

Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
 Data File : VU031676.D
 Acq On : 01 May 2019 12:00
 Operator : JC/SP
 Sample : K2604-06DL 25X
 Misc : 5.01g/5mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 GAJ80DL

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\SOMULM042719WMA.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	1.396	88	95	108	rVB	458171	488898	0.91%	0.298%
2	1.679	176	183	196	rBV	393186	477083	0.89%	0.291%
3	2.270	356	367	381	rBV	794365	1207684	2.24%	0.737%
4	2.473	427	430	437	rVB5	2188	2396	0.00%	0.001%
5	2.698	497	500	509	rVB6	3953	4740	0.01%	0.003%
6	2.888	556	559	562	rVB2	1555	867	0.00%	0.001%
7	2.984	585	589	591	rVB3	1350	887	0.00%	0.001%
8	3.019	595	600	604	rBV5	1758	1824	0.00%	0.001%
9	3.093	620	623	627	rVB2	1490	1043	0.00%	0.001%
10	3.299	681	687	692	rBV6	2320	3463	0.01%	0.002%
11	3.367	703	708	711	rBV2	980	920	0.00%	0.001%
12	3.441	728	731	736	rBV3	977	885	0.00%	0.001%
13	3.608	780	783	785	rBV3	1607	1173	0.00%	0.001%
14	3.656	795	798	804	rVB5	984	979	0.00%	0.001%
15	3.730	818	821	824	rVB	1632	978	0.00%	0.001%
16	3.791	835	840	844	rBV3	1043	1050	0.00%	0.001%
17	3.843	849	856	860	rVB3	1678	1660	0.00%	0.001%
18	3.875	860	866	871	rBV5	1319	1740	0.00%	0.001%
19	3.913	876	878	881	rBV2	1241	833	0.00%	0.001%
20	3.974	892	897	899	rBV3	1249	997	0.00%	0.001%
21	4.129	942	945	949	rBV3	1075	794	0.00%	0.000%
22	4.187	949	963	999	rBV	366687	1166363	2.16%	0.712%
23	4.643	1089	1105	1128	rVV	569159	1370762	2.54%	0.836%
24	5.074	1237	1239	1242	rBV	1826	1028	0.00%	0.001%
25	5.132	1253	1257	1259	rBV4	1351	1034	0.00%	0.001%
26	5.170	1267	1269	1274	rVB	1554	1151	0.00%	0.001%
27	5.334	1296	1320	1363	rBV2	1209184	3796025	7.04%	2.316%
28	5.749	1447	1449	1453	rVB2	1743	935	0.00%	0.001%
29	5.768	1453	1455	1459	rBV6	1478	1096	0.00%	0.001%
30	5.814	1465	1469	1473	rBV5	1736	1273	0.00%	0.001%
31	5.881	1476	1490	1512	rBV	1163273	2360966	4.38%	1.440%
32	6.177	1574	1582	1595	rBV10	13744	28302	0.05%	0.017%
33	6.325	1614	1628	1654	rBV	809978	1680928	3.12%	1.025%
34	6.508	1681	1685	1686	rBV4	1611	1276	0.00%	0.001%

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Integration Parameters: LSCINT.P

Integrator: RTE
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Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM042719WMA.M
 Title : VOC Analysis

35	6.727	1749	1753	1757	rVB7	1431	1373	0.00%	0.001%
36	6.752	1757	1761	1765	rBV4	1340	1372	0.00%	0.001%
37	6.813	1777	1780	1783	rBV3	1121	791	0.00%	0.000%
38	6.839	1783	1788	1789	rBV4	1306	1214	0.00%	0.001%
39	6.871	1794	1798	1801	rBV4	1099	1007	0.00%	0.001%
40	6.894	1801	1805	1808	rBV4	2245	1825	0.00%	0.001%
41	6.920	1808	1813	1814	rBV5	1064	912	0.00%	0.001%
42	7.029	1844	1847	1851	rBV4	1694	1077	0.00%	0.001%
43	7.100	1865	1869	1871	rBV4	1575	1111	0.00%	0.001%
44	7.132	1876	1879	1884	rBV2	1188	772	0.00%	0.000%
45	7.154	1884	1886	1890	rVB3	1104	830	0.00%	0.001%
46	7.225	1893	1908	1941	rBV	448388	862846	1.60%	0.526%
47	7.502	1992	1994	1997	rBV5	1223	785	0.00%	0.000%
48	7.563	1998	2013	2029	rBV	1477584	2730965	5.07%	1.666%
49	7.849	2091	2102	2128	rBV	298859	556365	1.03%	0.339%
50	8.309	2233	2245	2294	rBV	1400729	2622605	4.87%	1.600%
51	8.710	2367	2370	2372	rBV3	2597	1551	0.00%	0.001%
52	8.801	2396	2398	2400	rBV3	1449	733	0.00%	0.000%
53	8.813	2400	2402	2404	rVV3	1671	836	0.00%	0.001%
54	8.849	2410	2413	2419	rVB6	3212	2521	0.00%	0.002%
55	9.087	2475	2487	2526	rBV	1709793	2971473	5.51%	1.813%
56	9.247	2527	2537	2553	rBV	222258	374825	0.70%	0.229%
57	9.370	2563	2575	2592	rBV	1564386	2673814	4.96%	1.631%
58	9.704	2677	2679	2682	rBV4	2515	1927	0.00%	0.001%
59	9.788	2690	2705	2738	rBV3	2607655	4877379	9.05%	2.975%
60	10.427	2892	2904	2918	rBV	1124448	1819174	3.37%	1.110%
61	10.585	2946	2953	2965	rVB	51592	77241	0.14%	0.047%
62	10.678	2972	2982	2998	rBV	220539	490941	0.91%	0.299%
63	10.768	3000	3010	3018	rBV2	283798	452410	0.84%	0.276%
64	10.993	3064	3080	3092	rBV3	179957	401509	0.74%	0.245%
65	11.144	3110	3127	3138	rBV	977024	1615533	3.00%	0.986%
66	11.212	3138	3148	3157	rVV	1594745	2431369	4.51%	1.483%
67	11.260	3157	3163	3177	rVB	593007	918345	1.70%	0.560%
68	11.479	3221	3231	3249	rBV	1855526	2895434	5.37%	1.766%
69	11.566	3250	3258	3268	rVV	330715	497635	0.92%	0.304%
70	11.624	3268	3276	3293	rVB2	223814	360297	0.67%	0.220%
71	11.762	3309	3319	3328	rBV	217090	372100	0.69%	0.227%

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 Title : VOC Analysis

72	11.855	3338	3348	3365	rVB	1626437	2705809	5.02%	1.651%
73	11.977	3374	3386	3397	rBV	6081229	9525798	17.67%	5.811%
74	12.148	3432	3439	3440	rBV2	43275	47954	0.09%	0.029%
75	12.222	3455	3462	3466	rBV3	92004	137261	0.25%	0.084%
76	12.421	3516	3524	3535	rVB	387026	595504	1.10%	0.363%
77	12.482	3536	3543	3556	rVV4	102886	173060	0.32%	0.106%
78	12.704	3603	3612	3626	rBV4	268194	622351	1.15%	0.380%
79	12.820	3640	3648	3658	rBV	291617	476982	0.88%	0.291%
80	12.961	3686	3692	3702	rVB2	78031	110883	0.21%	0.068%
81	13.119	3733	3741	3750	rBV3	413948	619633	1.15%	0.378%
82	13.180	3751	3760	3780	rVB	1283148	1976936	3.67%	1.206%
83	13.286	3782	3793	3810	rVB2	1451338	2444371	4.53%	1.491%
84	13.739	3920	3934	3960	rBV2	33673105	53905636	100.00%	32.884%
85	13.858	3964	3971	3990	rVB2	343649	700151	1.30%	0.427%
86	14.135	4051	4057	4065	rBV	200895	273789	0.51%	0.167%
87	14.334	4113	4119	4128	rBV4	184840	260462	0.48%	0.159%
88	14.871	4275	4286	4311	rVB	10655318	16525856	30.66%	10.081%
89	15.054	4332	4343	4362	rBV	5939338	9799360	18.18%	5.978%
90	15.601	4505	4513	4523	rVB	908130	1348999	2.50%	0.823%
91	15.794	4566	4573	4581	rBV	546707	770293	1.43%	0.470%
92	15.906	4597	4608	4622	rBV	2483815	4277211	7.93%	2.609%
93	16.054	4644	4654	4660	rBV	2359625	3679841	6.83%	2.245%
94	16.186	4686	4695	4707	rVB	769872	1211005	2.25%	0.739%
95	16.279	4709	4724	4745	rVB	800304	1942193	3.60%	1.185%
96	16.427	4762	4770	4781	rBV	388559	589442	1.09%	0.360%
97	16.517	4788	4798	4815	rBV	1815402	3202905	5.94%	1.954%
98	17.019	4946	4954	4976	rVB3	661299	1705968	3.16%	1.041%
99	17.160	4991	4998	5008	rVB	553288	842378	1.56%	0.514%
100	17.221	5010	5017	5028	rVV3	543786	820413	1.52%	0.500%

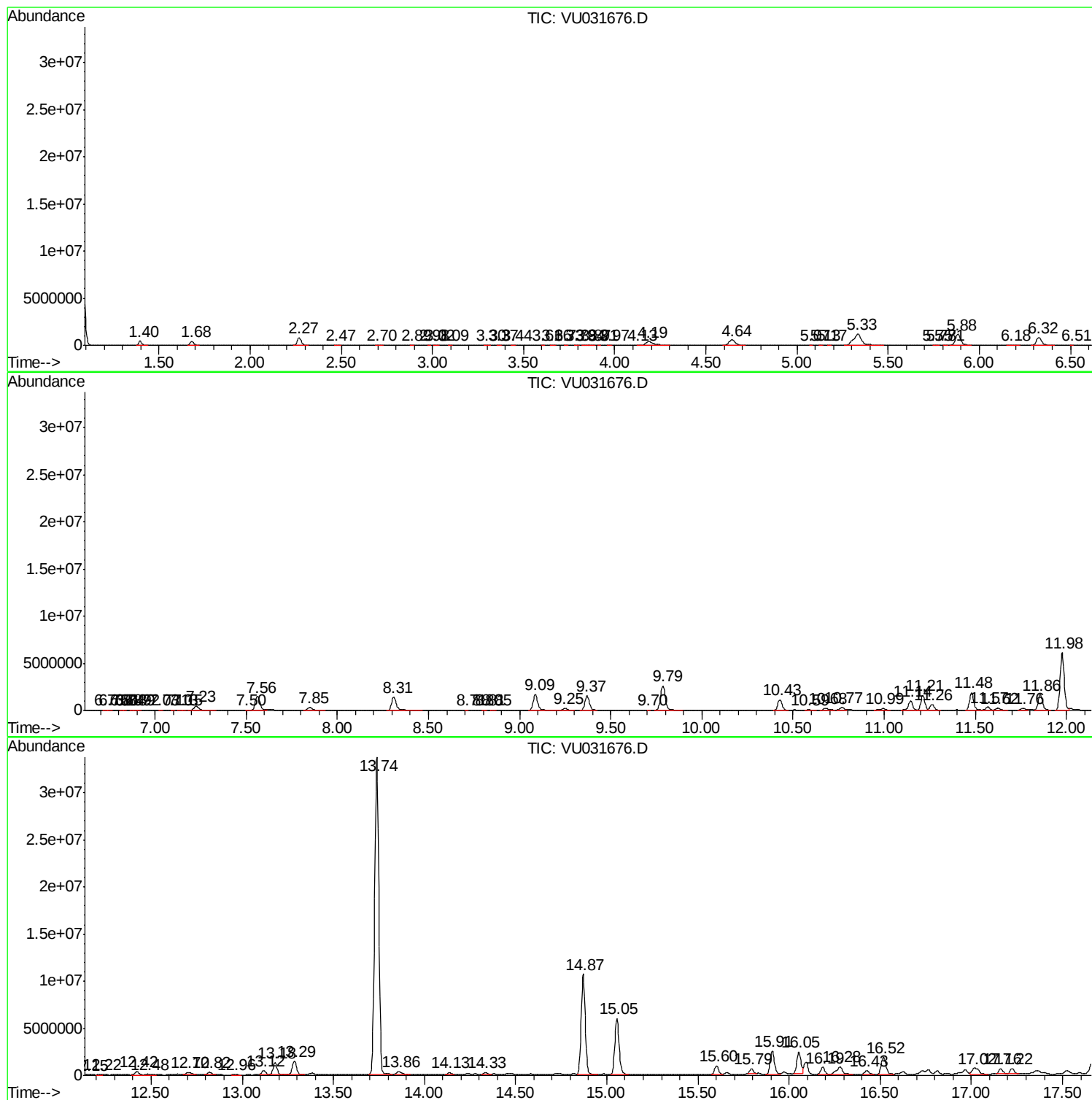
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Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
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ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAJ80DL

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM042719WMA.M
Quant Title : VOC Analysis

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



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ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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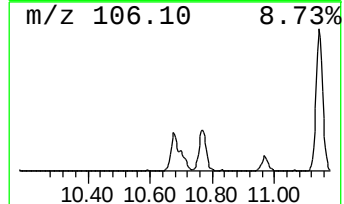
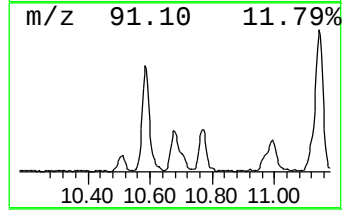
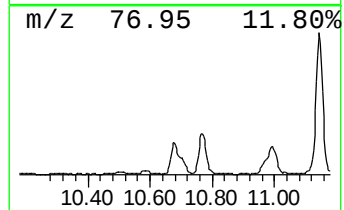
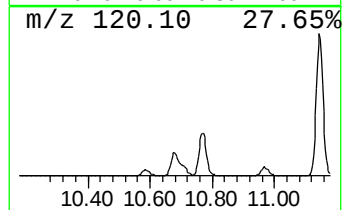
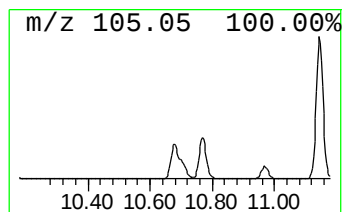
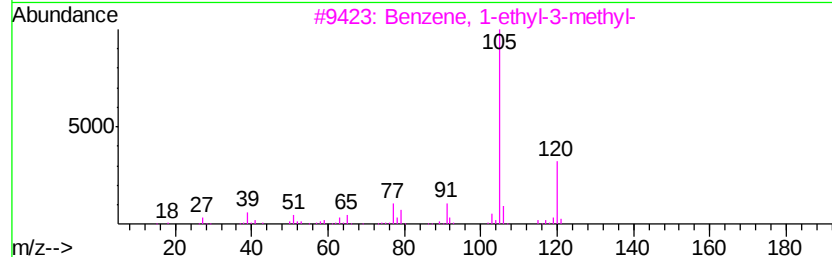
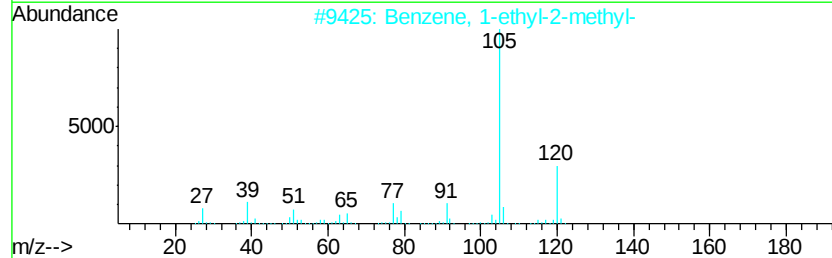
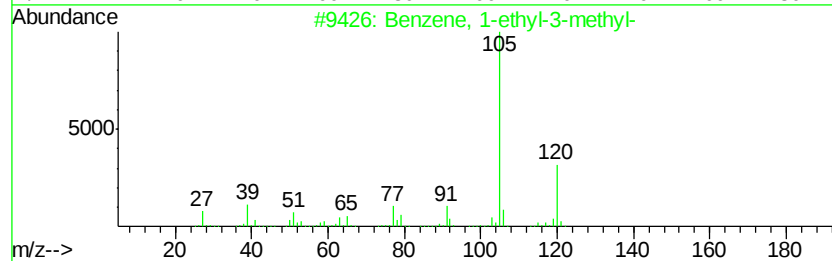
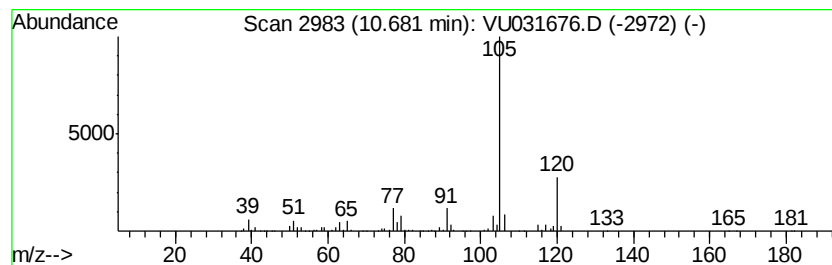
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM042719WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.68	8.48 ug/L	490941	1,4-Dichlorobenzene-d4	11.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94



