

Data Path : W:\HPCHEM1\MSVOA U\DATA\VU050818\
 Data File : VU023741.D
 Acq On : 08 May 2018 16:58
 Operator : MD/SY
 Sample : J2725-10 50X
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 17 Sample Multiplier: 1

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 : W:\HPCHEM1\MSVOA_U\METHOD\SOMULM050818WMA.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	1.088	136	155	177	rBV	1063445	1207687	90.74%	7.524%
2	1.182	180	184	191	rVB	6957	7176	0.54%	0.045%
3	1.397	245	251	268	rBV	187979	192220	14.44%	1.198%
4	1.680	331	339	353	rVB	131160	170133	12.78%	1.060%
5	2.272	512	523	536	rBV	310869	472146	35.47%	2.941%
6	2.413	565	567	570	rBV2	319	173	0.01%	0.001%
7	2.619	625	631	633	rBV2	612	644	0.05%	0.004%
8	2.702	647	657	659	rBV3	732	917	0.07%	0.006%
9	2.783	673	682	686	rVB2	229	269	0.02%	0.002%
10	2.841	686	700	710	rBV	1504	3330	0.25%	0.021%
11	3.030	752	759	763	rBV	274	277	0.02%	0.002%
12	3.223	814	819	821	rBV	204	179	0.01%	0.001%
13	3.281	832	837	840	rBV2	164	189	0.01%	0.001%
14	3.390	867	871	880	rVB3	327	419	0.03%	0.003%
15	3.497	901	904	907	rVB	243	159	0.01%	0.001%
16	3.789	992	995	1001	rVB2	190	190	0.01%	0.001%
17	3.850	1010	1014	1018	rVB	217	189	0.01%	0.001%
18	3.979	1051	1054	1059	rVB2	166	148	0.01%	0.001%
19	4.178	1100	1116	1150	rBV	125637	330017	24.80%	2.056%
20	4.390	1179	1182	1187	rVB3	439	378	0.03%	0.002%
21	4.651	1246	1263	1291	rBV	209239	494454	37.15%	3.080%
22	4.889	1330	1337	1345	rVV3	634	1053	0.08%	0.007%
23	4.998	1368	1371	1378	rVV4	459	380	0.03%	0.002%
24	5.037	1379	1383	1387	rVB2	162	150	0.01%	0.001%
25	5.146	1410	1417	1419	rBV2	147	157	0.01%	0.001%
26	5.345	1454	1479	1513	rBV2	436041	1330957	100.00%	8.292%
27	5.548	1540	1542	1546	rBV2	250	150	0.01%	0.001%
28	5.606	1557	1560	1567	rBV2	210	201	0.02%	0.001%
29	5.725	1591	1597	1600	rBV2	185	184	0.01%	0.001%
30	5.741	1600	1602	1606	rVV2	174	164	0.01%	0.001%
31	5.763	1606	1609	1611	rVV	187	146	0.01%	0.001%
32	5.789	1613	1617	1623	rVB2	250	303	0.02%	0.002%
33	5.889	1633	1648	1672	rBV	403458	802664	60.31%	5.000%
34	6.185	1732	1740	1748	rBV6	1118	1983	0.15%	0.012%

