,4-Dichlorobenzene-d4	11.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-propenyl-	118	C9H10	000637-50-3	95
2			Benzene, 1-propenyl-	118	C9H10	000637-50-3	94
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	94
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	93
5			Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00uL/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

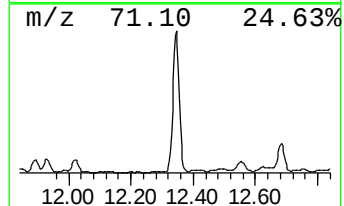
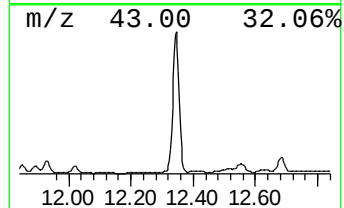
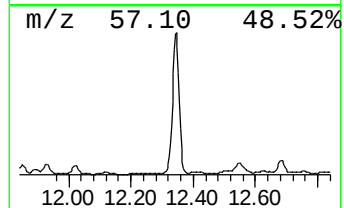
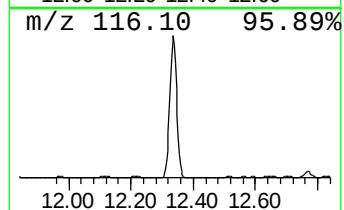
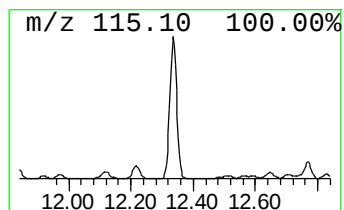
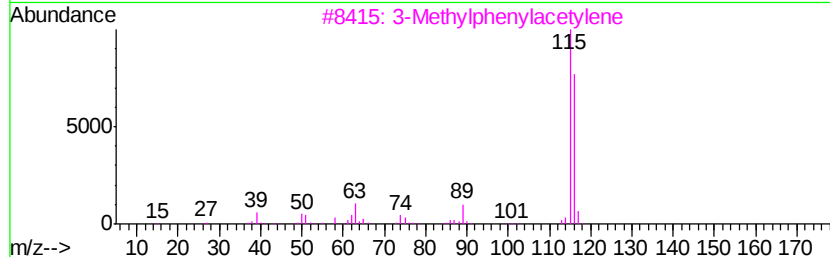
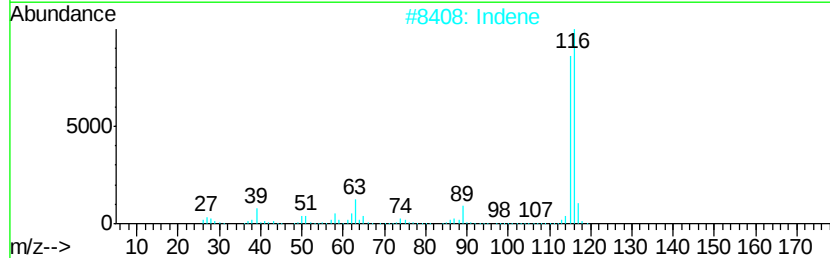
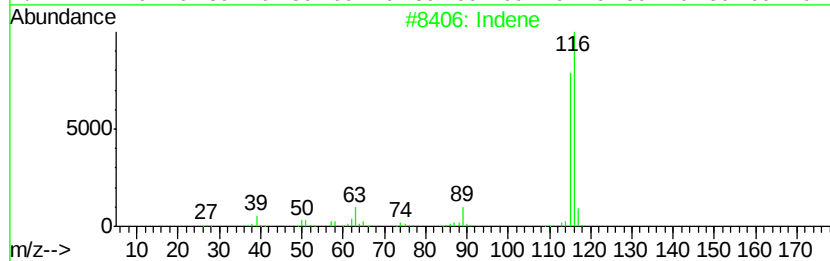
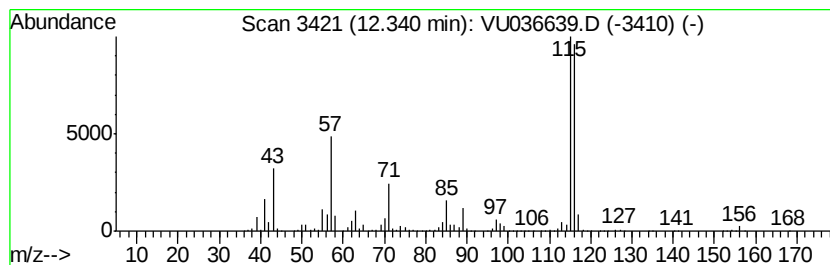
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Indene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.34	305.46 ug/L	5943980	1,4-Dichlorobenzene-d4	11.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indene	116	C9H8	000095-13-6	93
2		Indene	116	C9H8	000095-13-6	93
3		3-Methylphenylacetylene	116	C9H8	000766-82-5	81
4		Benzene, 1-propynyl-	116	C9H8	000673-32-5	81
5		Benzene, 1-propynyl-	116	C9H8	000673-32-5	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00uL/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

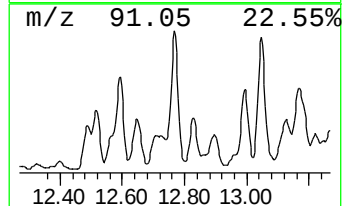
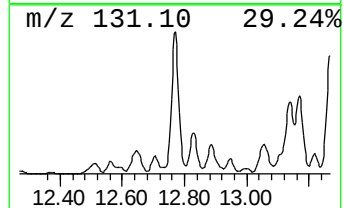
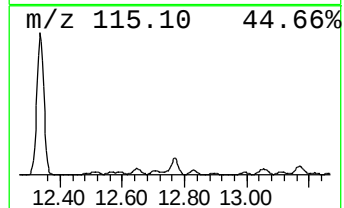
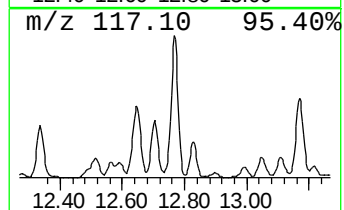
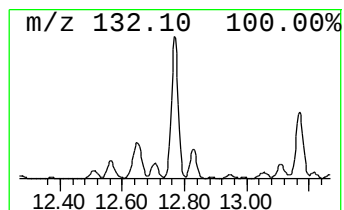
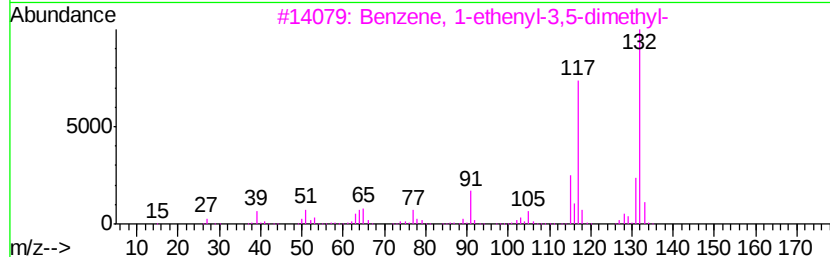
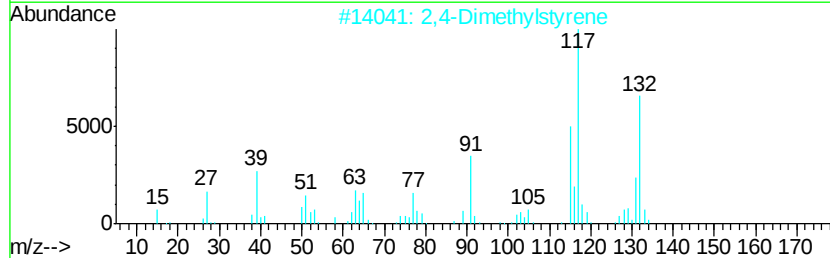
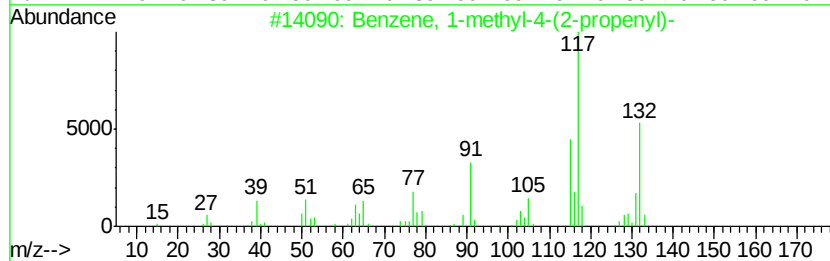
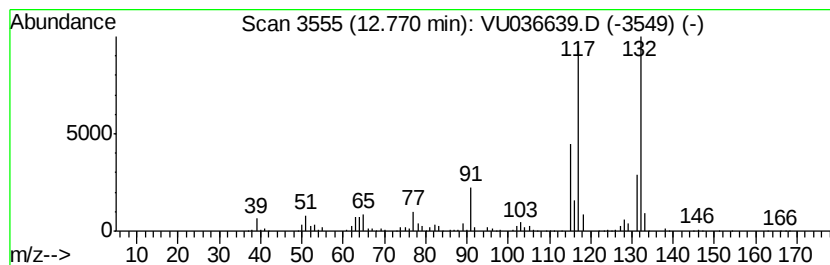
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Benzene, 1-methyl-4-(2-prop... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	62.03 ug/L	1206990	1,4-Dichlorobenzene-d4	11.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	94
2		2,4-Dimethylstyrene	132	C10H12	002234-20-0	94
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	94
4		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	94
5		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00uL/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