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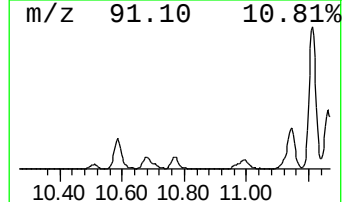
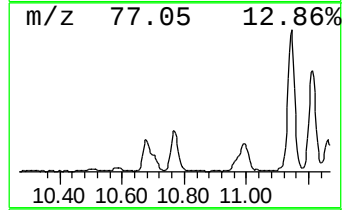
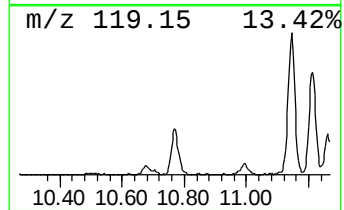
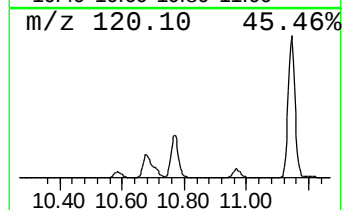
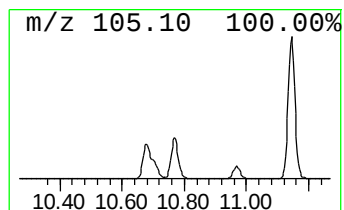
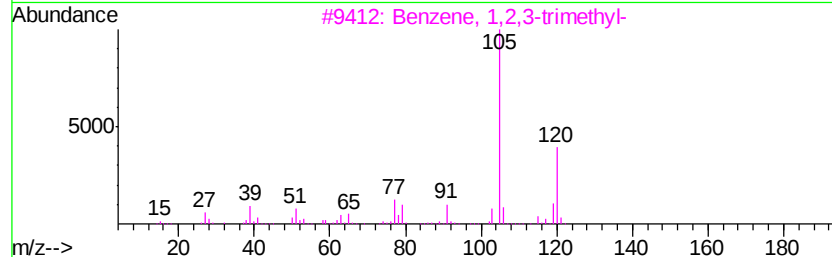
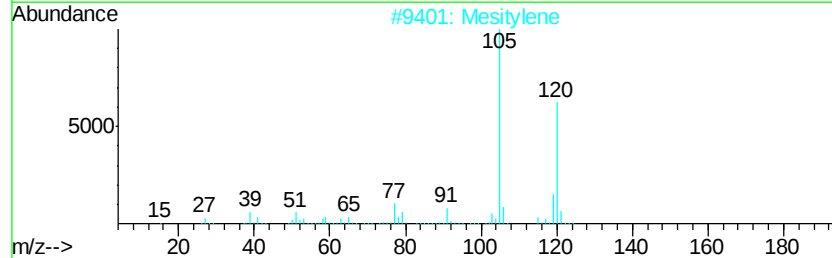
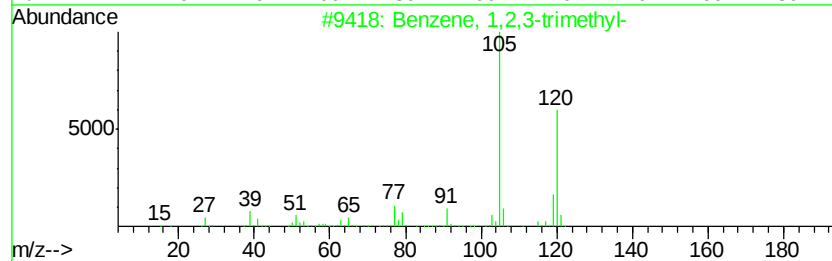
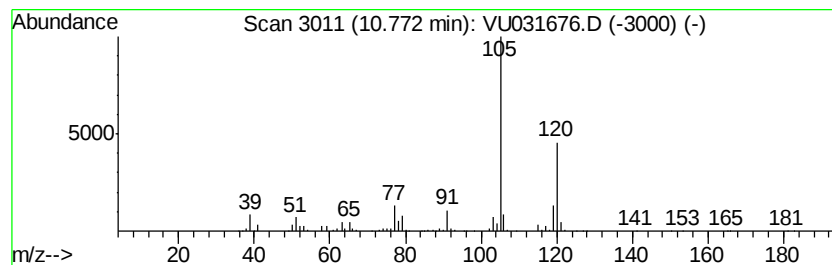
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM042719WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 Benzene, 1,2,3-trimethyl- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.77	7.81 ug/L	452410	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Mesitylene	120	C9H12	000108-67-8	95
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	93



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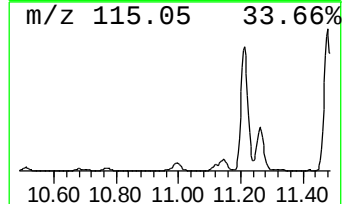
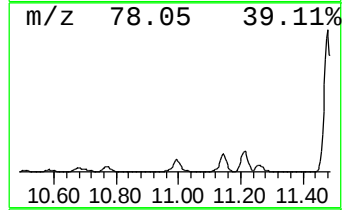
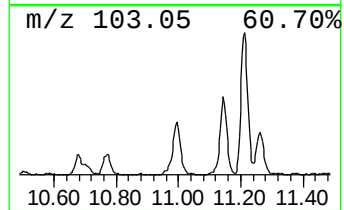
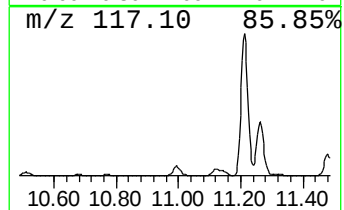
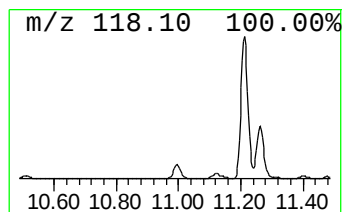
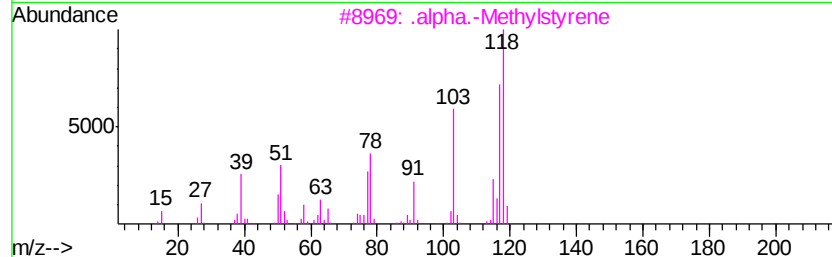
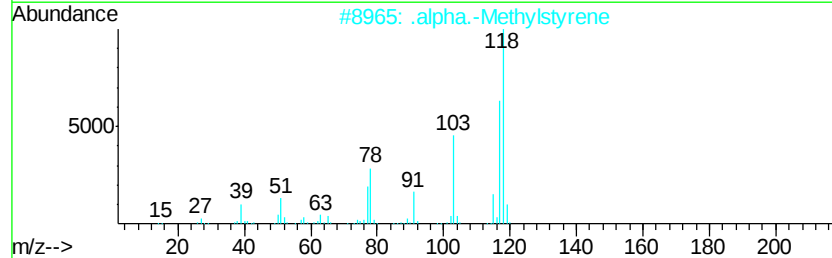
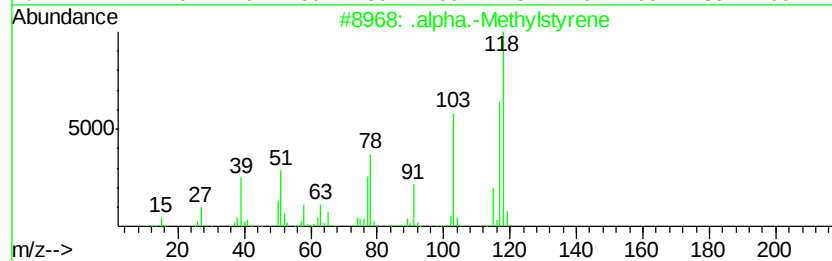
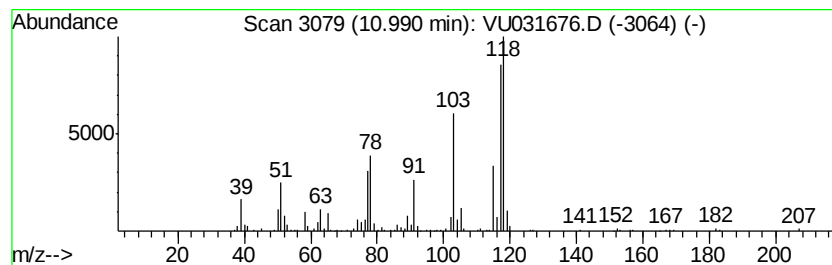
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ClientSampleId :
GAJ80DL

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM042719WMA.M
Quant Title : VOC Analysis

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

Peak Number 4 .alpha.-Methylstyrene Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.99	6.93 ug/L	401509	1,4-Dichlorobenzene-d4	11.48	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Methylstyrene	118	C9H10	000098-83-9	95
2	.alpha.-Methylstyrene	118	C9H10	000098-83-9	94
3	.alpha.-Methylstyrene	118	C9H10	000098-83-9	93
4	Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	022250-74-4	53
5	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAJ80DL

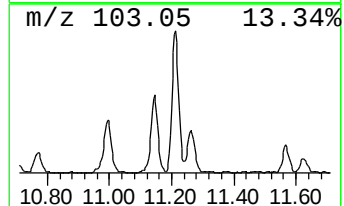
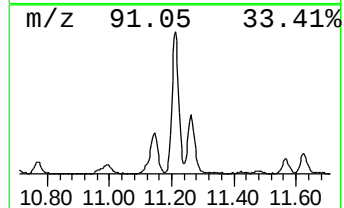
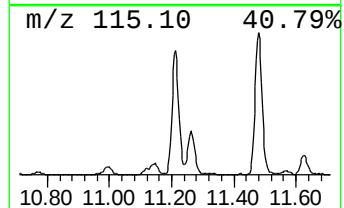
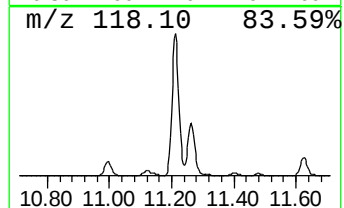
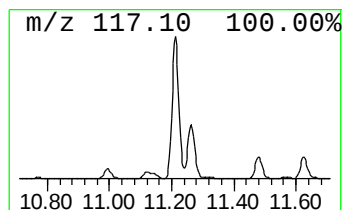
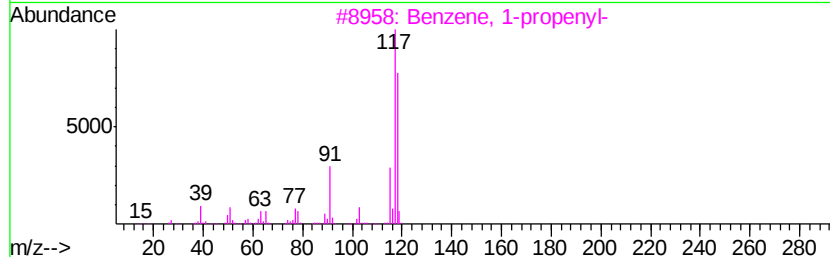
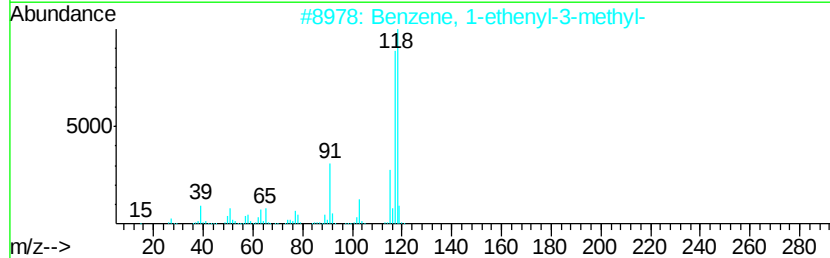
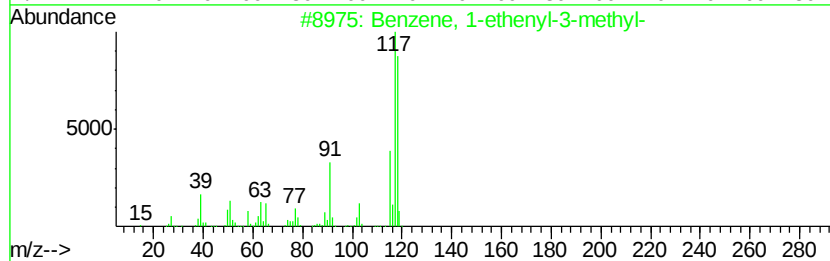
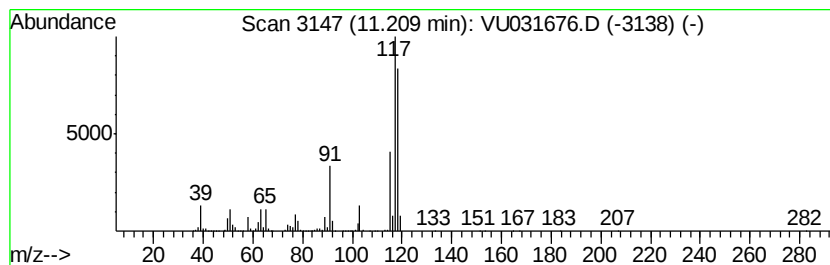
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Quant Title : VOC Analysis