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 Title : VOC Analysis

35	6.336	1772	1787	1821	rBV	288608	595030	44.71%	3.707%
36	6.570	1855	1860	1865	rVB2	196	194	0.01%	0.001%
37	6.869	1949	1953	1958	rVB2	355	423	0.03%	0.003%
38	7.014	1993	1998	2001	rBV2	153	150	0.01%	0.001%
39	7.230	2053	2065	2089	rBV	155454	280356	21.06%	1.747%
40	7.387	2111	2114	2117	rVV2	243	152	0.01%	0.001%
41	7.570	2157	2171	2186	rBV	589625	1038317	78.01%	6.469%
42	7.786	2236	2238	2242	rVB2	496	296	0.02%	0.002%
43	7.805	2242	2244	2248	rBV3	314	189	0.01%	0.001%
44	7.853	2248	2259	2277	rBV	107724	183302	13.77%	1.142%
45	8.056	2315	2322	2325	rBV2	150	161	0.01%	0.001%
46	8.181	2357	2361	2364	rBV	281	253	0.02%	0.002%
47	8.262	2380	2386	2389	rBV2	142	157	0.01%	0.001%
48	8.313	2389	2402	2440	rBV	416558	714864	53.71%	4.454%
49	8.619	2493	2497	2499	rBV3	407	331	0.02%	0.002%
50	9.004	2614	2617	2619	rBV	274	203	0.02%	0.001%
51	9.094	2631	2645	2683	rBV	594859	998700	75.04%	6.222%
52	9.255	2683	2695	2717	rVB	102338	165700	12.45%	1.032%
53	9.377	2722	2733	2747	rBV	64562	110083	8.27%	0.686%
54	9.548	2783	2786	2796	rBV4	618	805	0.06%	0.005%
55	9.783	2848	2859	2877	rVV	51446	82756	6.22%	0.516%
56	9.873	2884	2887	2890	rBV	272	152	0.01%	0.001%
57	10.027	2932	2935	2938	rBV2	214	203	0.02%	0.001%
58	10.172	2970	2980	2990	rBV	7374	10963	0.82%	0.068%
59	10.265	3005	3009	3013	rBV2	506	557	0.04%	0.003%
60	10.435	3050	3062	3083	rBV	399834	635058	47.71%	3.956%
61	10.593	3103	3111	3124	rVB2	2068	3480	0.26%	0.022%
62	10.686	3130	3140	3157	rBV2	10044	22473	1.69%	0.140%
63	10.776	3160	3168	3178	rVB	9194	13110	0.99%	0.082%
64	10.869	3194	3197	3204	rVB	273	236	0.02%	0.001%
65	10.972	3221	3229	3237	rBV	4726	8539	0.64%	0.053%
66	11.152	3271	3285	3298	rBV	55964	86963	6.53%	0.542%
67	11.223	3299	3307	3317	rVV2	3709	5559	0.42%	0.035%
68	11.271	3317	3322	3329	rVB2	1249	1699	0.13%	0.011%
69	11.326	3336	3339	3342	rBV2	583	437	0.03%	0.003%
70	11.406	3353	3364	3377	rBV	27141	44746	3.36%	0.279%
71	11.487	3377	3389	3406	rVV	649111	1028193	77.25%	6.406%