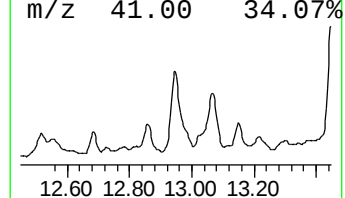
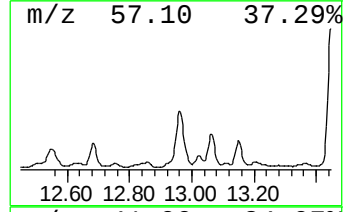
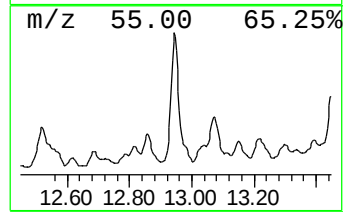
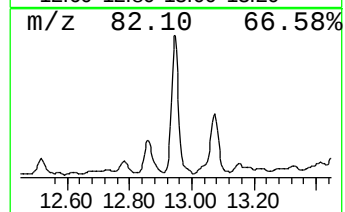
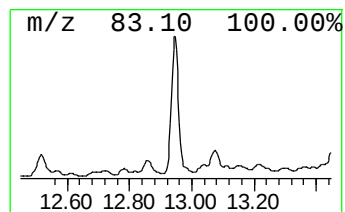
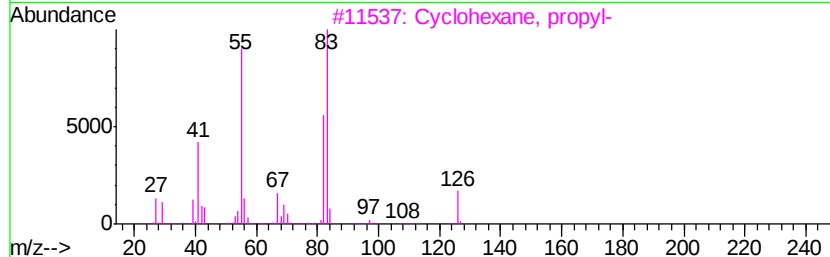
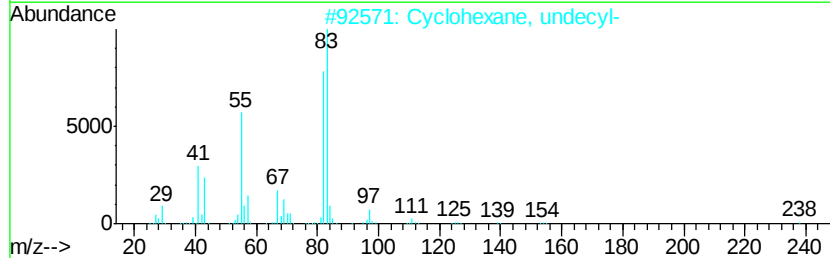
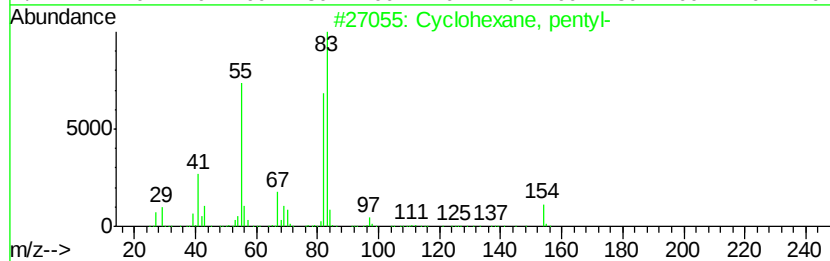
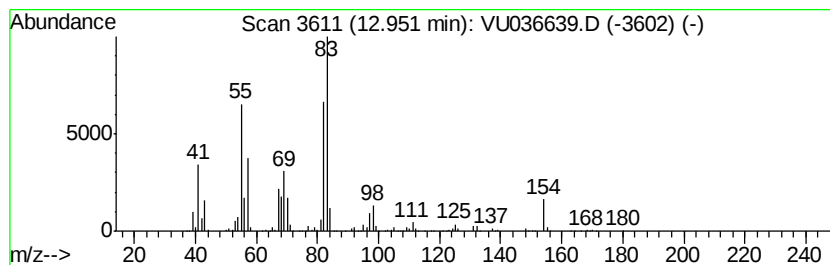
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 (DEL) Alkane: Cyclic12.95 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.95	71.64 ug/L	1393990	1,4-Dichlorobenzene-d4	11.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, pentyl-	154	C11H22	004292-92-6	74
2		Cyclohexane, undecyl-	238	C17H34	054105-66-7	64
3		Cyclohexane, propyl-	126	C9H18	001678-92-8	59
4		Cyclohexane, (4-methylpentyl)-	168	C12H24	061142-20-9	59
5		Cyclohexane, hexyl-	168	C12H24	004292-75-5	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00µL/5.0mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