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

Peak Number 6 Benzene, 1-ethenyl-3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.21	41.99 ug/L	2431370	1,4-Dichlorobenzene-d4	11.48

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-propenyl-	118	C9H10	000637-50-3	95
4		Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	94
5		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µL/5mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

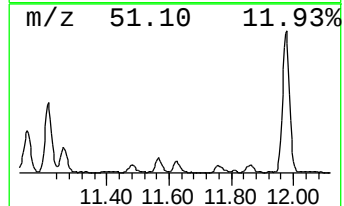
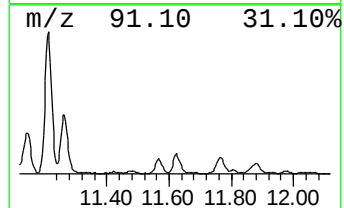
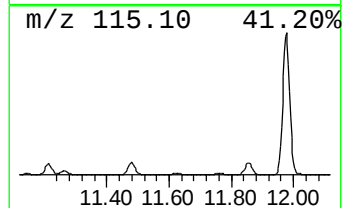
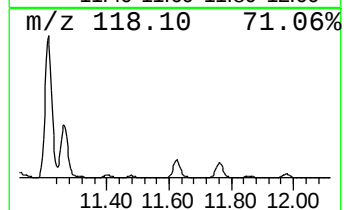
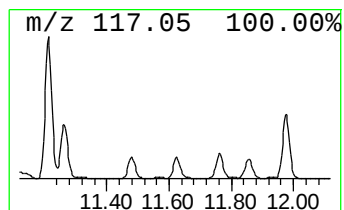
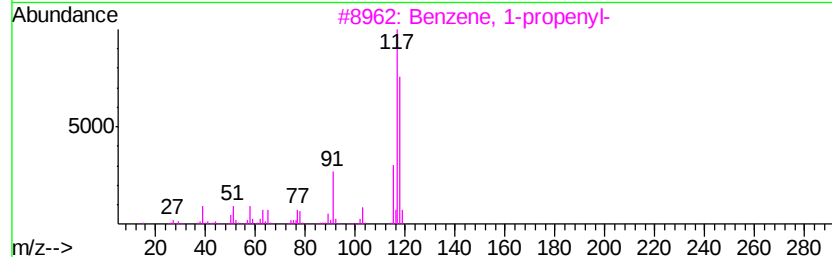
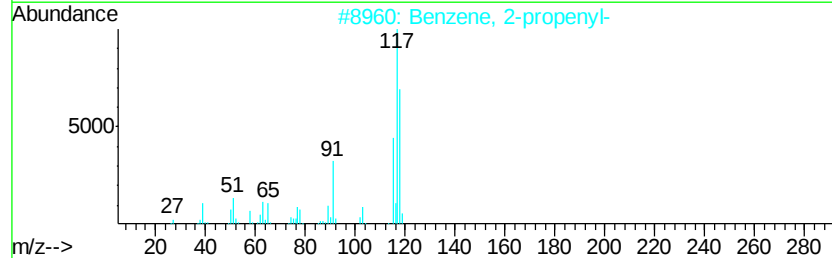
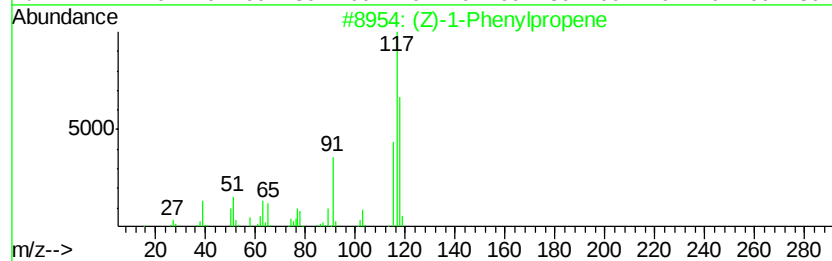
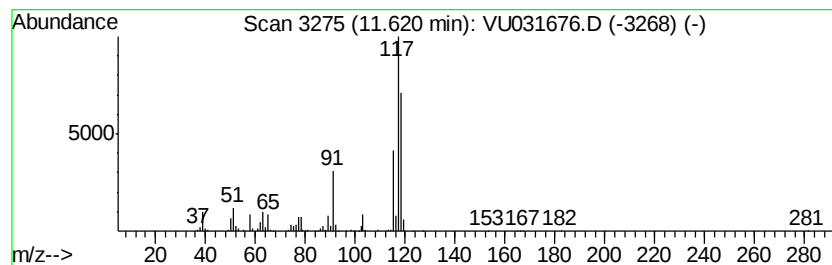
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MSVOA_U
ClientSampleID :
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Quant Title : VOC Analysis

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

Peak Number 9 (Z)-1-Phenylpropene Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.62	6.22 ug/L	360297	1,4-Dichlorobenzene-d4	11.48	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	95
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	93
3	Benzene, 1-propenyl-	118	C9H10	000637-50-3	93
4	Benzene, 2-propenyl-	118	C9H10	000300-57-2	93
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAJ80DL

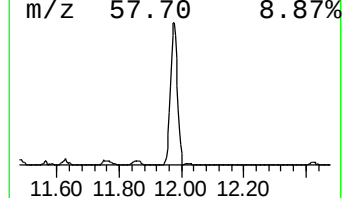
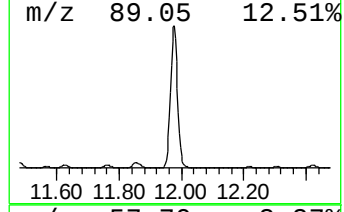
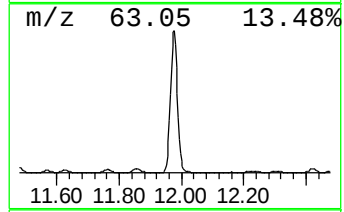
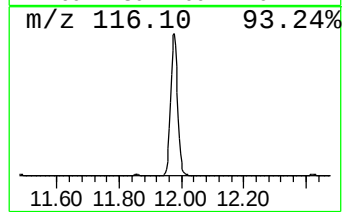
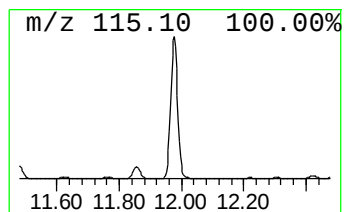
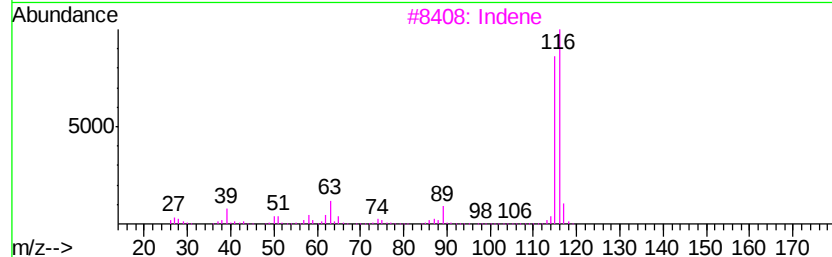
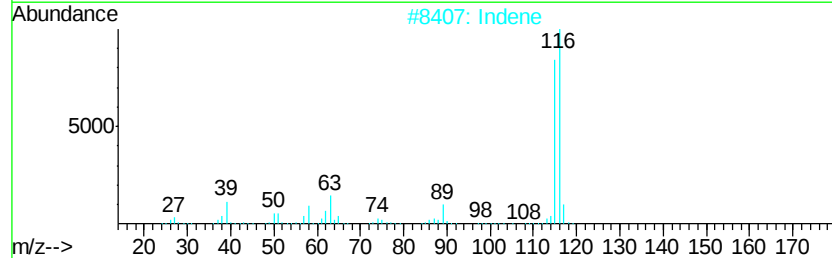
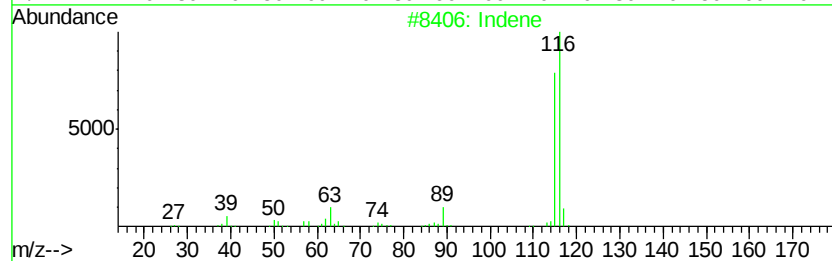
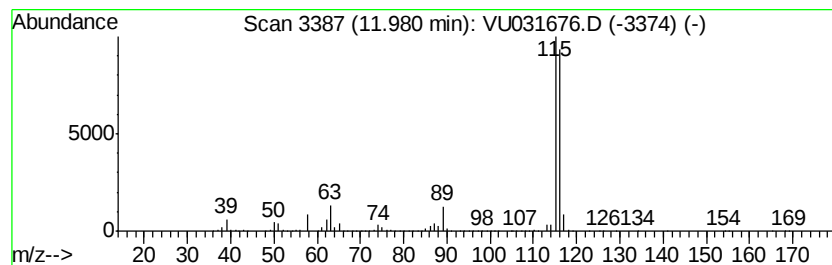
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Quant Title : VOC Analysis

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

Peak Number 11 Indene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	164.50 ug/L	9525800	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indene	116	C9H8	000095-13-6	97
2		Indene	116	C9H8	000095-13-6	96
3		Indene	116	C9H8	000095-13-6	95
4		Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5		3-Methylphenylacetylene	116	C9H8	000766-82-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01g/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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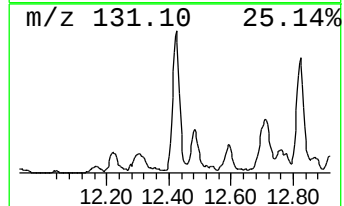
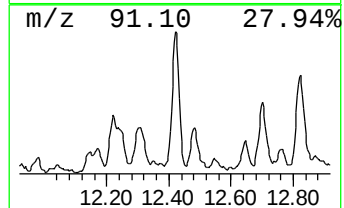
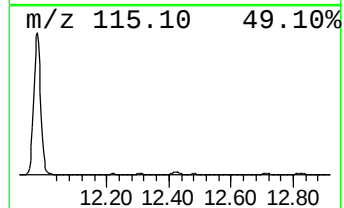
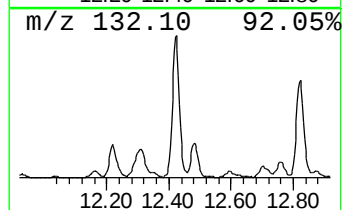
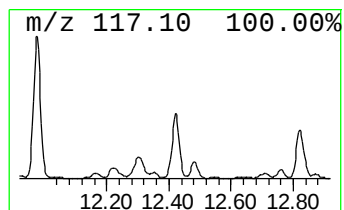
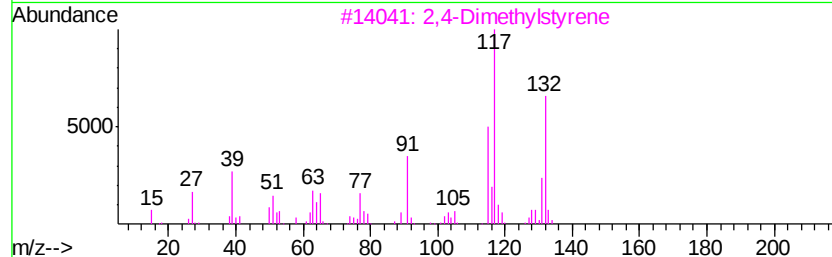
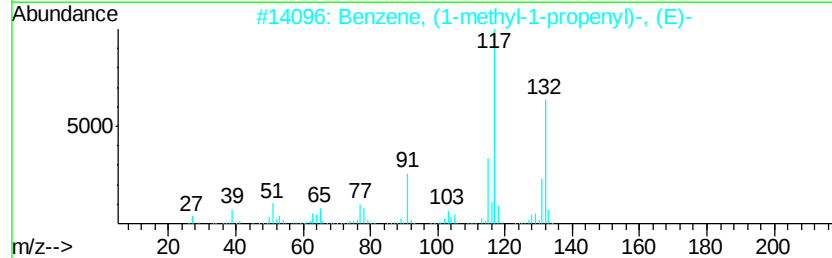
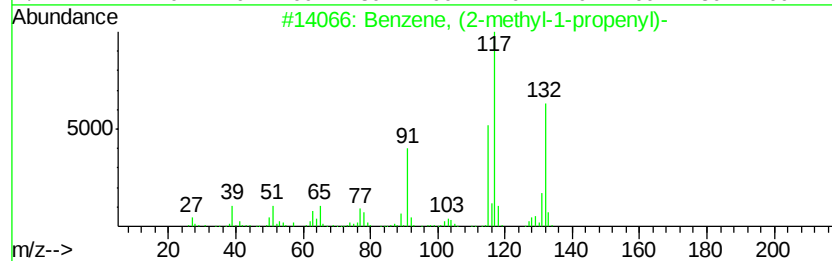
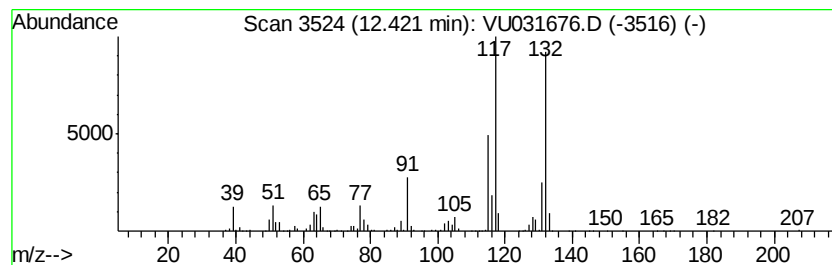
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Quant Title : VOC Analysis

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

Peak Number 12 Benzene, (2-methyl-1-propen... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.42	10.28 ug/L	595504	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	95
2		Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	94
3		2,4-Dimethylstyrene	132	C10H12	002234-20-0	94
4		Benzene, 1-methyl-4-(1-methyleth...	132	C10H12	001195-32-0	94
5		o-Isopropenyltoluene	132	C10H12	007399-49-7	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