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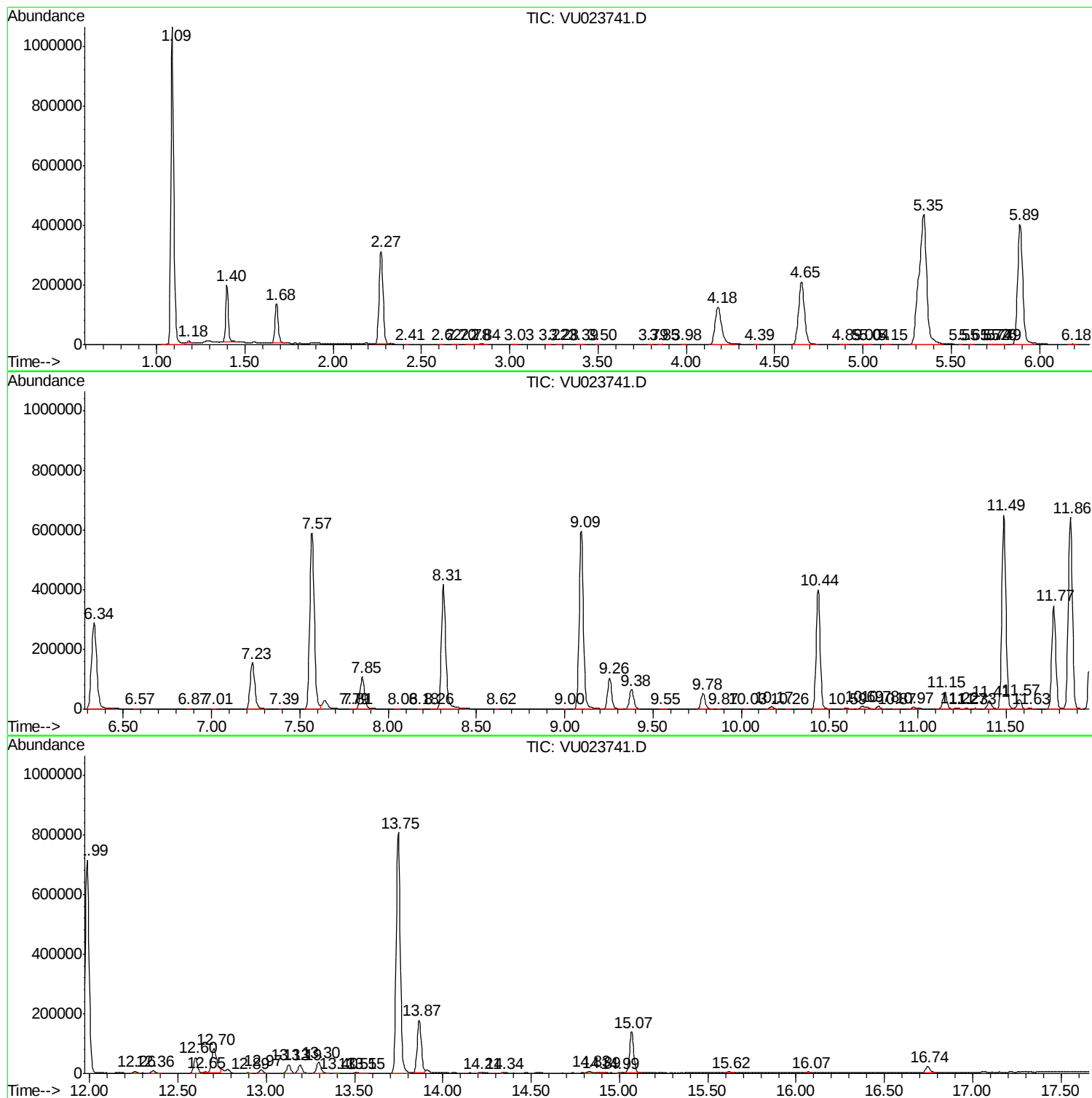
72	11.573	3408	3416	3428	rVV	32070	50164	3.77%	0.313%
73	11.631	3428	3434	3444	rVB4	1678	2630	0.20%	0.016%
74	11.770	3464	3477	3494	rBV	346271	540389	40.60%	3.367%
75	11.863	3494	3506	3526	rVB	640335	1033139	77.62%	6.436%
76	11.985	3531	3544	3570	rBV	714056	1111587	83.52%	6.925%
77	12.255	3621	3628	3635	rBV2	5012	6483	0.49%	0.040%
78	12.358	3651	3660	3672	rBV	7332	11056	0.83%	0.069%
79	12.599	3725	3735	3745	rBV	51887	77084	5.79%	0.480%
80	12.651	3746	3751	3756	rBV4	3634	4427	0.33%	0.028%
81	12.702	3756	3767	3779	rBV	79906	164557	12.36%	1.025%
82	12.895	3824	3827	3833	rVB4	974	877	0.07%	0.005%
83	12.969	3842	3850	3861	rVB3	10681	15360	1.15%	0.096%
84	13.126	3889	3899	3909	rBV3	27794	42378	3.18%	0.264%