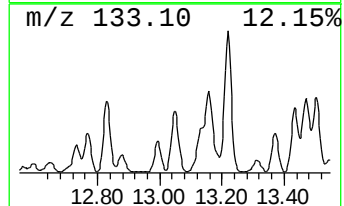
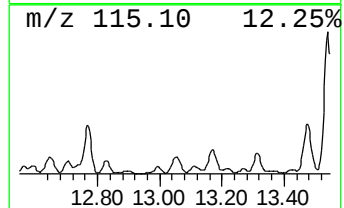
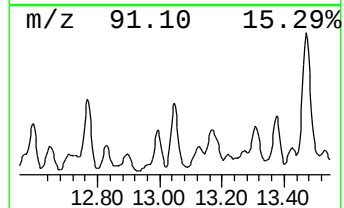
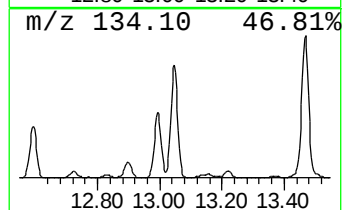
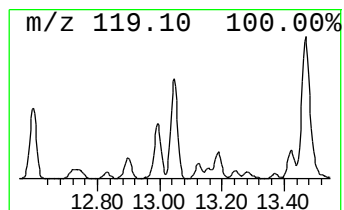
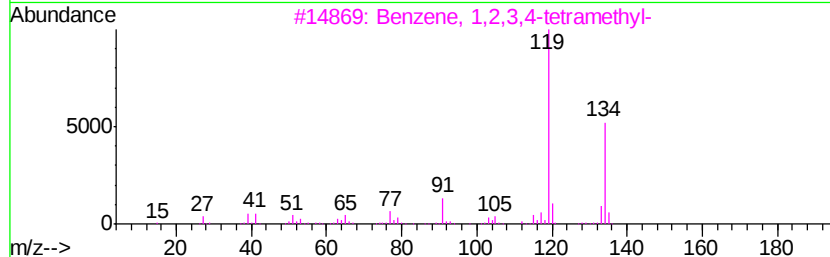
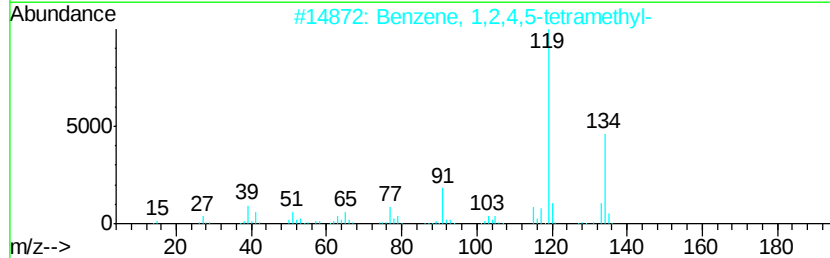
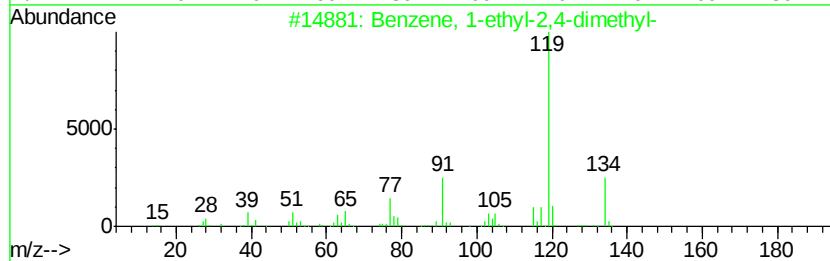
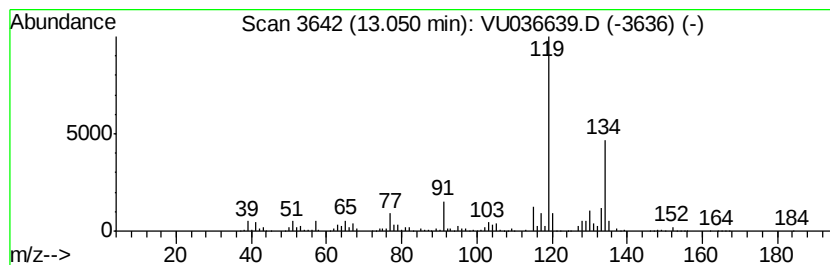
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.05	117.74 ug/L	2291110	1,4-Dichlorobenzene-d4	11.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00µ/5.0mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

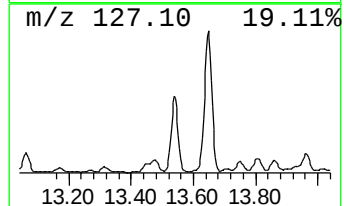
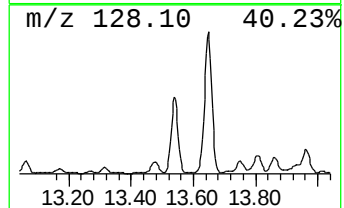
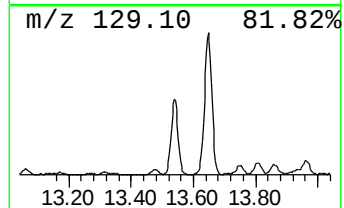
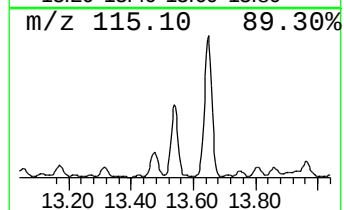
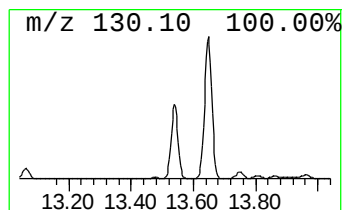
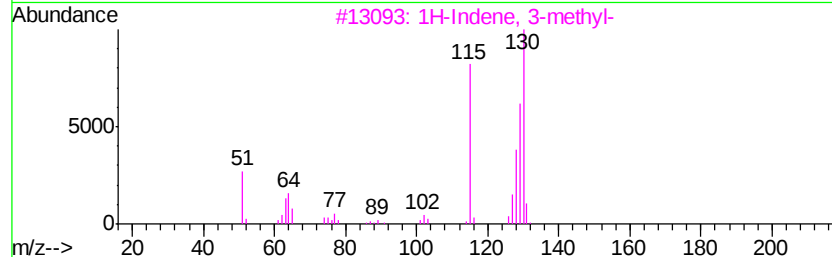
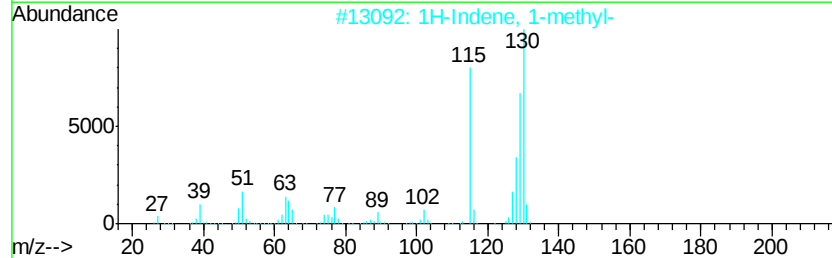
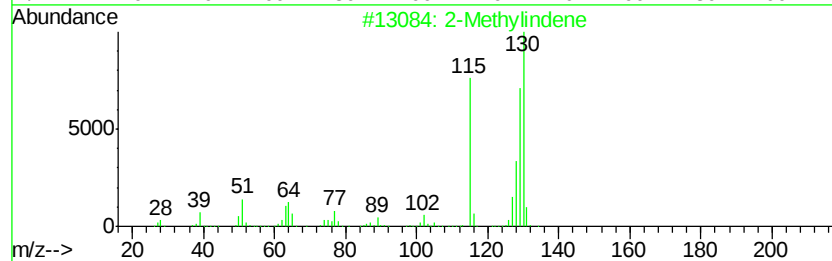
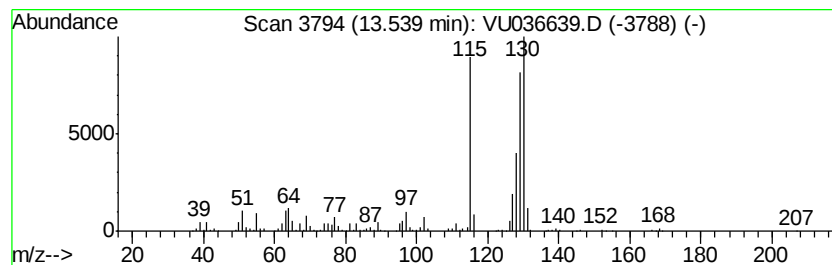
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 2-Methylindene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	107.84 ug/L	2098480	1,4-Dichlorobenzene-d4	11.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Methylindene	130	C10H10	002177-47-1	94
2		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	93
3		1H-Indene, 3-methyl-	130	C10H10	000767-60-2	91
4		Benzene, 1-butynyl-	130	C10H10	000622-76-4	91
5		1,4-Dihydronaphthalene	130	C10H10	000612-17-9	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00µ/5.0mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