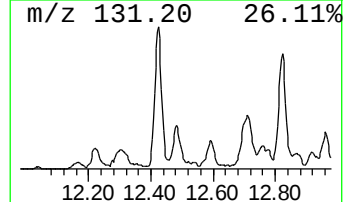
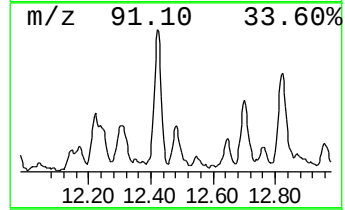
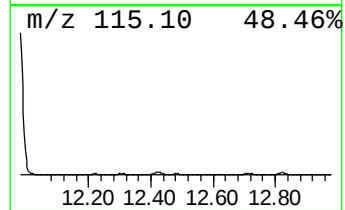
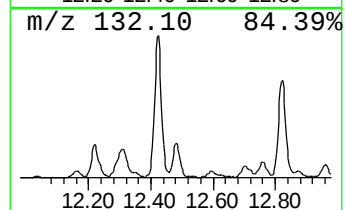
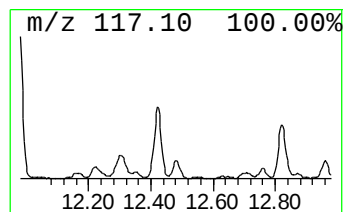
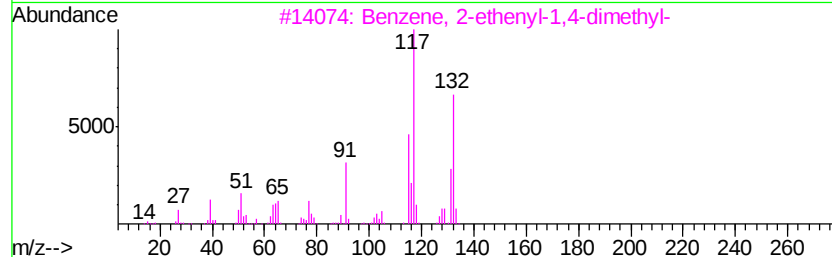
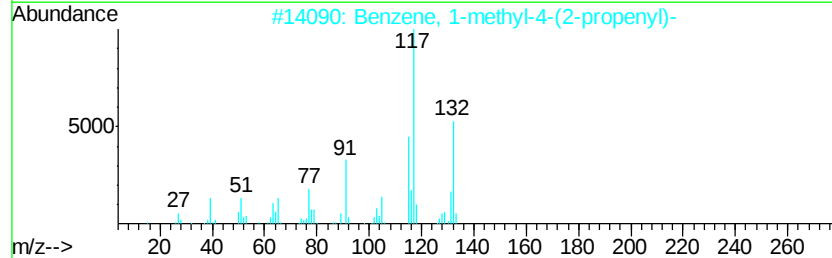
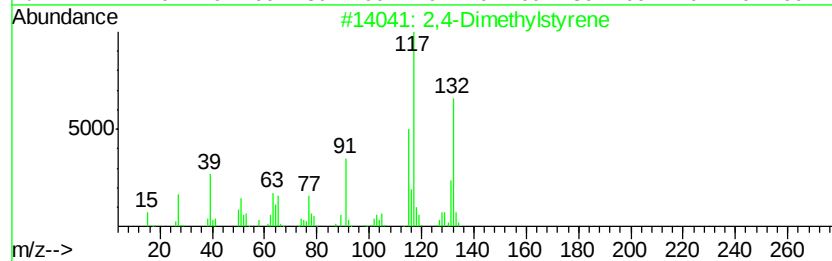
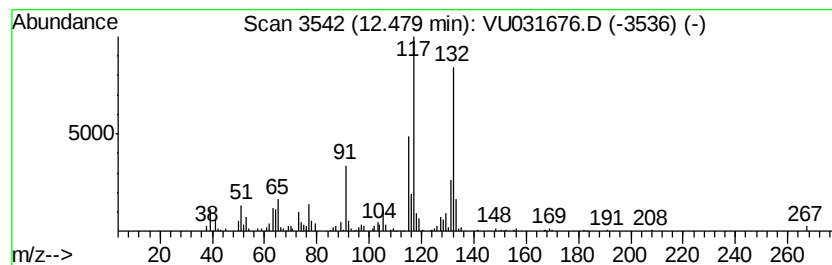
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ClientSampleId :
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Quant Title : VOC Analysis

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

Peak Number 13 2,4-Dimethylstyrene Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.48	2.99 ug/L	173060	1,4-Dichlorobenzene-d4	11.48
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2,4-Dimethylstyrene	132 C10H12	002234-20-0	97
2	Benzene, 1-methyl-4-(2-propenyl)-	132 C10H12	003333-13-9	97
3	Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6	96
4	Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6	96
5	Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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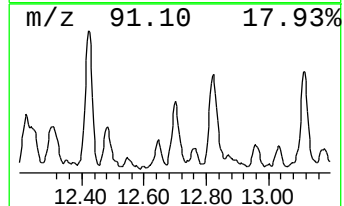
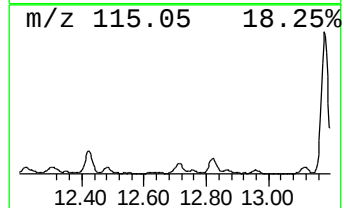
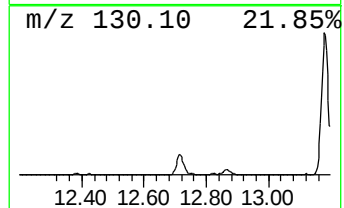
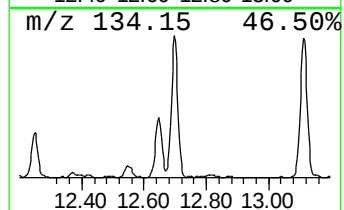
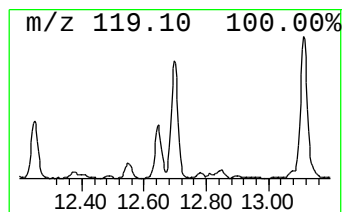
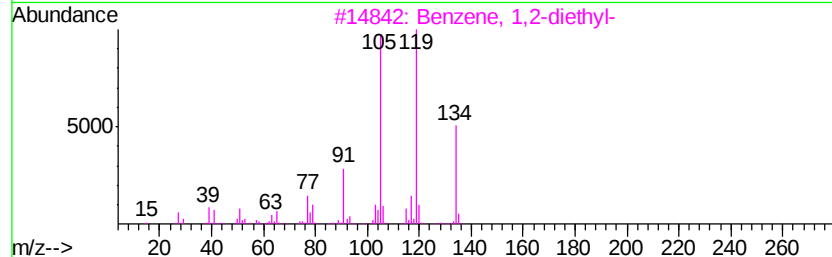
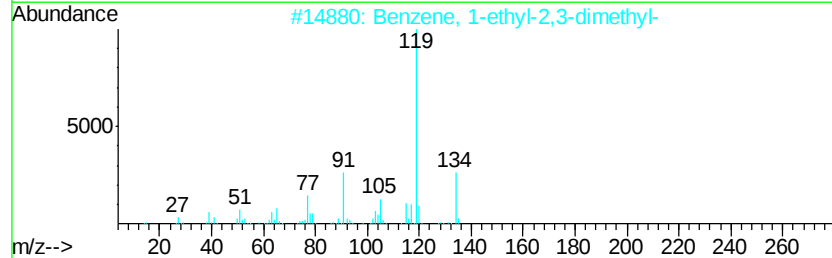
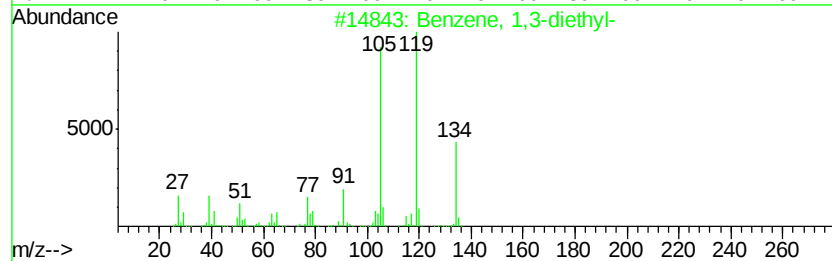
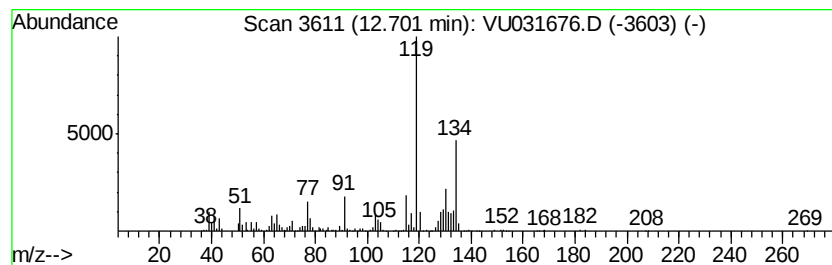
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Quant Title : VOC Analysis

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

Peak Number 14 Benzene, 1,3-diethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	10.75 ug/L	622351	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	55
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	55
3		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	55
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	55
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01g/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAJ80DL

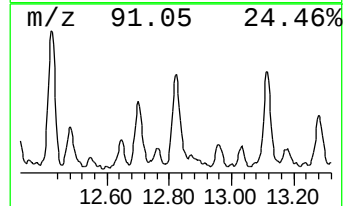
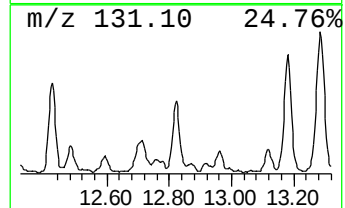
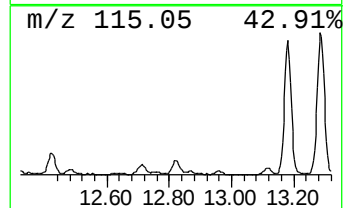
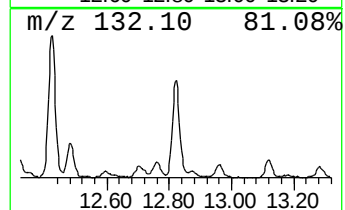
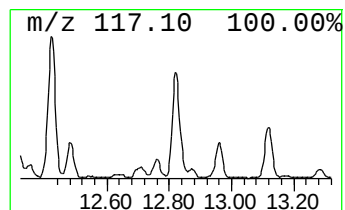
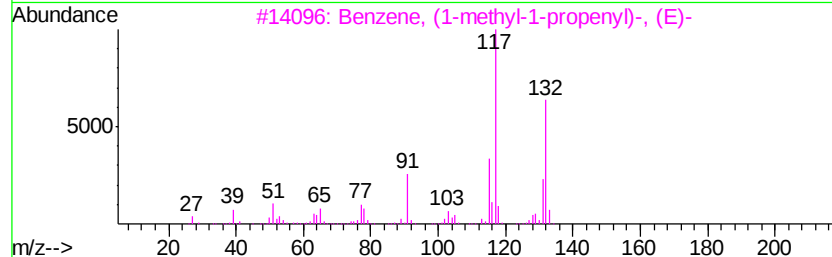
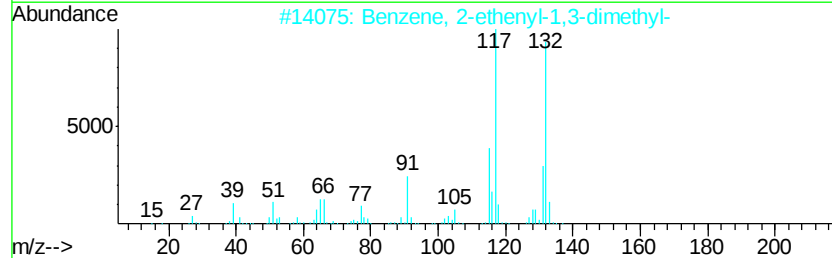
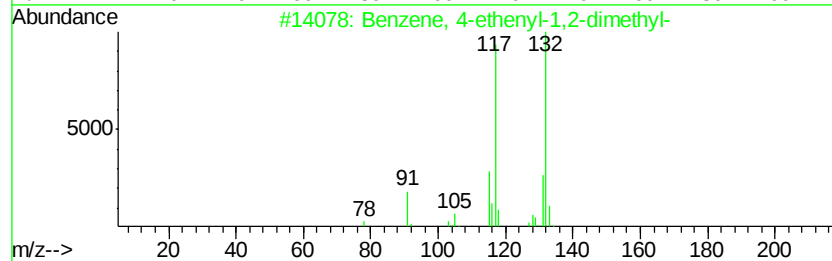
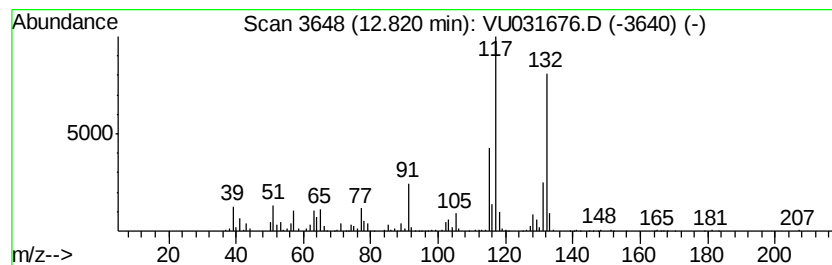
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Quant Title : VOC Analysis

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

Peak Number 15 Benzene, 4-ethenyl-1,2-dime... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.82	8.24 ug/L	476982	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethenyl-1,2-dimethyl-	132	C10H12	027831-13-6	95
2		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	95
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	94
4		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	94
5		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAJ80DL