85	13.191	3909	3919	3932	rVB	28242	42198	3.17%	0.263%
86	13.297	3940	3952	3969	rBV	37546	61286	4.60%	0.382%
87	13.397	3977	3983	3987	rBV3	626	682	0.05%	0.004%
88	13.506	4014	4017	4027	rVB3	1809	2086	0.16%	0.013%
89	13.551	4027	4031	4038	rBV3	581	779	0.06%	0.005%
90	13.747	4079	4092	4113	rBV	805525	1273157	95.66%	7.932%
91	13.866	4118	4129	4139	rBV	174400	275982	20.74%	1.719%
92	14.210	4231	4236	4239	rBV4	796	947	0.07%	0.006%
93	14.335	4272	4275	4277	rBV2	662	412	0.03%	0.003%
94	14.827	4420	4428	4438	rBV2	5996	8164	0.61%	0.051%
95	14.885	4440	4446	4457	rVB4	2293	3914	0.29%	0.024%
96	14.995	4477	4480	4483	rBV2	871	752	0.06%	0.005%
97	15.069	4491	4503	4519	rBV	137370	225348	16.93%	1.404%
98	15.618	4668	4674	4684	rVB3	3820	5366	0.40%	0.033%
99	16.072	4809	4815	4819	rBV4	2507	2700	0.20%	0.017%
100	16.744	5016	5024	5038	rBV3	20037	33233	2.50%	0.207%