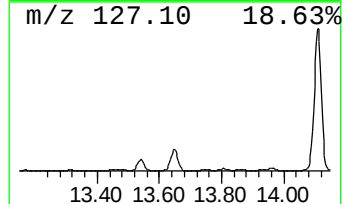
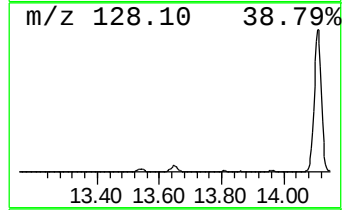
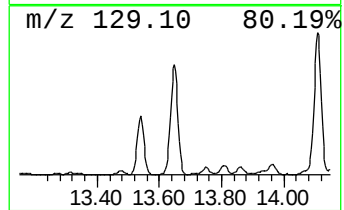
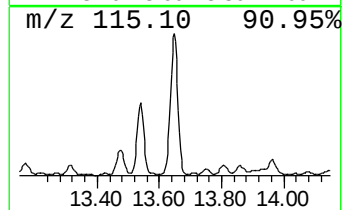
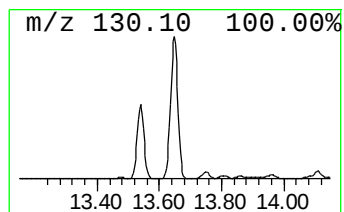
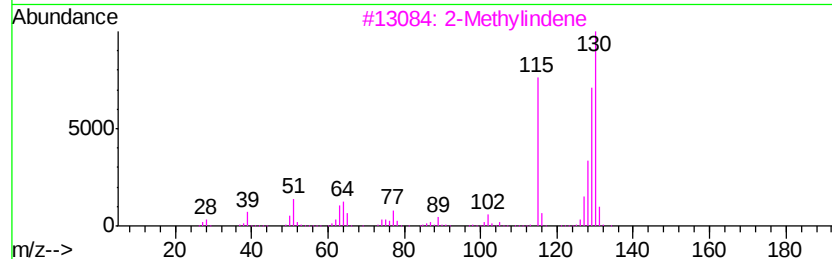
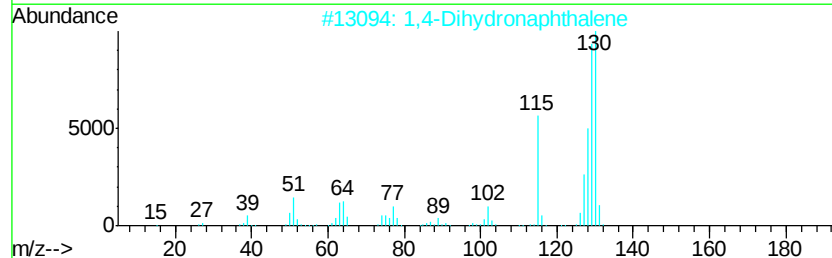
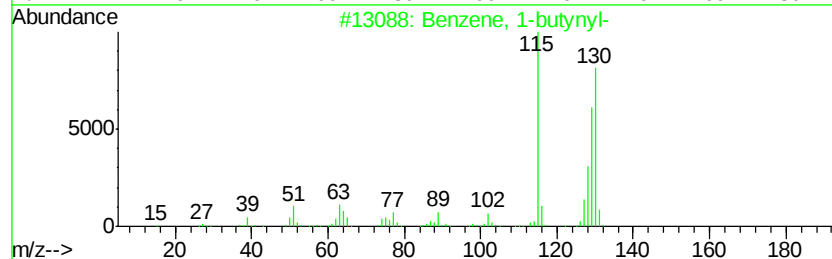
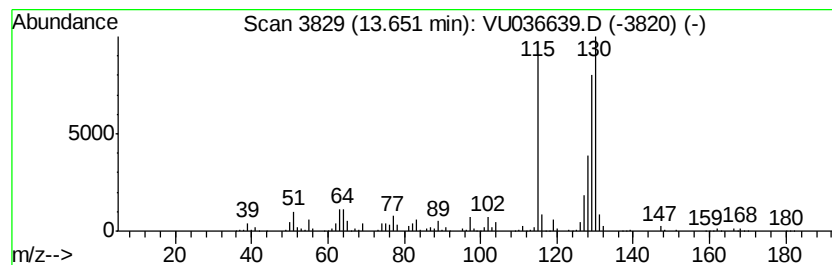
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 Benzene, 1-butylnyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.65	323.45 ug/L	6293990	1,4-Dichlorobenzene-d4	11.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-butylnyl-	130	C10H10	000622-76-4	95
2		1,4-Dihydronaphthalene	130	C10H10	000612-17-9	95
3		2-Methylindene	130	C10H10	002177-47-1	94
4		Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	94
5		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00µL/5.0mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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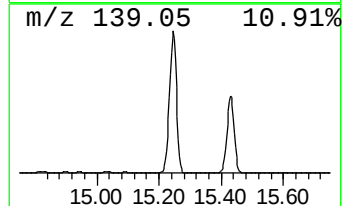
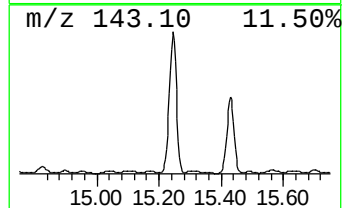
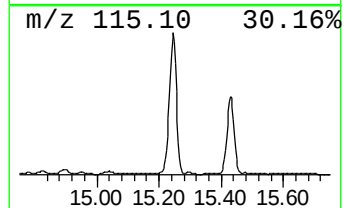
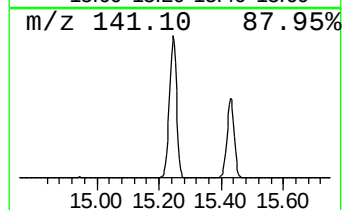
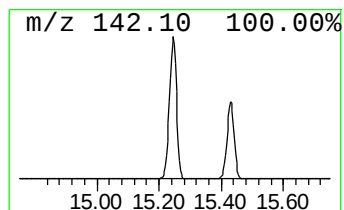
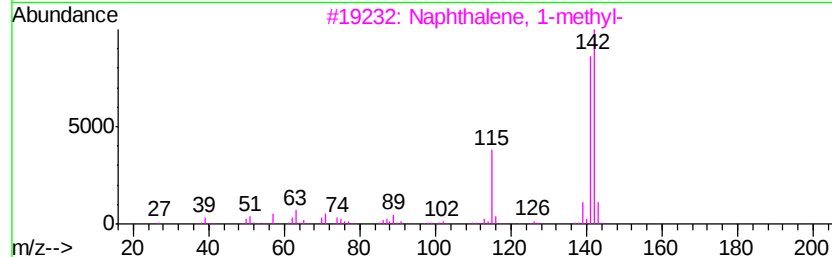
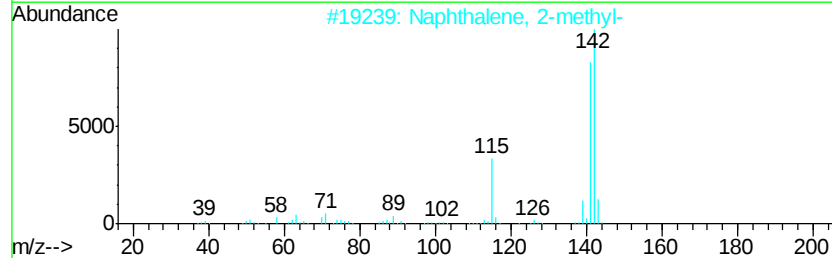
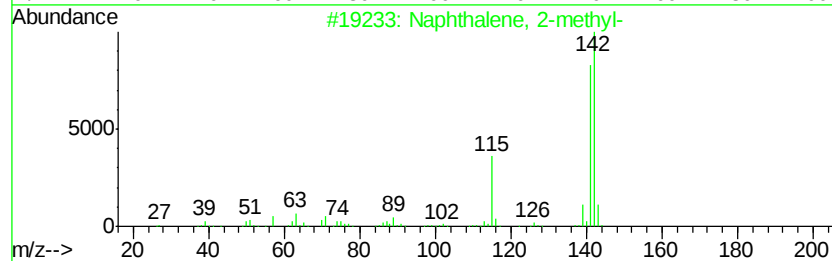
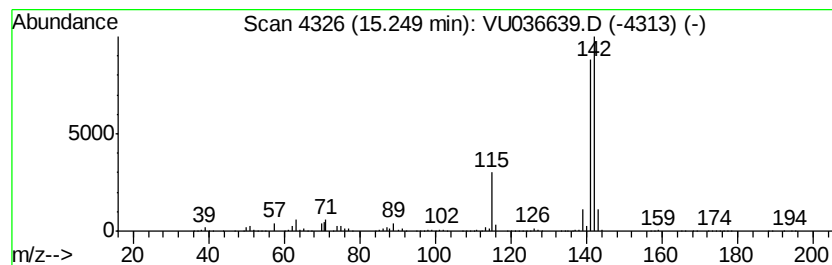
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 23 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.25	1398.14 ug/L	27206400	1,4-Dichlorobenzene-d4	11.84