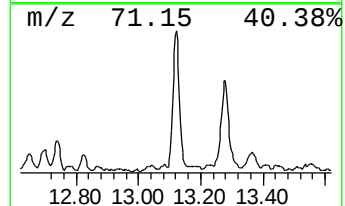
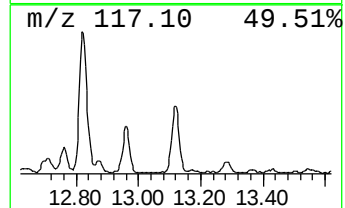
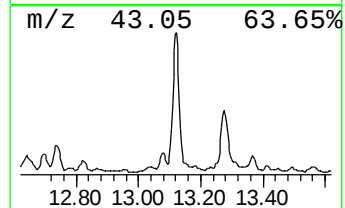
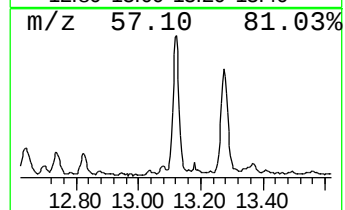
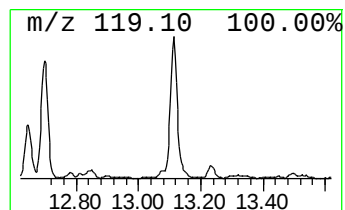
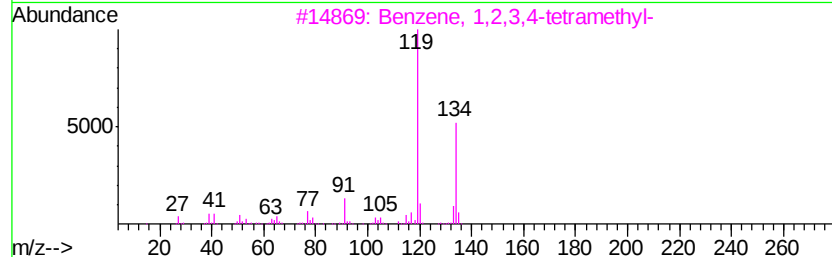
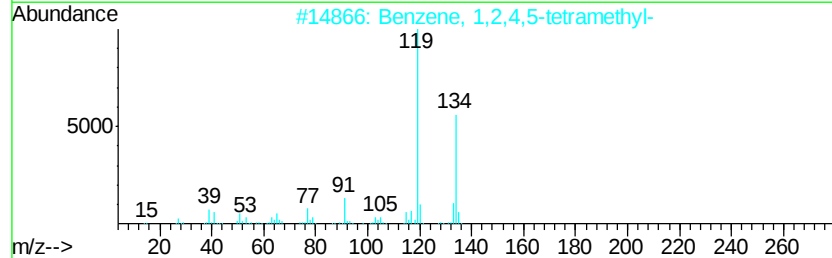
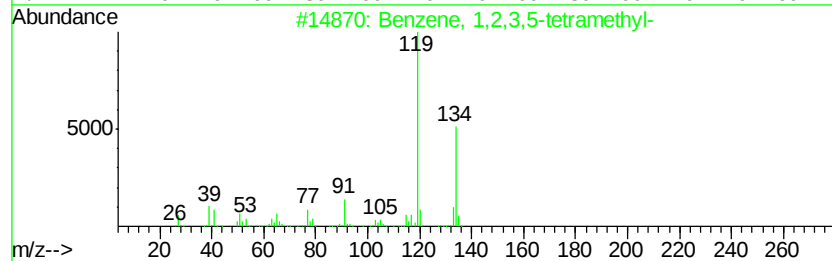
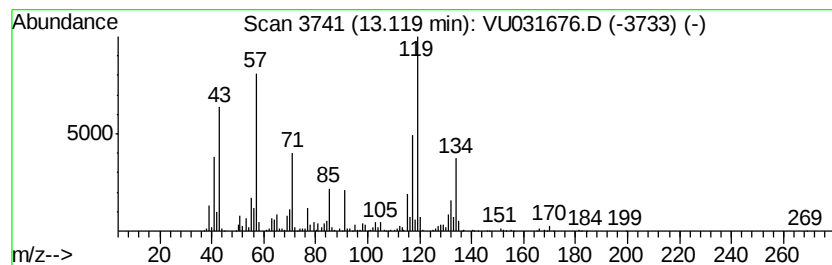
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Quant Title : VOC Analysis

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

Peak Number 16 unknown-01 Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	10.70 ug/L	619633	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	46
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	41
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	41
4		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	41
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	41



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAJ80DL

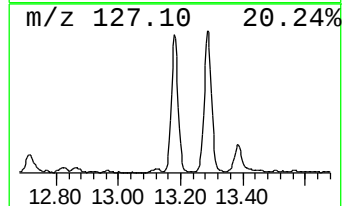
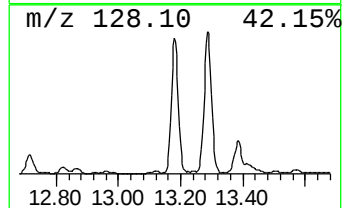
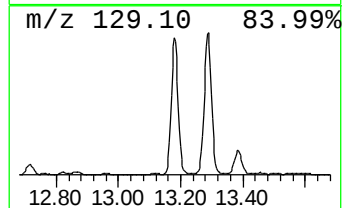
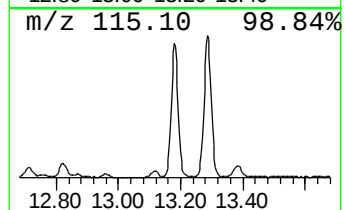
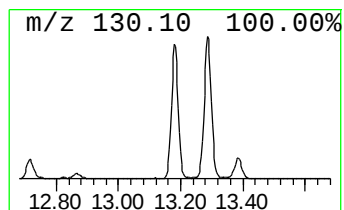
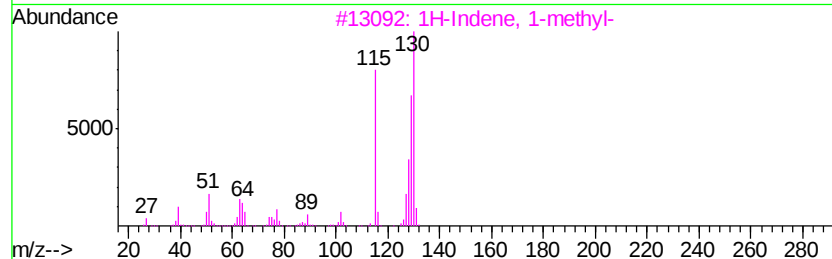
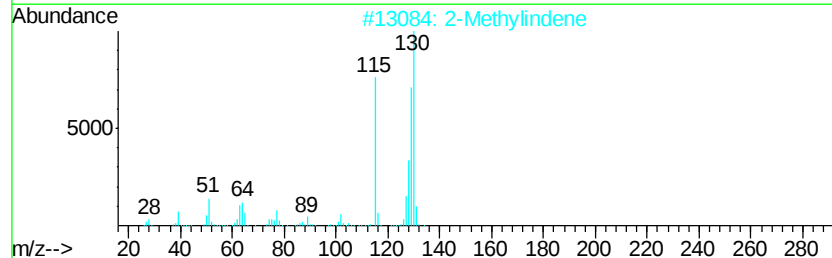
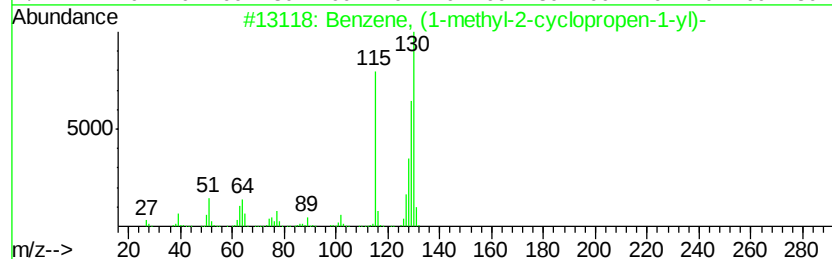
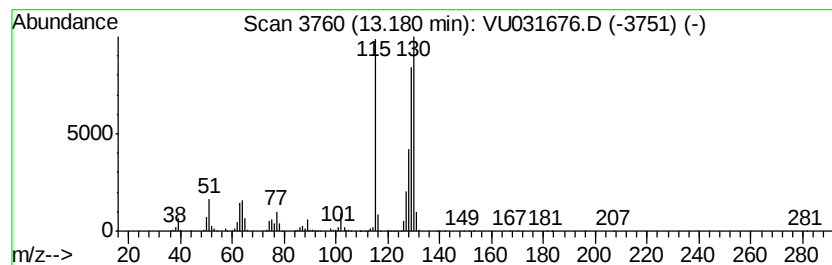
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Quant Title : VOC Analysis

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

Peak Number 17 Benzene, (1-methyl-2-cyclop... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.18	34.14 ug/L	1976940	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	97
2		2-Methylindene	130	C10H10	002177-47-1	97
3		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
4		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	94
5		1H-Indene, 3-methyl-	130	C10H10	000767-60-2	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01g/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAJ80DL

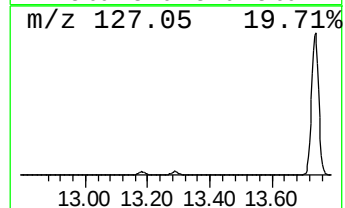
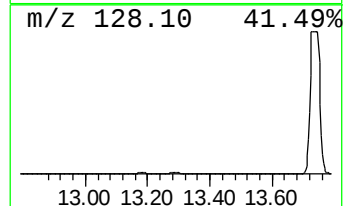
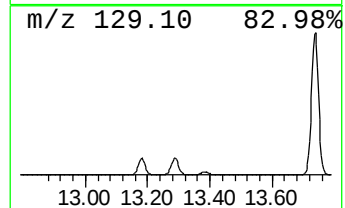
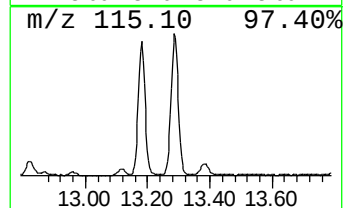
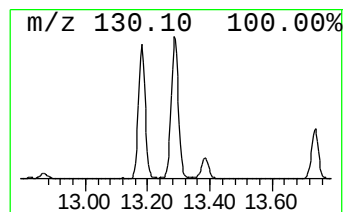
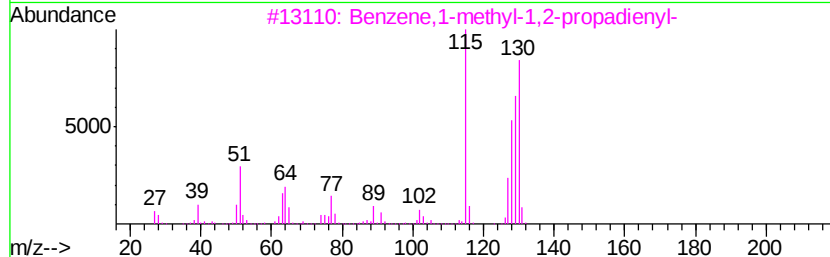
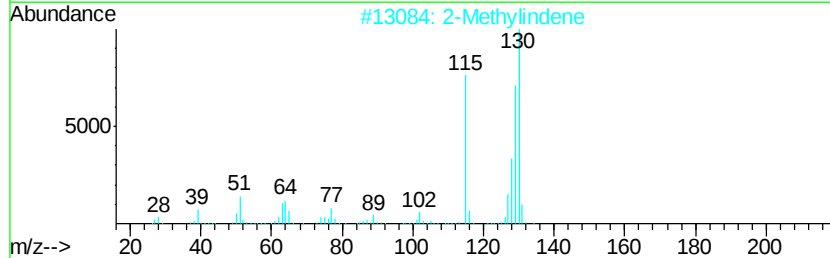
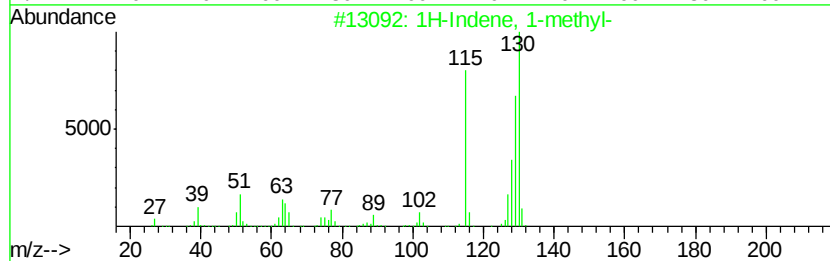
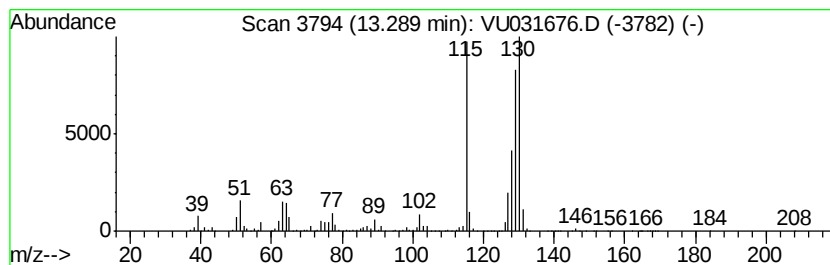
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Quant Title : VOC Analysis

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

Peak Number 18 1H-Indene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.29	42.21 ug/L	2444370	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
2		2-Methylindene	130	C10H10	002177-47-1	96
3		Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	96
4		Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	96
5		Benzene, 1-butyryl-	130	C10H10	000622-76-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

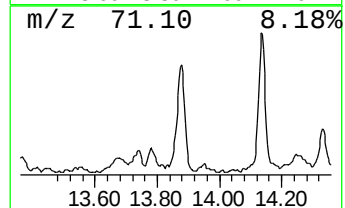
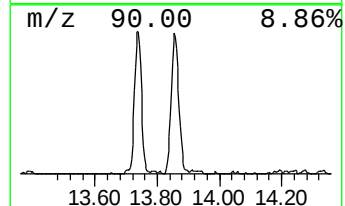
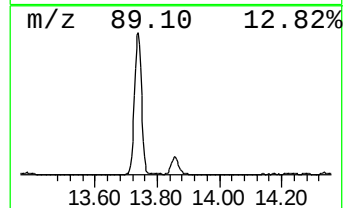
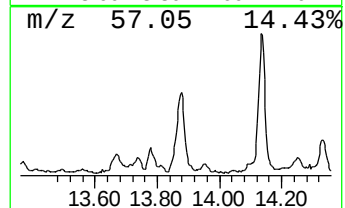
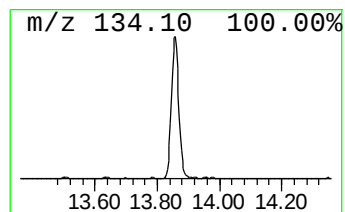
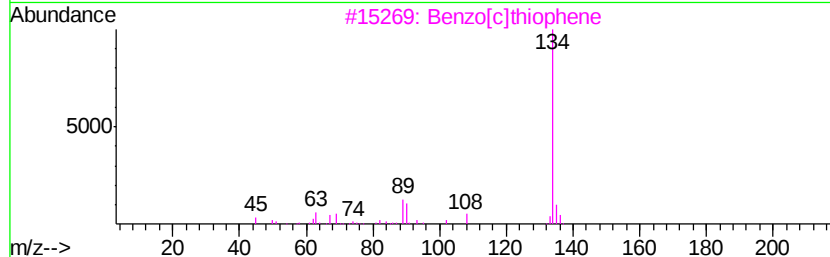
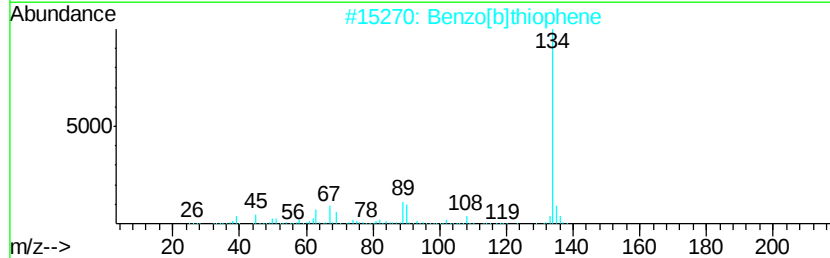
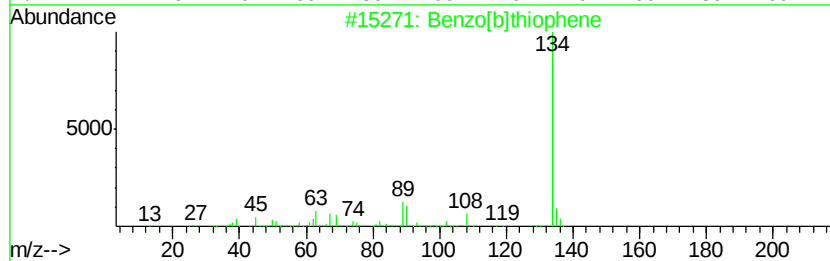
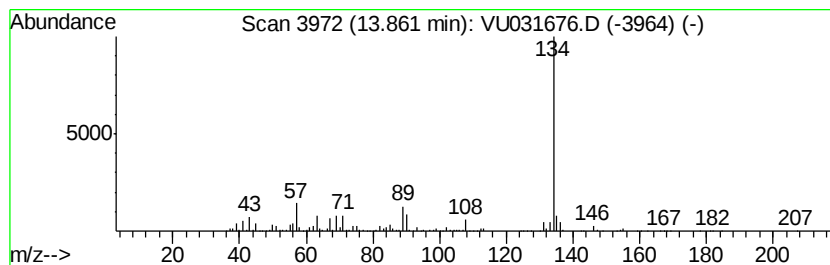
Instrument :
MSVOA_U
ClientSampleId :
GAJ80DL