Sum of corrected areas: 16051713

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Quant Method : W:\HPCHEM1\MSVOA_U\METHOD\SOMULM050818WMA.M
Quant Title : VOC Analysis

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



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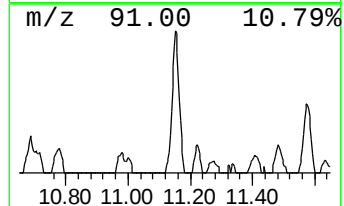
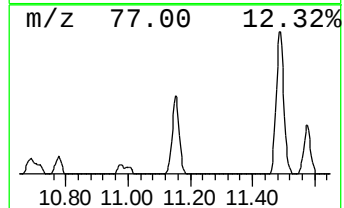
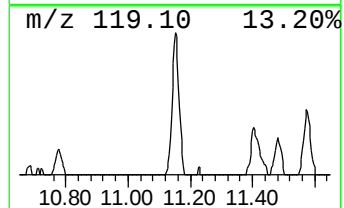
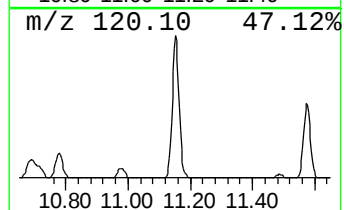
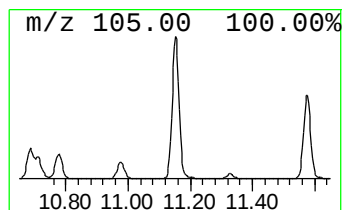
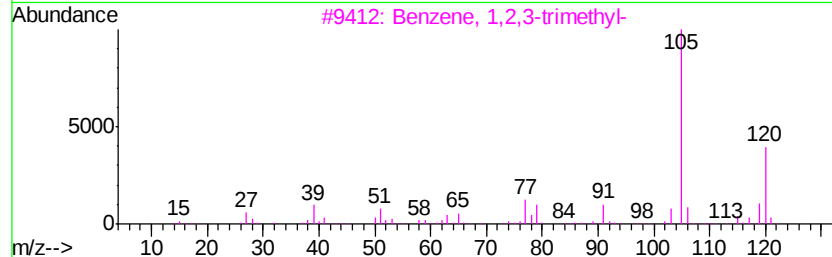
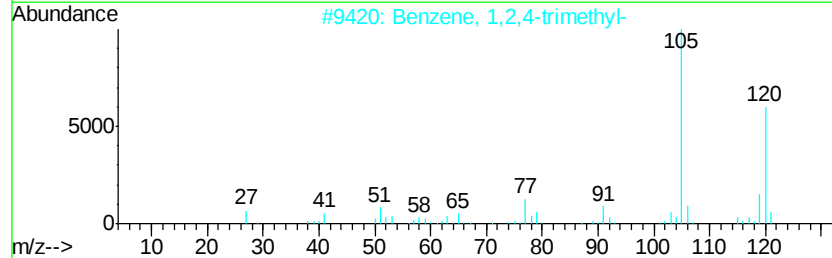
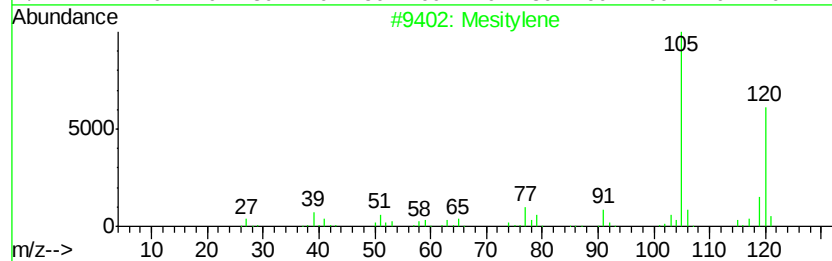
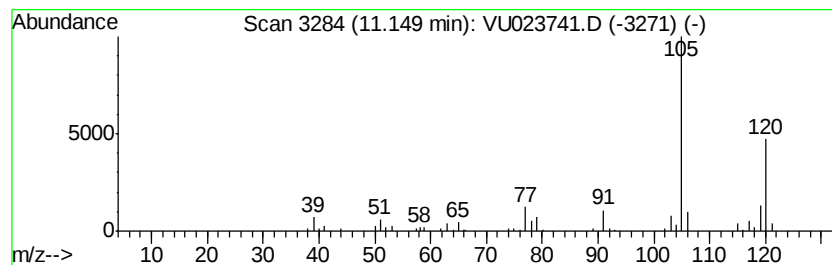
Quant Method : W:\HPCHEM1\MSVOA_U\METHOD\SOMULM050818WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 Mesitylene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.15	4.23 ug/L	86963	1,4-Dichlorobenzene-d4	11.49