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00µ/5.0mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

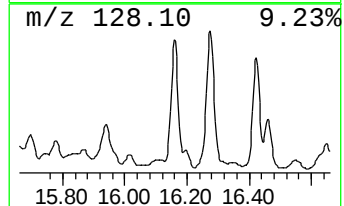
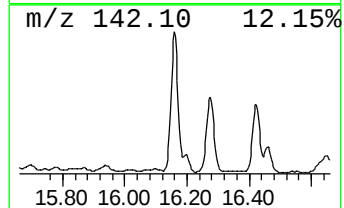
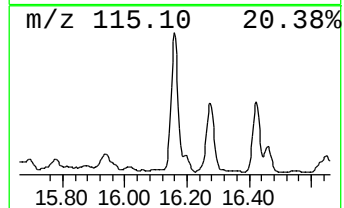
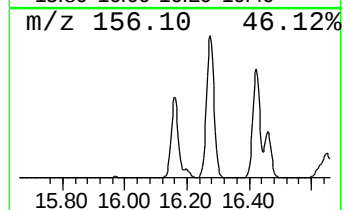
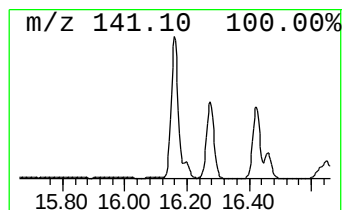
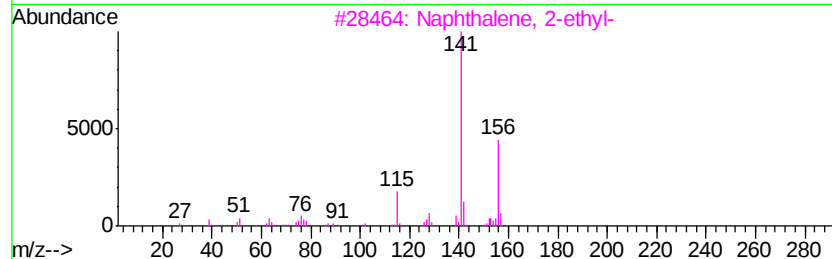
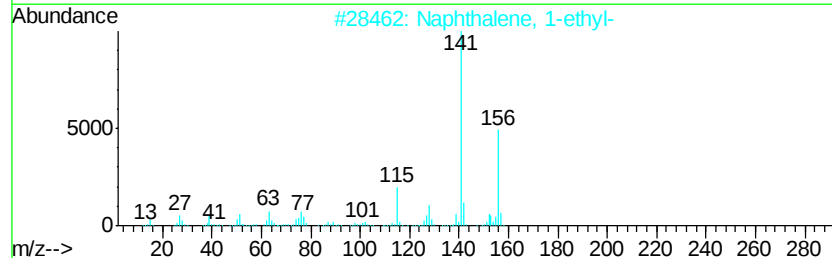
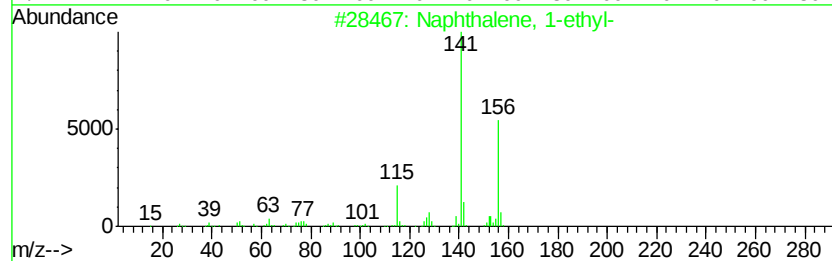
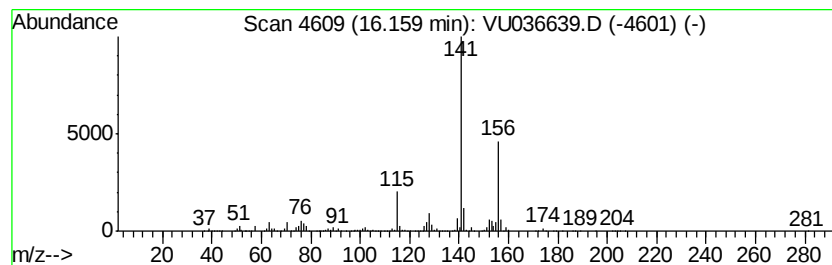
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 25 Naphthalene, 1-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.16	116.14 ug/L	2259910	1,4-Dichlorobenzene-d4	11.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	96
2		Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	96
3		Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	96
4		Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	96
5		Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00µ/5.0mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