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM042719WMA.M
Quant Title : VOC Analysis

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

Peak Number 20 Benzo[b]thiophene Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	12.09 ug/L	700151	1,4-Dichlorobenzene-d4	11.48
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzo[b]thiophene	134 C8H6S	000095-15-8 95
2		Benzo[b]thiophene	134 C8H6S	000095-15-8 94
3		Benzo[c]thiophene	134 C8H6S	000270-82-6 94
4		Benzo[b]thiophene	134 C8H6S	000095-15-8 94
5		Cyclopenta[c]thiapyran	134 C8H6S	000270-63-3 91



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAJ80DL

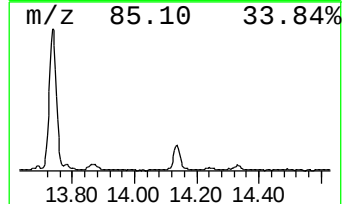
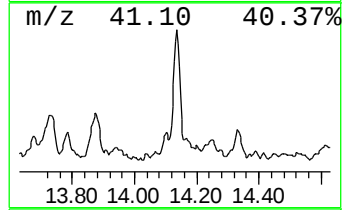
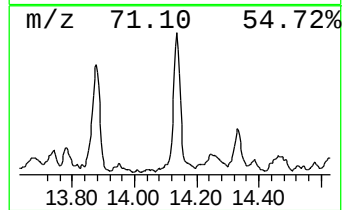
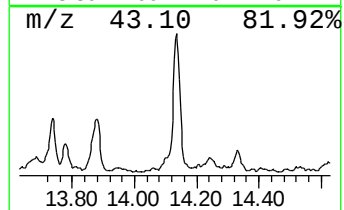
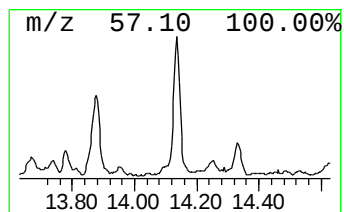
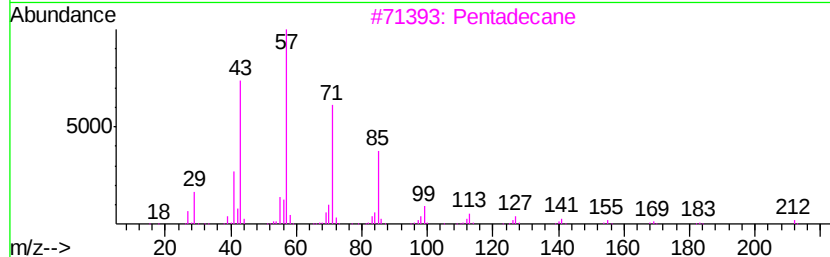
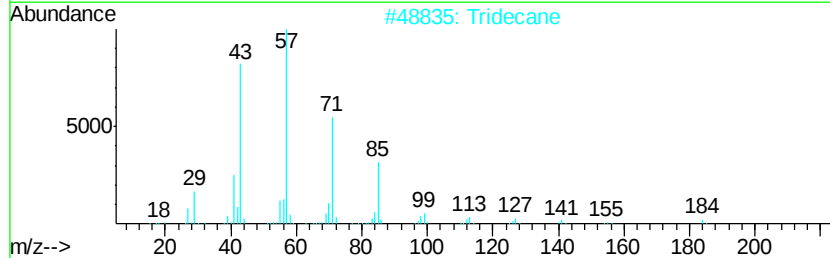
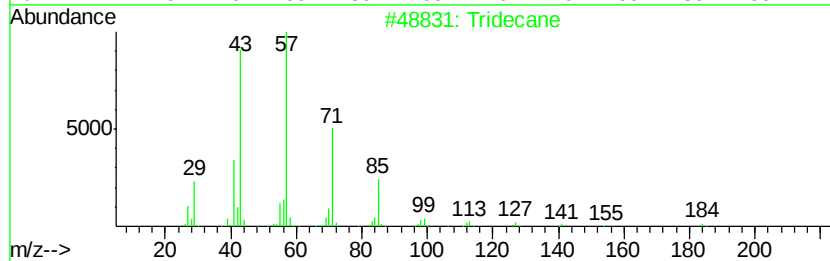
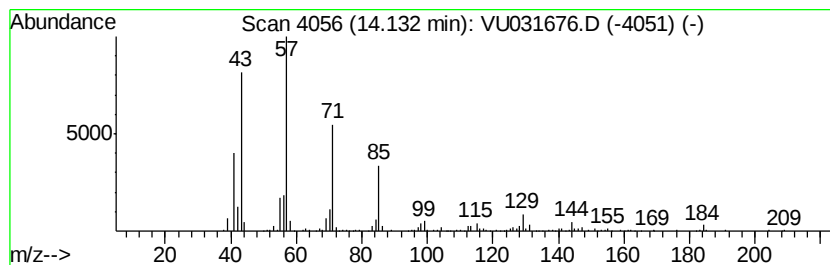
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Quant Title : VOC Analysis

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.13	4.73 ug/L	273789	1,4-Dichlorobenzene-d4	11.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	91
2			Tridecane	184	C13H28	000629-50-5	87
3			Pentadecane	212	C15H32	000629-62-9	80
4			Tetradecane	198	C14H30	000629-59-4	80
5			Tetradecane	198	C14H30	000629-59-4	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µL/5mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

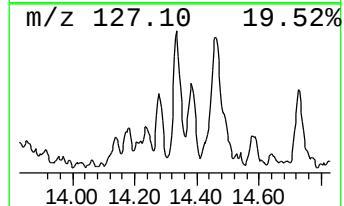
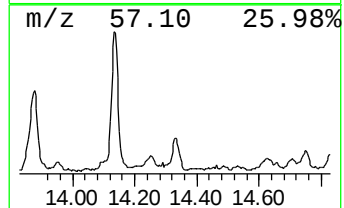
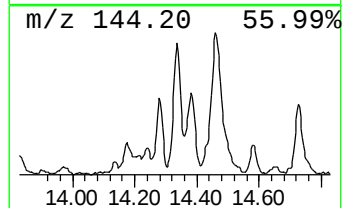
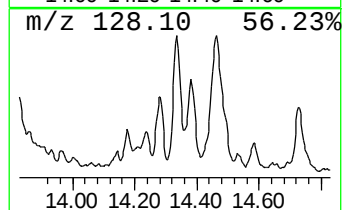
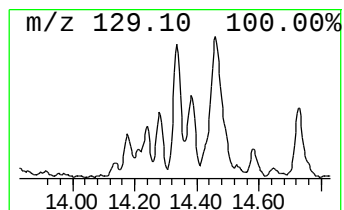
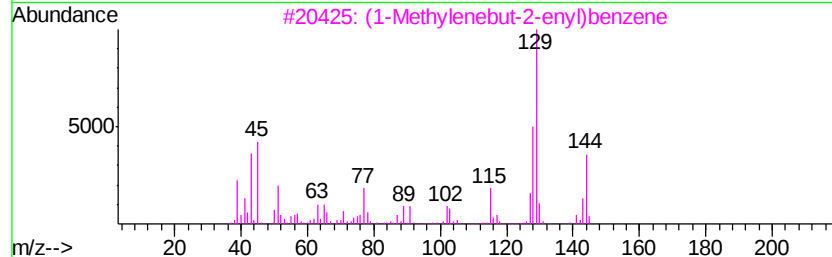
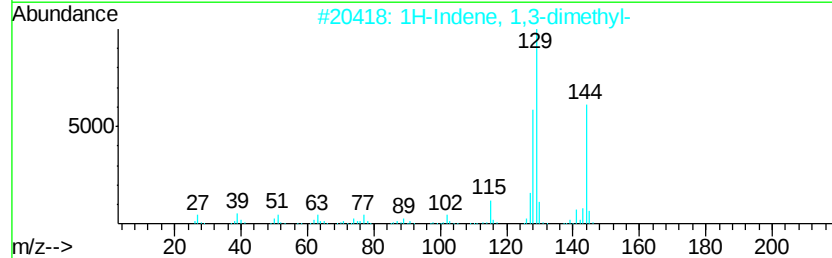
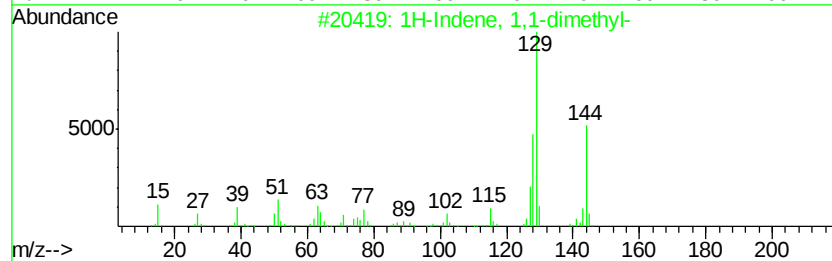
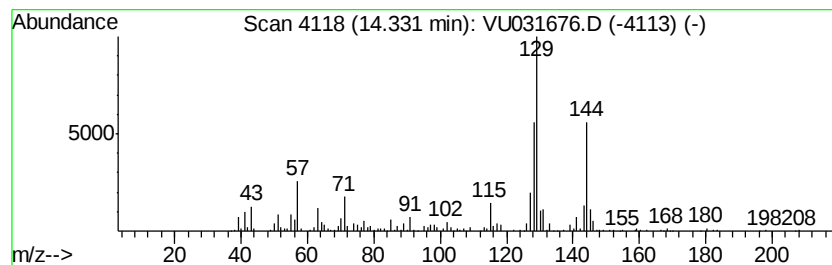
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ClientSampleId :
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM042719WMA.M
Quant Title : VOC Analysis

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

Peak Number 22 1H-Indene, 1,1-dimethyl- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.33	4.50 ug/L	260462	1,4-Dichlorobenzene-d4	11.48	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 1,1-dimethyl-	144	C11H12	018636-55-0	95
2	1H-Indene, 1,3-dimethyl-	144	C11H12	002177-48-2	91
3	(1-Methylenebut-2-enyl)benzene	144	C11H12	070588-46-4	90
4	(1-Methylbuta-1,3-dienvl)benzene	144	C11H12	054758-36-0	87
5	Benzene, (1-methyl-2-butynyl)-	144	C11H12	087712-69-4	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µ/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

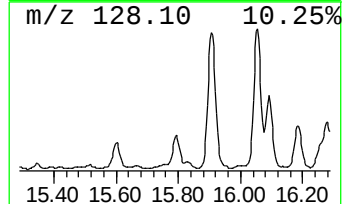
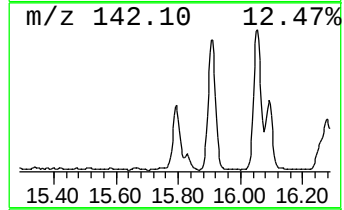
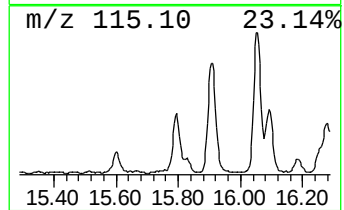
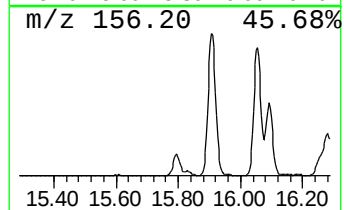
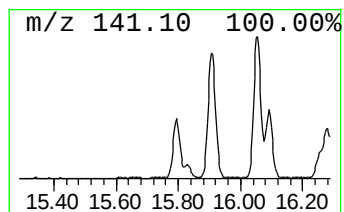
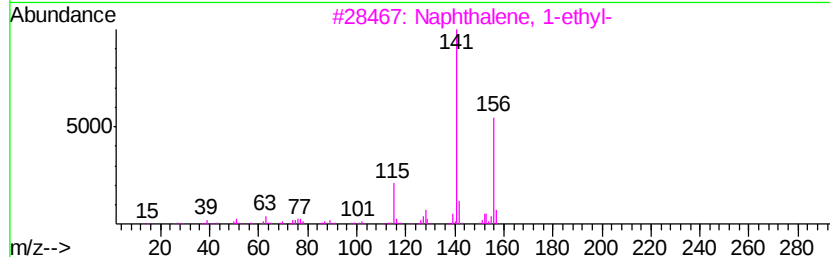
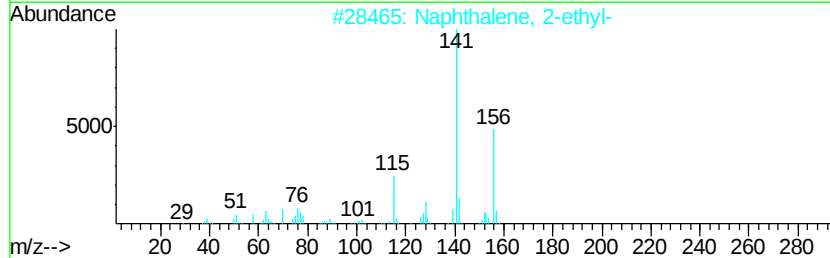
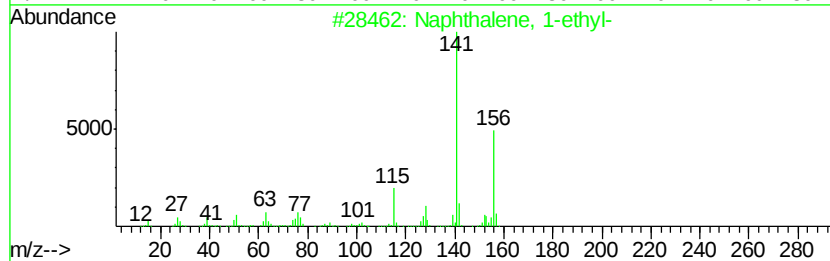
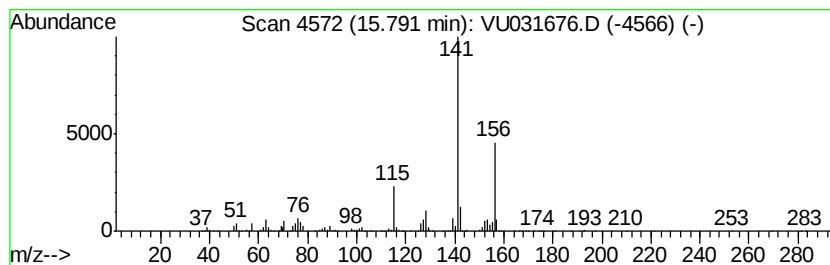
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MSVOA_U
ClientSampleId :
GAJ80DL