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



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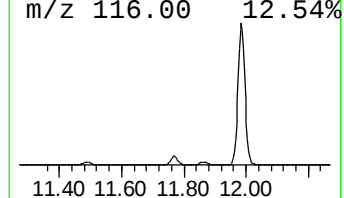
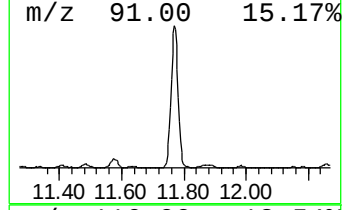
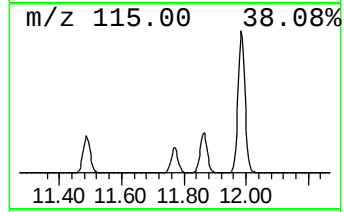
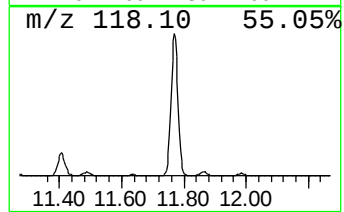
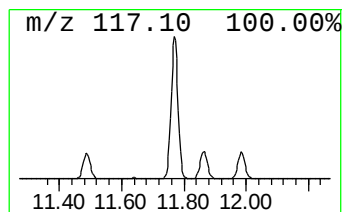
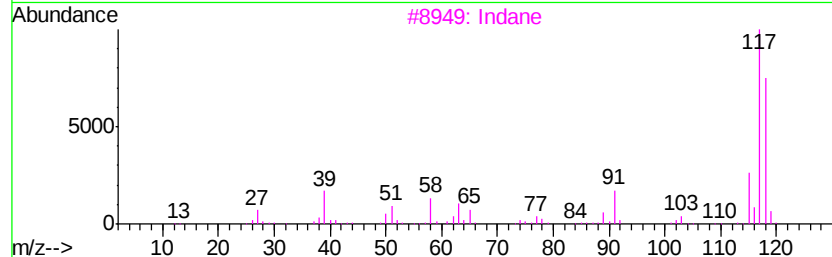
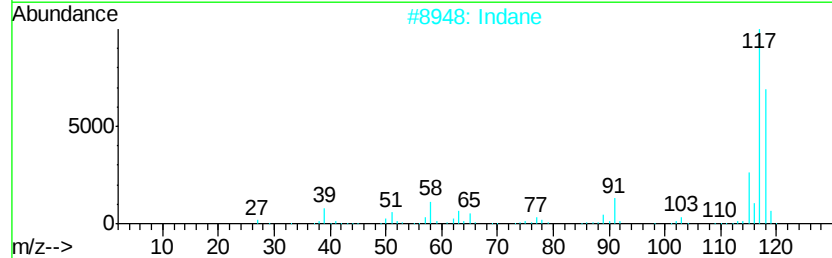
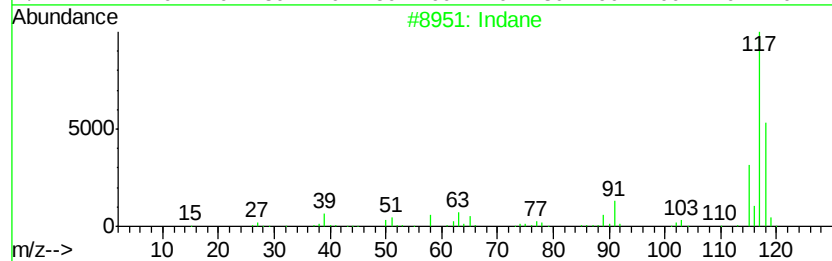
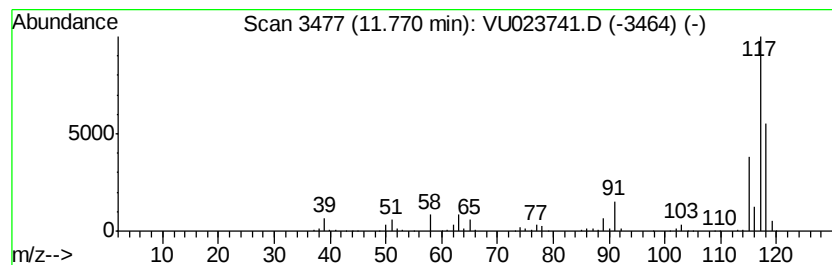
Quant Method : W:\HPCHEM1\MSVOA_U\METHOD\SOMULM050818WMA.M
Quant Title : VOC Analysis