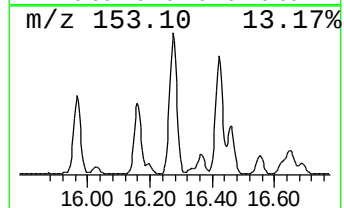
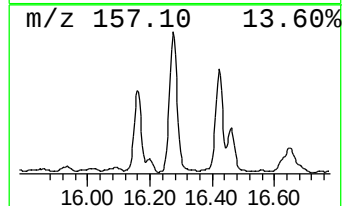
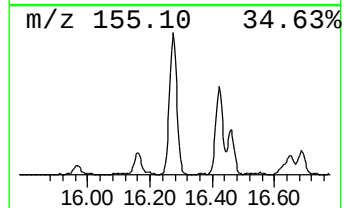
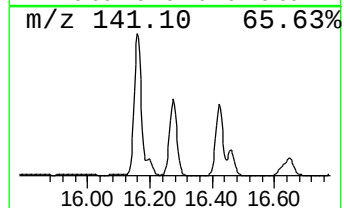
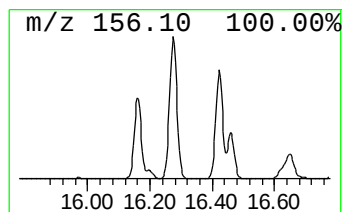
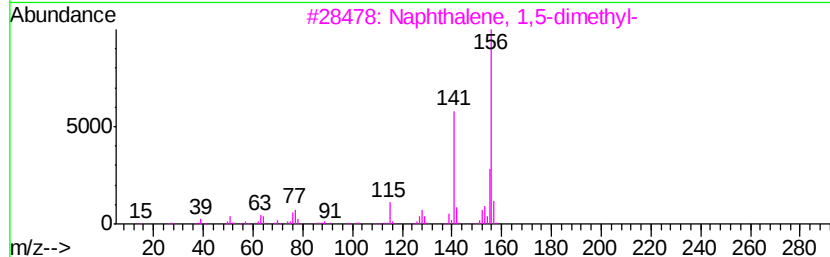
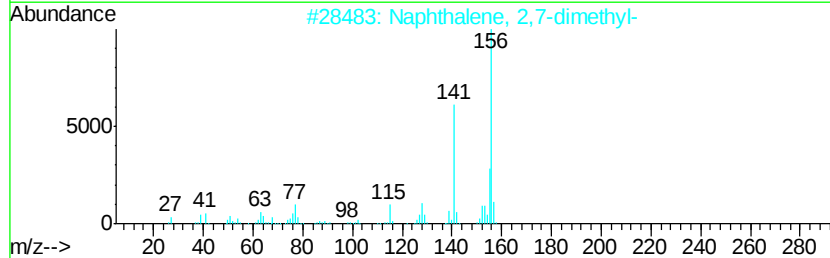
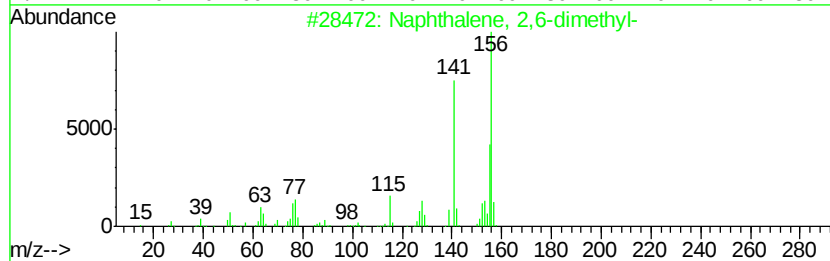
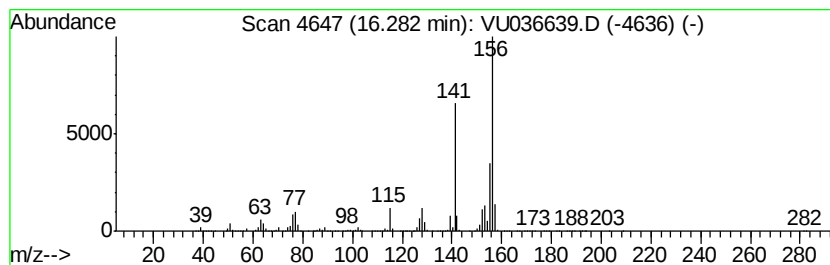
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 26 Naphthalene, 2,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.28	143.05 ug/L	2783640	1,4-Dichlorobenzene-d4	11.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
4		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00µL/5.0mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

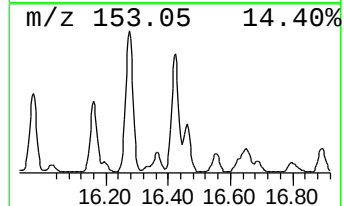
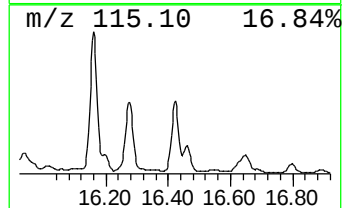
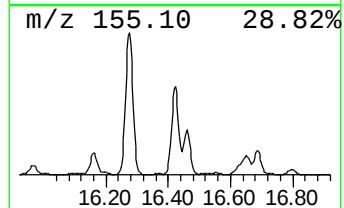
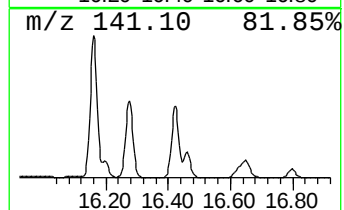
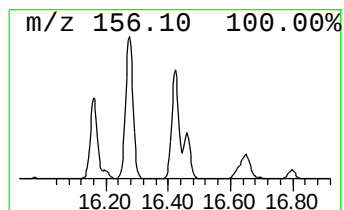
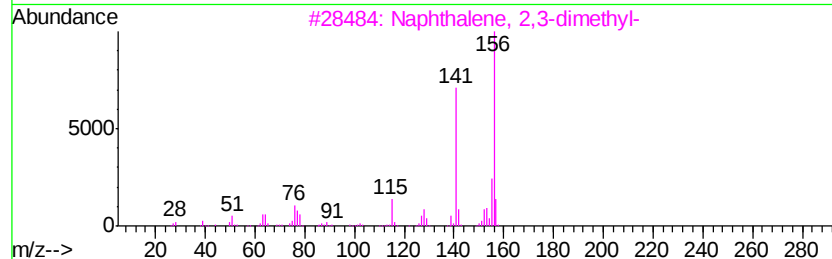
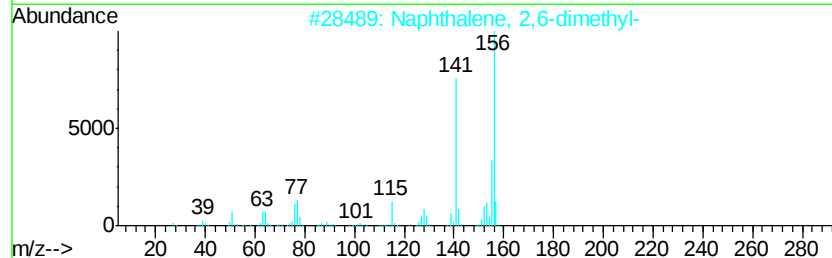
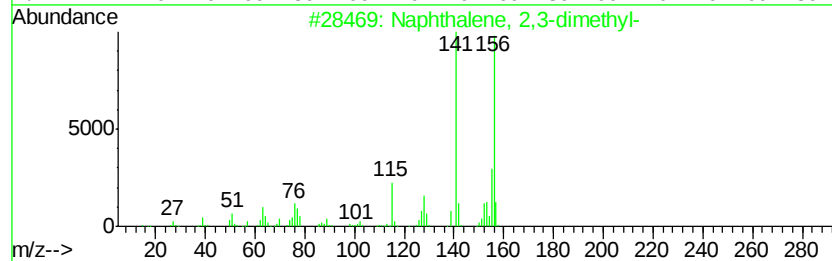
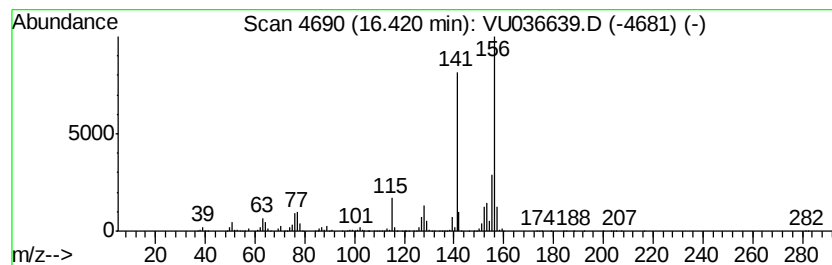
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 27 Naphthalene, 2,3-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.42	104.95 ug/L	2042240	1,4-Dichlorobenzene-d4	11.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
5			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00µL/5.0mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