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Quant Title : VOC Analysis

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

Peak Number 26 Naphthalene, 1-ethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.79	13.30 ug/L	770293	1,4-Dichlorobenzene-d4	11.48	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	96
2	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	96
3	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	95
4	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	95
5	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
 Data File : VU031676.D
 Acq On : 01 May 2019 12:00
 Operator : JC/SP
 Sample : K2604-06DL 25X
 Misc : 5.01µ/5mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 GAJ80DL

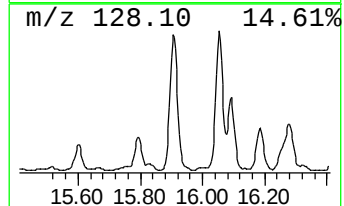
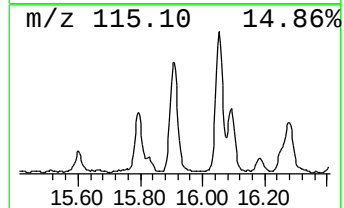
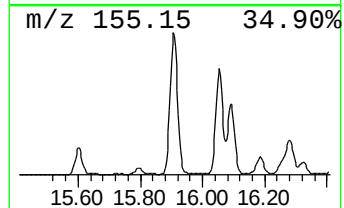
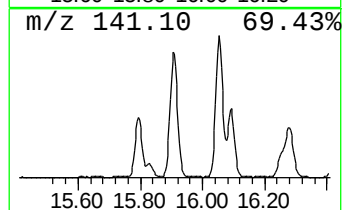
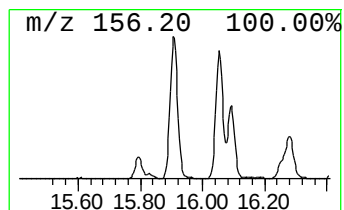
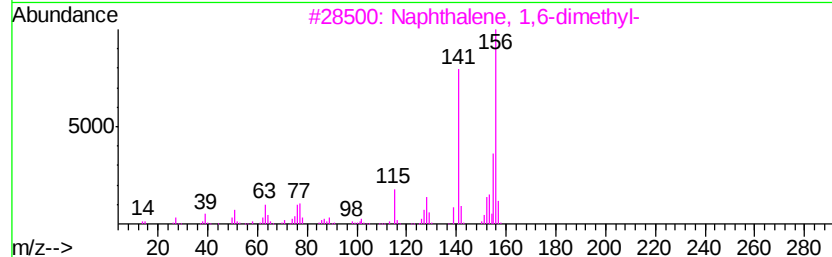
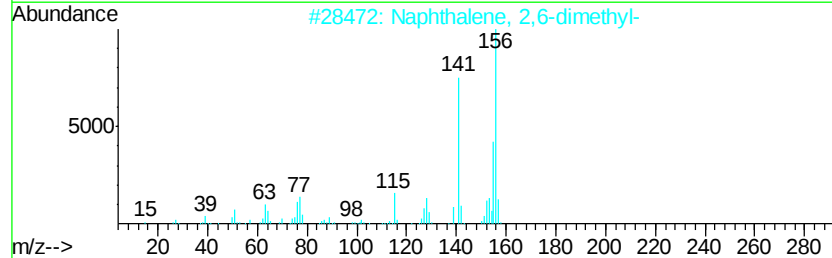
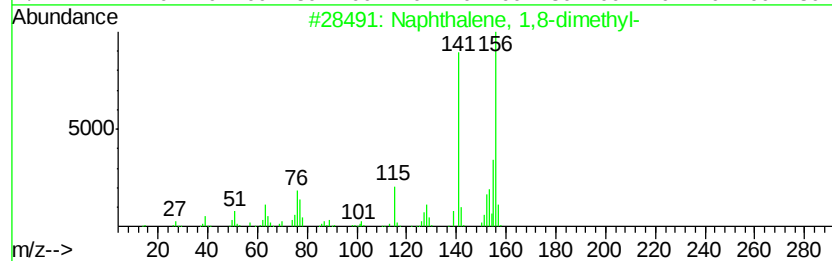
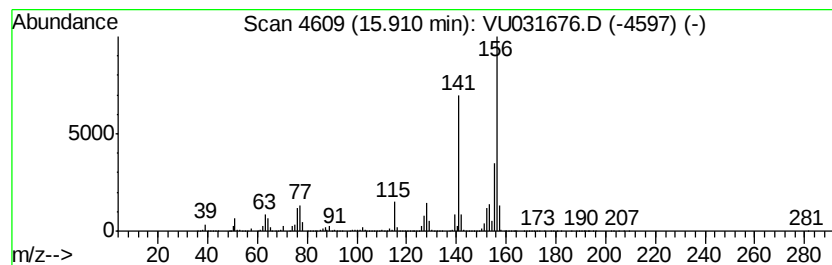
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 Quant Title : VOC Analysis

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

 Peak Number 27 Naphthalene, 1,8-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	73.86 ug/L	4277210	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	98
2		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
3		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
4		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
5		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µ/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAJ80DL

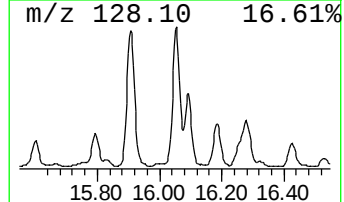
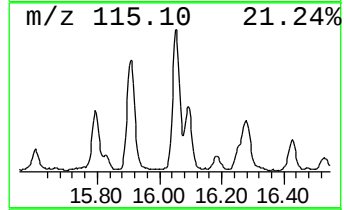
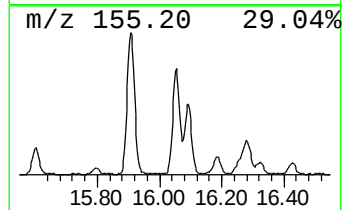
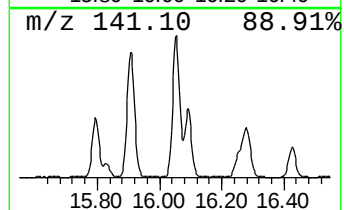
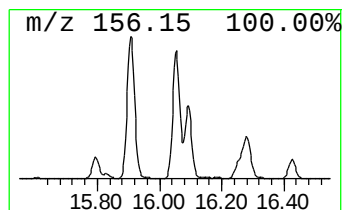
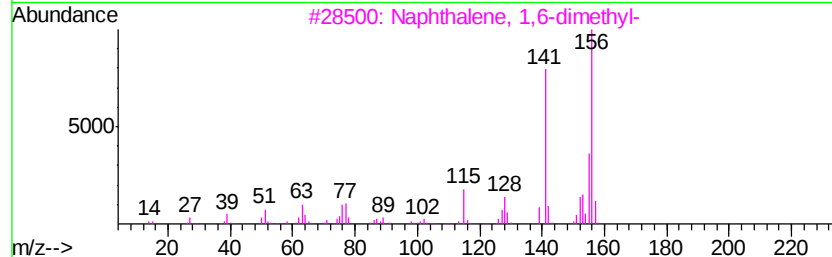
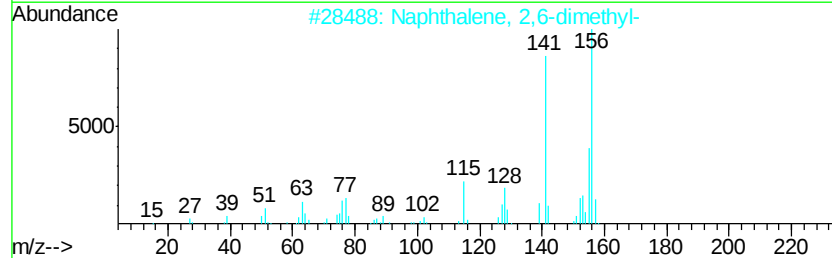
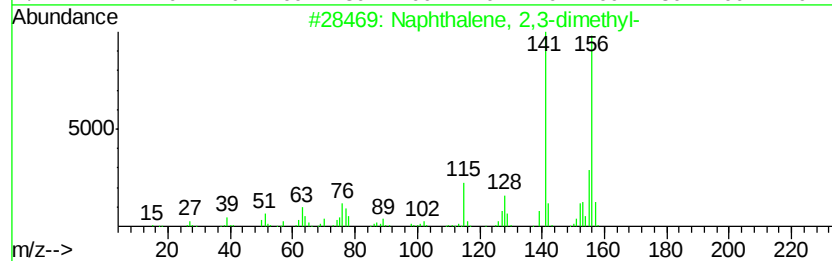
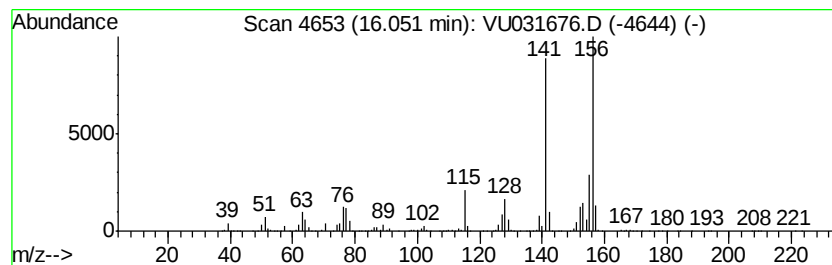
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Quant Title : VOC Analysis

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

Peak Number 28 Naphthalene, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.05	63.55 ug/L	3679840	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
3		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
4		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
5		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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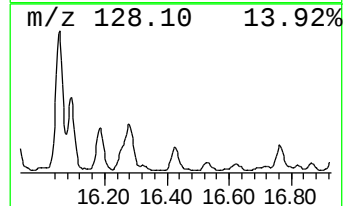
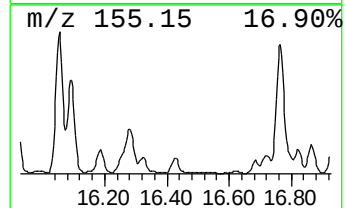
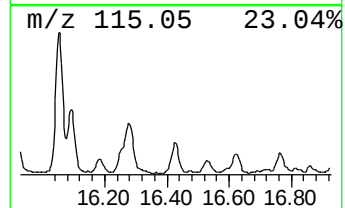
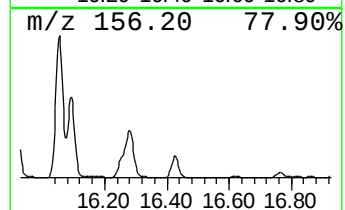
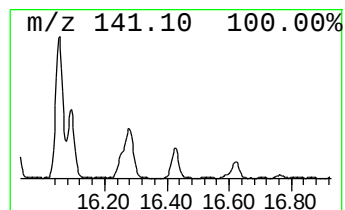
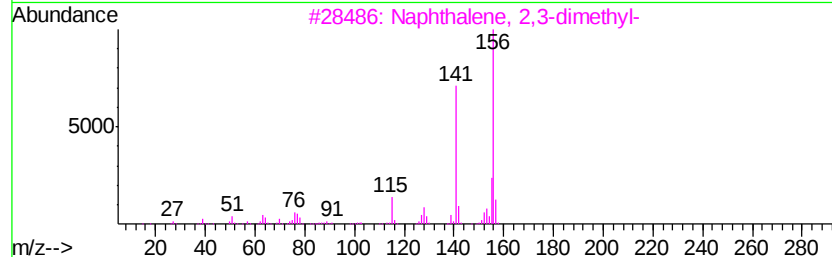
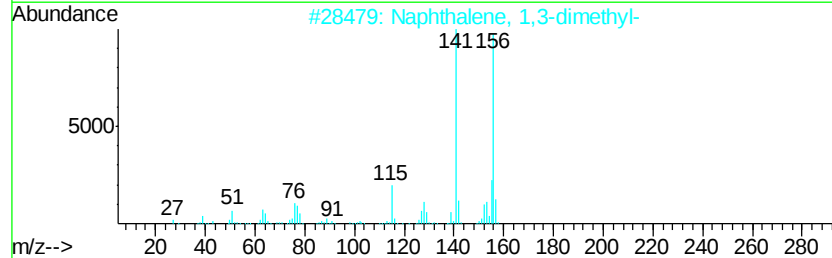
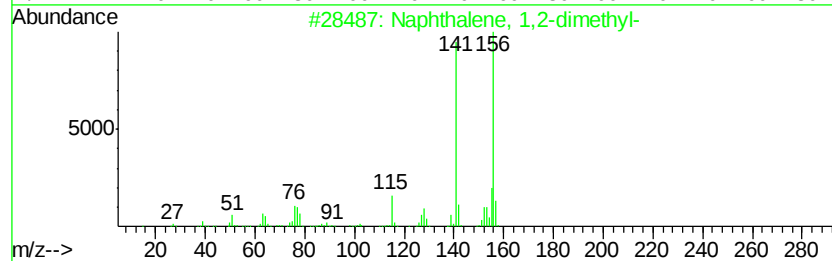
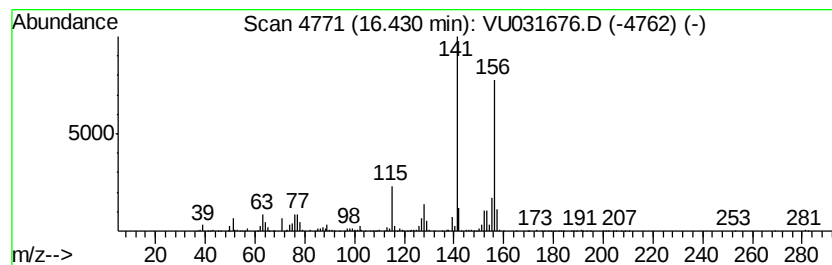
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Quant Title : VOC Analysis