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

Peak Number 4 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	26.28 ug/L	540389	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	95
2		Indane	118	C9H10	000496-11-7	95
3		Indane	118	C9H10	000496-11-7	89
4		Indane	118	C9H10	000496-11-7	87
5		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	68



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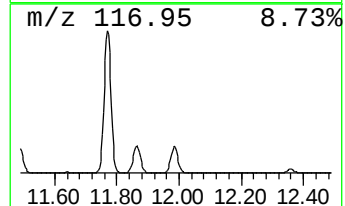
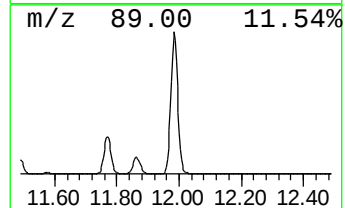
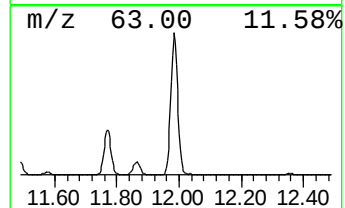
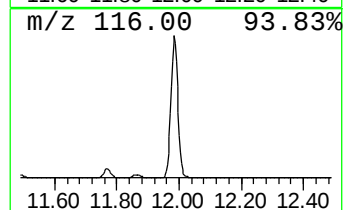
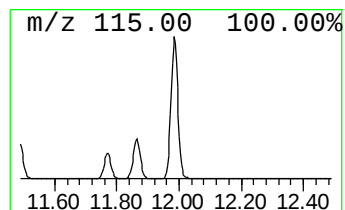
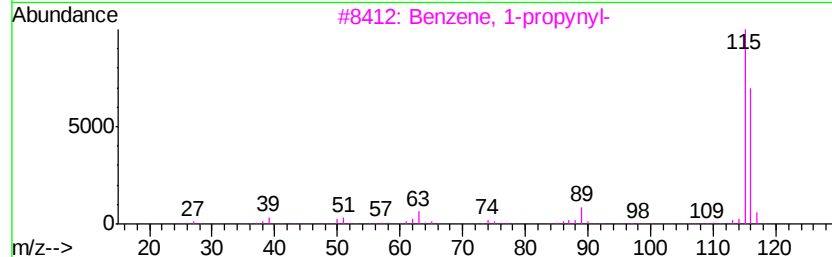
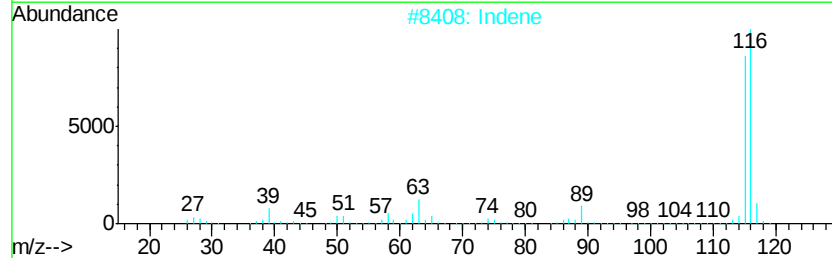
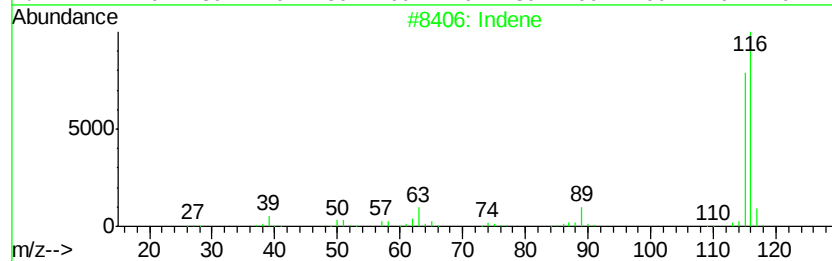
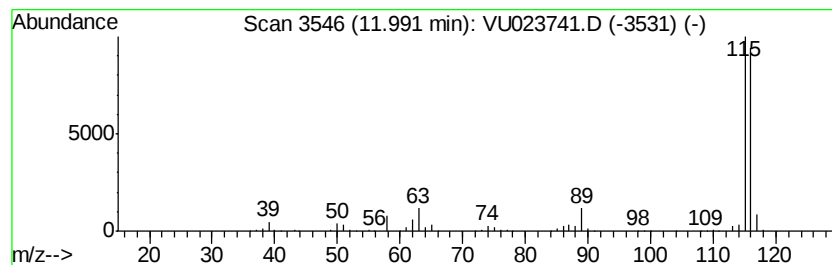
Quant Method : W:\HPCHEM1\MSVOA_U\METHOD\SOMULM050818WMA.M
Quant Title : VOC Analysis

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

Peak Number 5 Indene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.99	54.06 ug/L	1111590	1,4-Dichlorobenzene-d4	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene		116	C9H8	000095-13-6	97
2	Indene		116	C9H8	000095-13-6	95
3	Benzene, 1-propynyl-		116	C9H8	000673-32-5	94
4	Indene		116	C9H8	000095-13-6	94
5	3-Methylphenylacetylene		116	C9H8	000766-82-5	91



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ALS Vial : 17 Sample Multiplier: 1