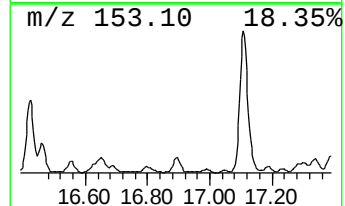
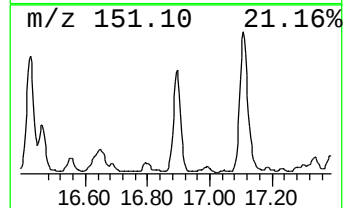
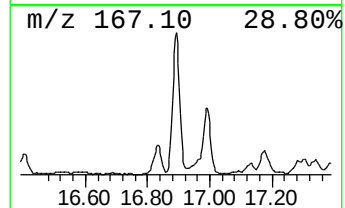
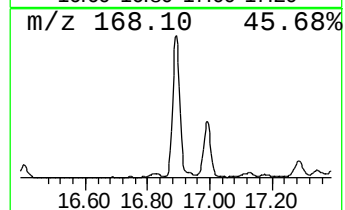
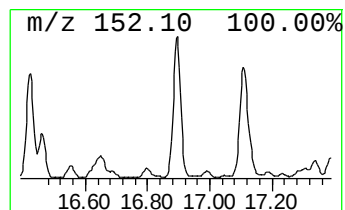
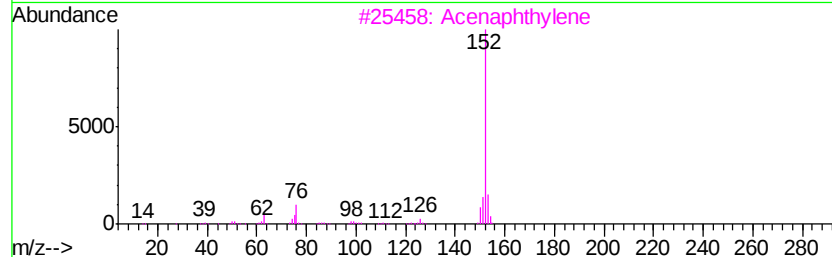
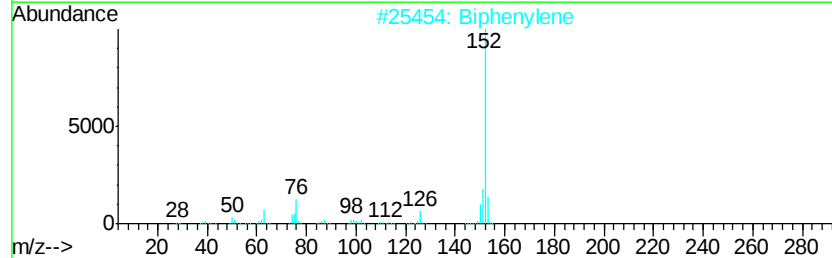
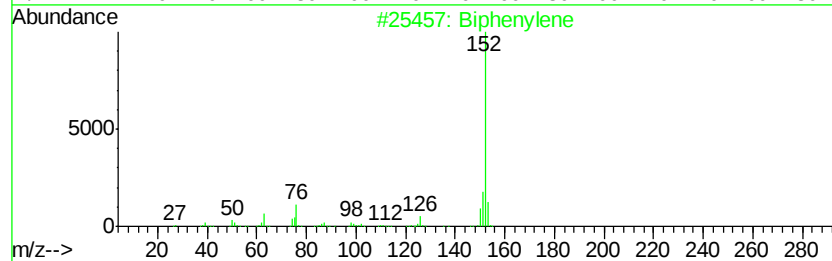
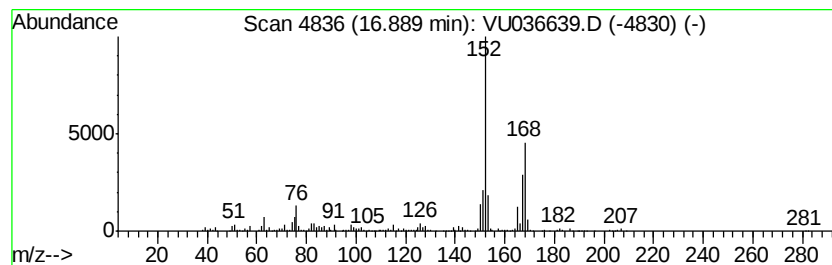
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 28 Biphenylene Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.89	15.68 ug/L	305047	1,4-Dichlorobenzene-d4	11.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Biphenylene	152	C12H8	000259-79-0	90
2			Biphenylene	152	C12H8	000259-79-0	78
3			Acenaphthylene	152	C12H8	000208-96-8	55
4			1-acenaphthenol, trifluoroacetat...	266	C14H9F3O2	1000376-45-2	49
5			(6Balpha,7beta,11alpha,11aalpha)...	334	C25H18O	1000241-69-8	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00µL/5.0mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

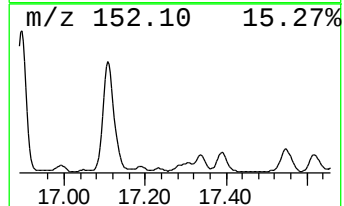
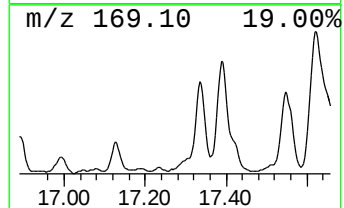
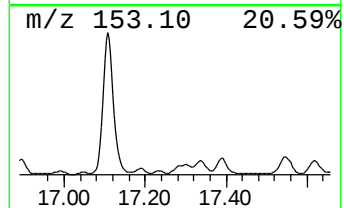
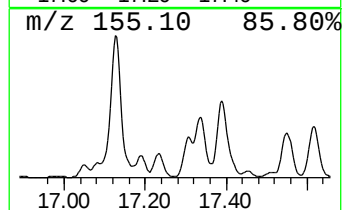
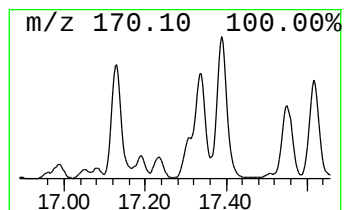
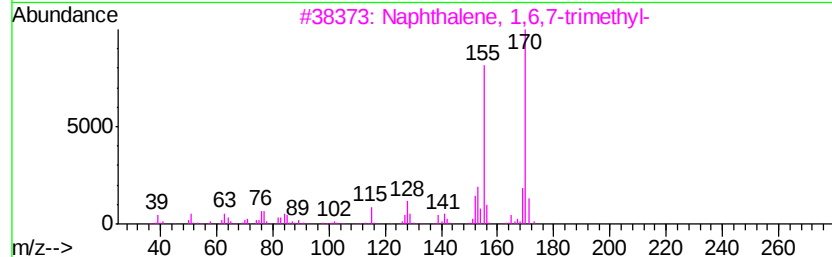
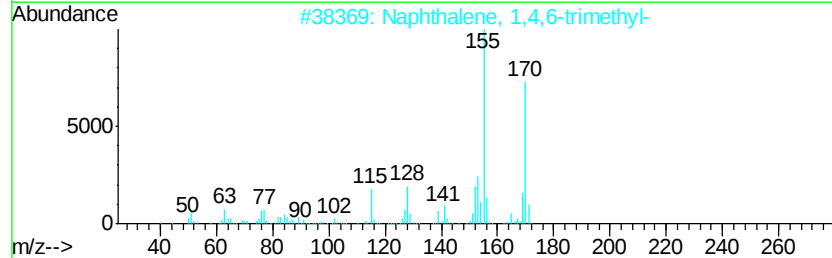
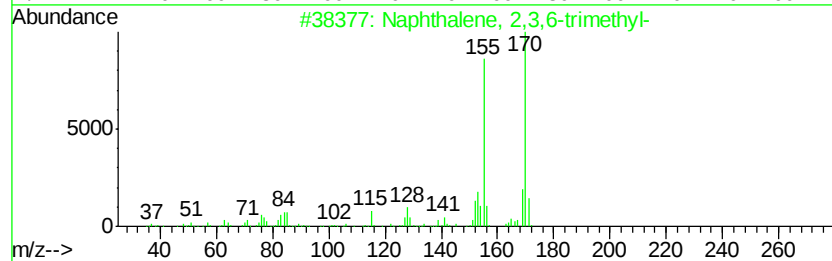
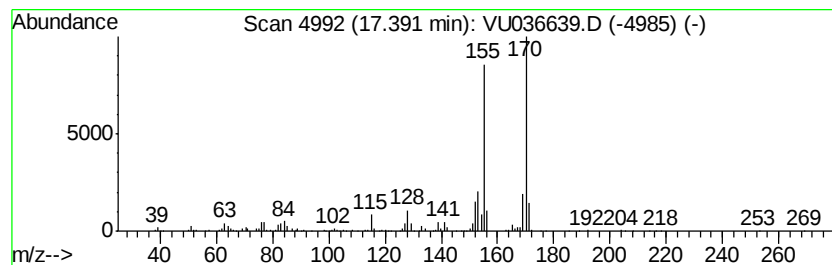
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 30 Naphthalene, 2,3,6-trimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.39	20.15 ug/L	392102	1,4-Dichlorobenzene-d4	11.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
2		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	96
3		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
4		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	95
5		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU013020\
Data File : VU036639.D
Acq On : 31 Jan 2020 00:11
Operator : JC/MD
Sample : L1324-08
Misc : 5.00µ/5.0mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAT65