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

Peak Number 31 Naphthalene, 1,2-dimethyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.43	10.18 ug/L	589442	1,4-Dichlorobenzene-d4	11.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96
2			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
5			Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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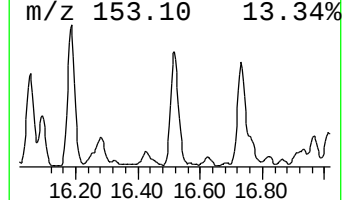
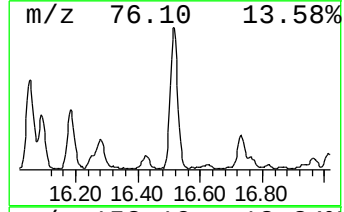
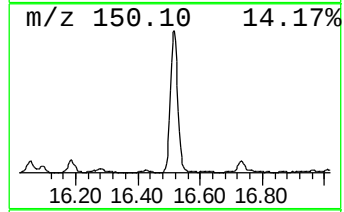
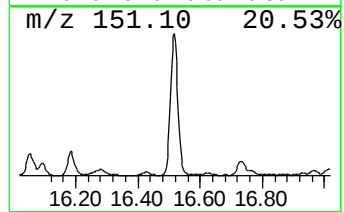
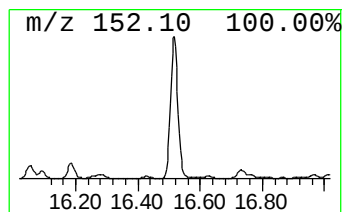
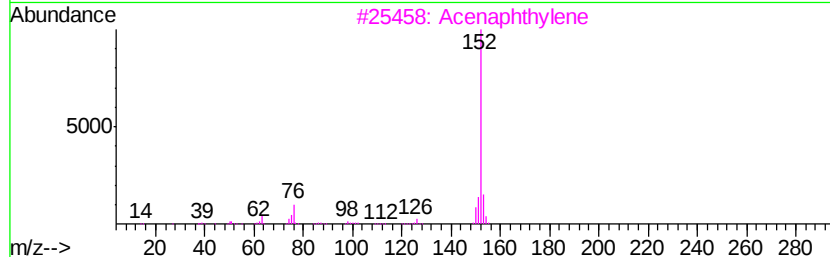
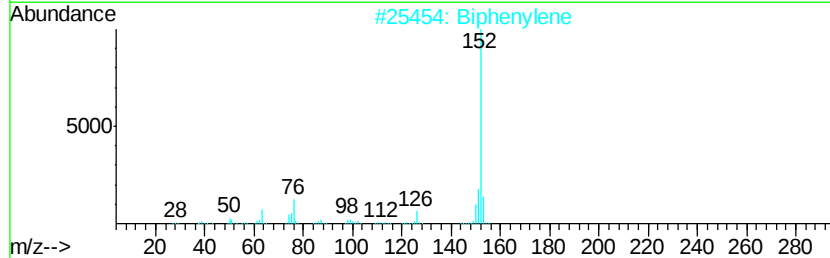
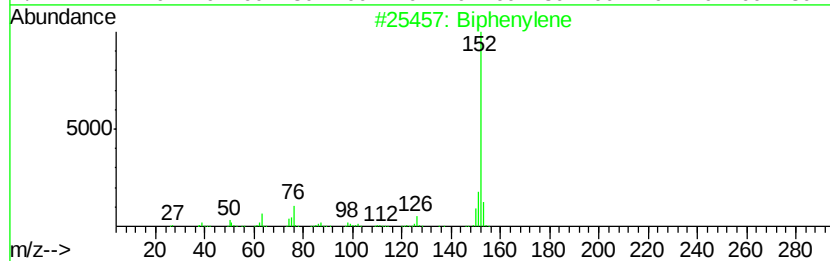
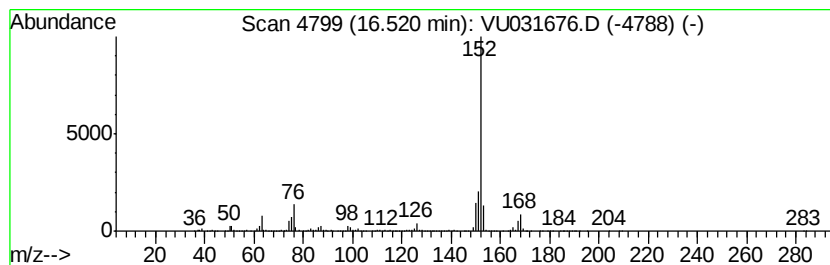
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Quant Title : VOC Analysis

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

Peak Number 32 Biphenylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.52	55.31 ug/L	3202910	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Biphenylene	152	C12H8	000259-79-0	95
2		Biphenylene	152	C12H8	000259-79-0	93
3		Acenaphthylene	152	C12H8	000208-96-8	87
4		Acenaphthylene	152	C12H8	000208-96-8	87
5		Biphenylene	152	C12H8	000259-79-0	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µ/5mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GAJ80DL

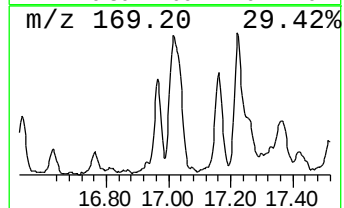
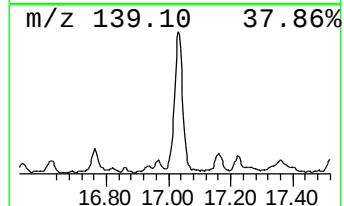
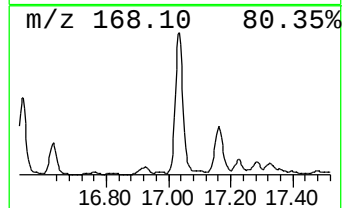
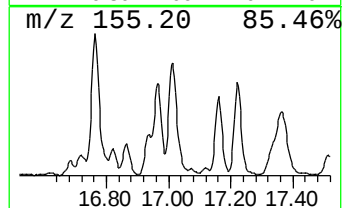
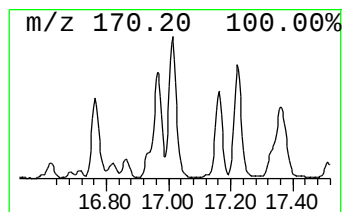
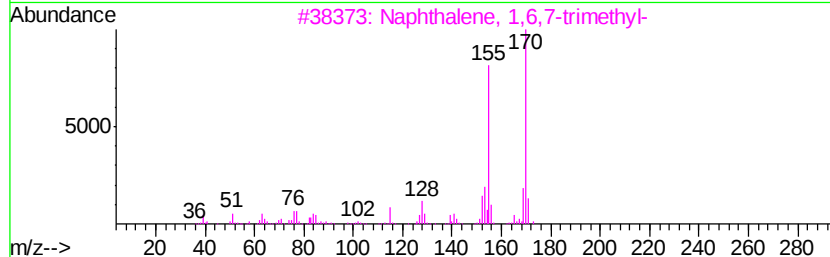
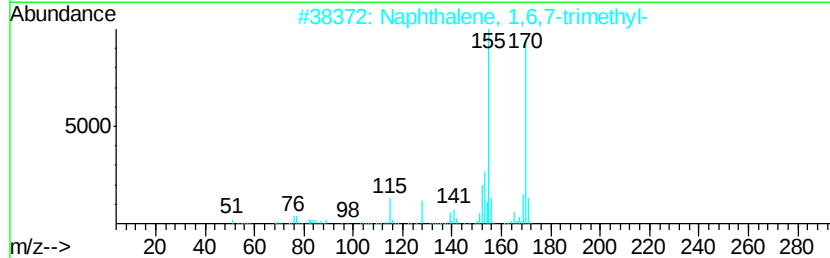
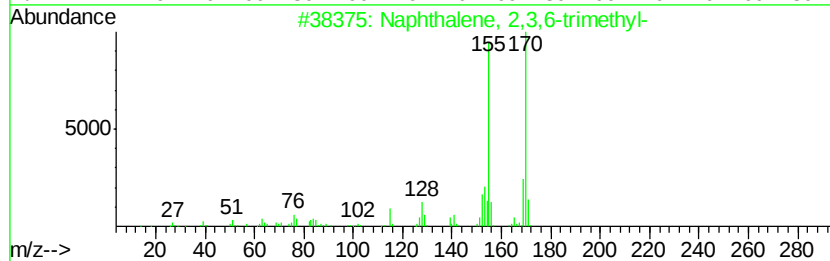
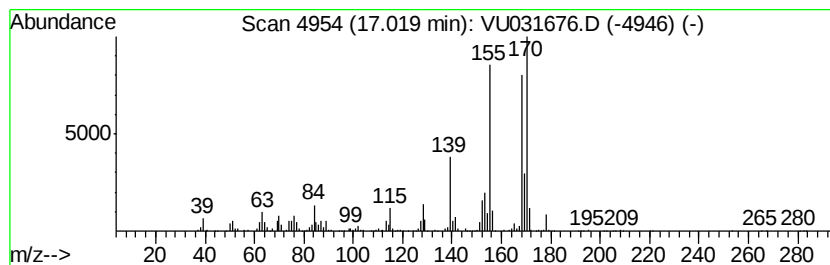
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Quant Title : VOC Analysis

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

Peak Number 33 Naphthalene, 2,3,6-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.02	29.46 ug/L	1705970	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
2		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
3		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
4		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	92
5		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU050119\
Data File : VU031676.D
Acq On : 01 May 2019 12:00
Operator : JC/SP
Sample : K2604-06DL 25X
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA_U/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
GAJ80DL

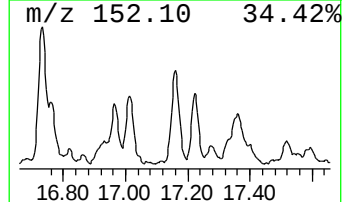
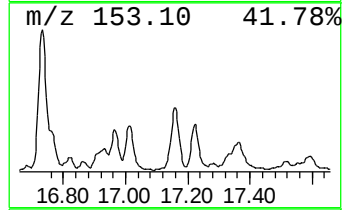
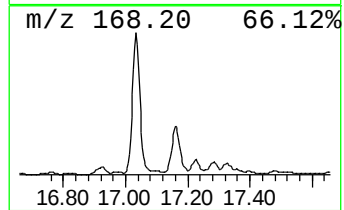
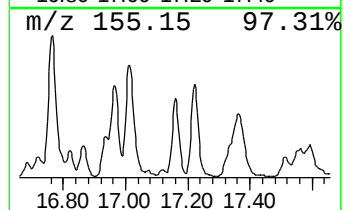
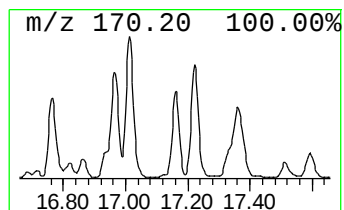
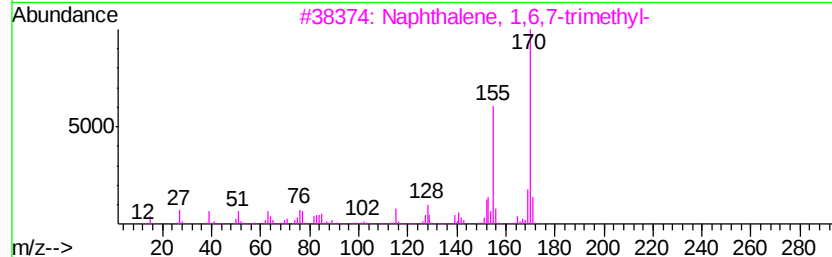
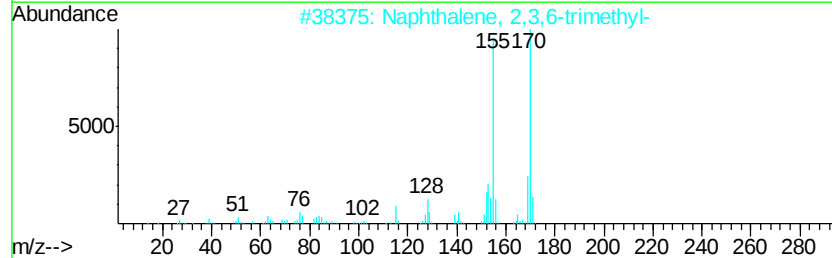
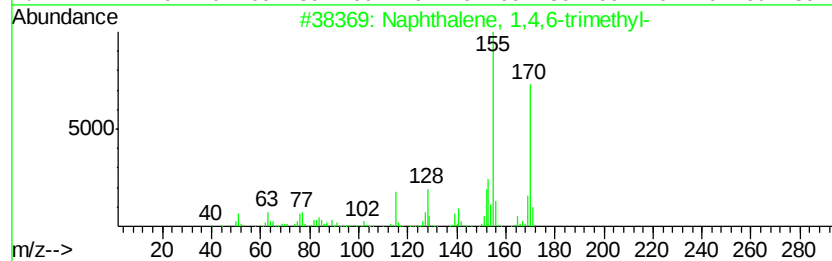
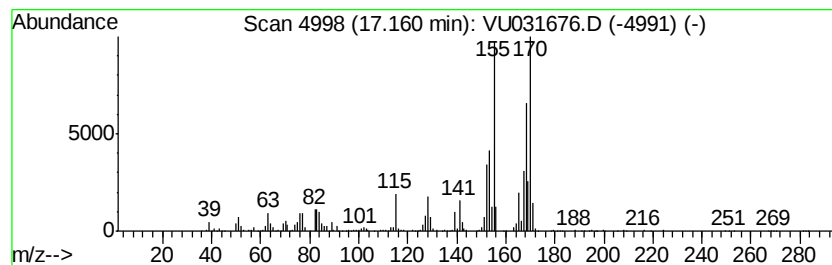
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM042719WMA.M
Quant Title : VOC Analysis

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

Peak Number 34 Naphthalene, 1,4,6-trimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.16	14.55 ug/L	842378	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	93
2		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	91
3		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	90
4		3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	83
5		1-(2,2-Dimethylcyclopropyl)-2-ph...	170	C13H14	1000294-91-1	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050119\
 Data File : VU031676.D
 Acq On : 01 May 2019 12:00
 Operator : JC/SP
 Sample : K2604-06DL 25X
 Misc : 5.01g/5mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 GAJ80DL

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM042719WMA.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
Benzene, 1-ethyl-...	10.68	8.5	ug/L	490941	3	11.48	2895430	50.0
Benzene, 1,2,3-tr...	10.77	7.8	ug/L	452410	3	11.48	2895430	50.0
.alpha.-Methylsty...	10.99	6.9	ug/L	401509	3	11.48	2895430	50.0
Benzene, 1-etheny...	11.21	42.0	ug/L	2431370	3	11.48	2895430	50.0
(Z)-1-Phenylpropene	11.62	6.2	ug/L	360297	3	11.48	2895430	50.0
Indene	11.98	164.5	ug/L	9525800	3	11.48	2895430	50.0
Benzene, (2-methy...	12.42	10.3	ug/L	595504	3	11.48	2895430	50.0
2,4-Dimethylstyrene	12.48	3.0	ug/L	173060	3	11.48	2895430	50.0
Benzene, 1,3-diet...	12.70	10.8	ug/L	622351	3	11.48	2895430	50.0
Benzene, 4-etheny...	12.82	8.2	ug/L	476982	3	11.48	2895430	50.0
unknown-01	13.12	10.7	ug/L	619633	3	11.48	2895430	50.0
Benzene, (1-methy...	13.18	34.1	ug/L	1976940	3	11.48	2895430	50.0
1H-Indene, 1-methyl-	13.29	42.2	ug/L	2444370	3	11.48	2895430	50.0
Benzo[b]thiophene	13.86	12.1	ug/L	700151	3	11.48	2895430	50.0
(DEL) Alkane: Str...	14.13	4.7	ug/L	273789	3	11.48	2895430	50.0
1H-Indene, 1,1-di...	14.33	4.5	ug/L	260462	3	11.48	2895430	50.0
Naphthalene, 1-et...	15.79	13.3	ug/L	770293	3	11.48	2895430	50.0
Naphthalene, 1,8-...	15.91	73.9	ug/L	4277210	3	11.48	2895430	50.0
Naphthalene, 2,3-...	16.05	63.5	ug/L	3679840	3	11.48	2895430	50.0
Naphthalene, 1,2-...	16.43	10.2	ug/L	589442	3	11.48	2895430	50.0
Biphenylene	16.52	55.3	ug/L	3202910	3	11.48	2895430	50.0
Naphthalene, 2,3,...	17.02	29.5	ug/L	1705970	3	11.48	2895430	50.0
Naphthalene, 1,4,...	17.16	14.6	ug/L	842378	3	11.48	2895430	50.0