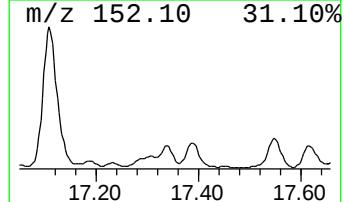
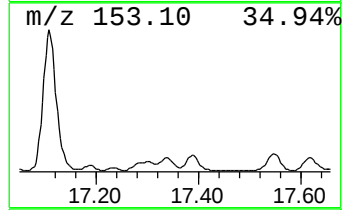
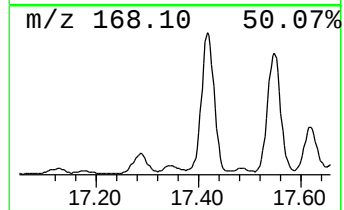
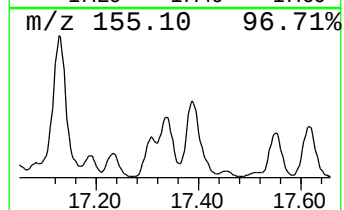
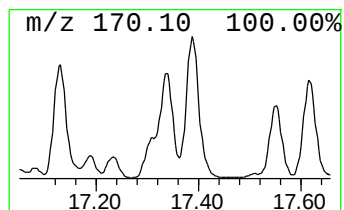
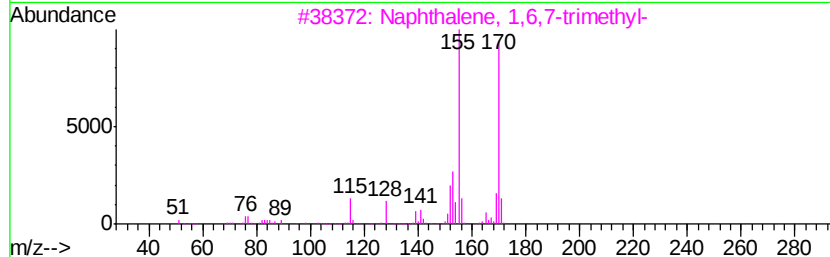
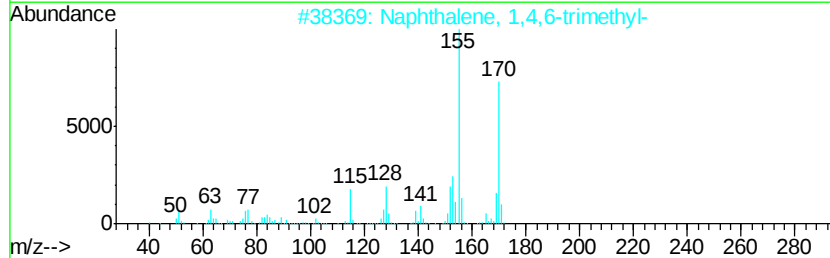
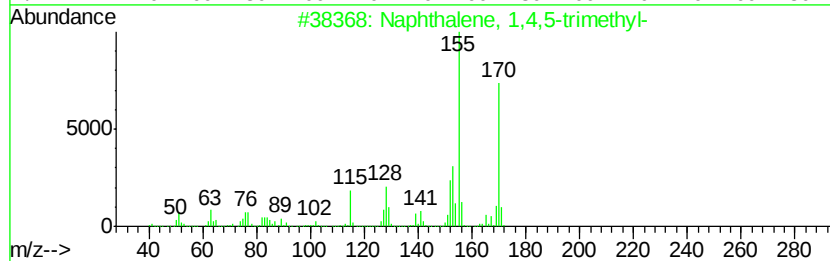
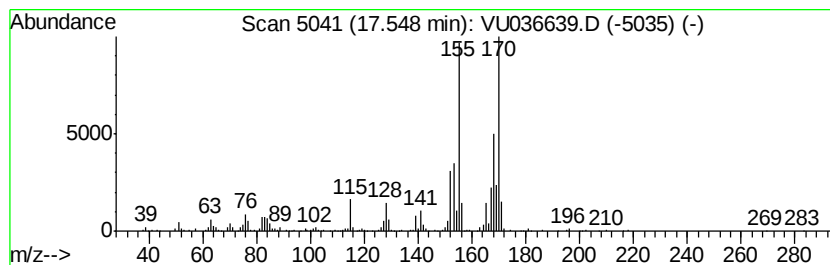
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 31 Naphthalene, 1,4,5-trimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.55	13.00 ug/L	252944	1,4-Dichlorobenzene-d4	11.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	95
2			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	94
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
4			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	89



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU013020\
 Data File : VU036639.D
 Acq On : 31 Jan 2020 00:11
 Operator : JC/MD
 Sample : L1324-08
 Misc : 5.00g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 GAT65

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM012920WMA.M
 Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Cyc...	10.06	2.5	ug/L	40414	2	9.45	791489	50.0
(DEL) Alkane: Str...	10.27	7.1	ug/L	111949	2	9.45	791489	50.0
(DEL) Alkane: Cyc...	10.38	2.6	ug/L	41701	2	9.45	791489	50.0
(DEL) Alkane: Str...	10.65	2.5	ug/L	49554	3	11.84	972951	50.0
Benzene, 1-ethyl-...	11.05	73.5	ug/L	1430390	3	11.84	972951	50.0
Benzene, 1,2,3-tr...	11.12	87.9	ug/L	1709860	3	11.84	972951	50.0
(DEL) Alkane: Str...	11.44	10.0	ug/L	195177	3	11.84	972951	50.0
Benzene, 1-etheny...	11.56	115.9	ug/L	2254580	3	11.84	972951	50.0
(DEL) Alkane: Cyc...	11.73	11.1	ug/L	215501	3	11.84	972951	50.0
Benzene, 1-propenyl-	11.97	11.3	ug/L	220148	3	11.84	972951	50.0
Indene	12.34	305.5	ug/L	5943980	3	11.84	972951	50.0
Benzene, 1-methyl...	12.77	62.0	ug/L	1206990	3	11.84	972951	50.0
(DEL) Alkane: Cyc...	12.95	71.6	ug/L	1393990	3	11.84	972951	50.0
Benzene, 1-ethyl-...	13.05	117.7	ug/L	2291110	3	11.84	972951	50.0
2-Methylindene	13.54	107.8	ug/L	2098480	3	11.84	972951	50.0
Benzene, 1-butyryl-	13.65	323.4	ug/L	6293990	3	11.84	972951	50.0
Naphthalene, 1-me...	15.25	1398.1	ug/L	27206400	3	11.84	972951	50.0
Naphthalene, 1-et...	16.16	116.1	ug/L	2259910	3	11.84	972951	50.0
Naphthalene, 2,6-...	16.28	143.1	ug/L	2783640	3	11.84	972951	50.0
Naphthalene, 2,3-...	16.42	105.0	ug/L	2042240	3	11.84	972951	50.0
Biphenylene	16.89	15.7	ug/L	305047	3	11.84	972951	50.0
Naphthalene, 2,3,...	17.39	20.1	ug/L	392102	3	11.84	972951	50.0
Naphthalene, 1,4,...	17.55	13.0	ug/L	252944	3	11.84	972951	50.0