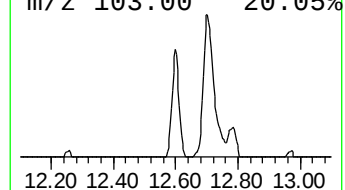
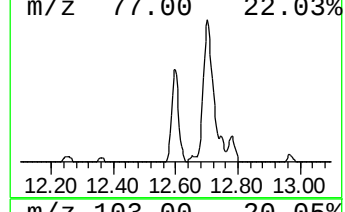
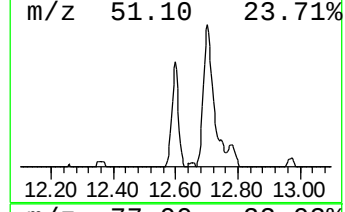
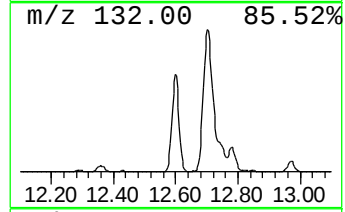
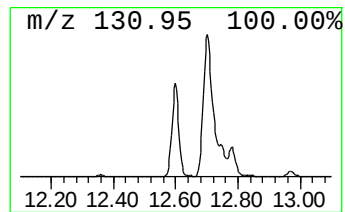
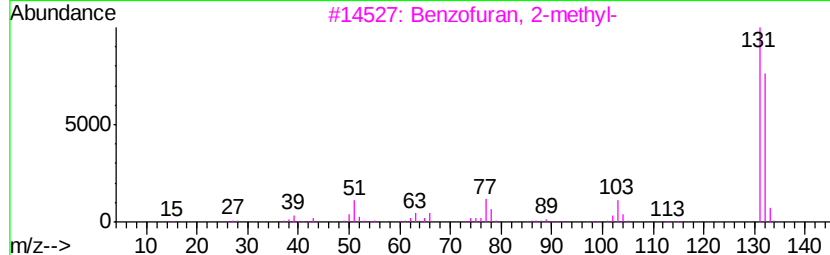
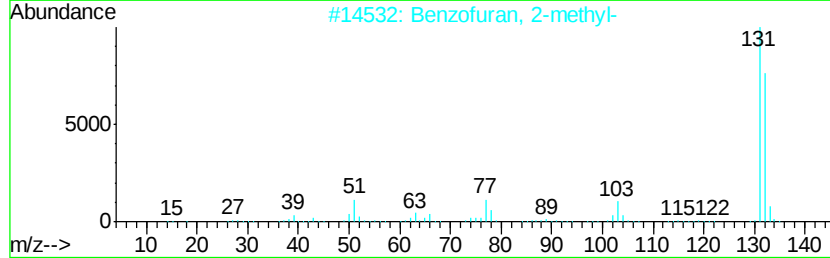
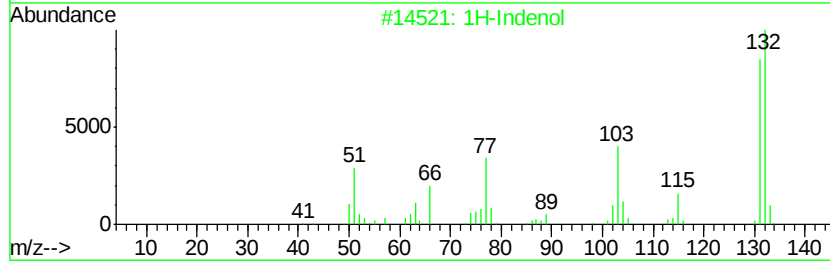
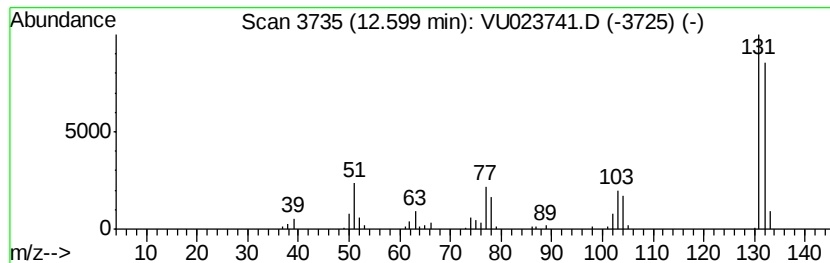
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Quant Title : VOC Analysis

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

Peak Number 6 1H-Indenol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.60	3.75 ug/L	77084	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indenol	132	C9H8O	056631-57-3	91
2		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
3		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
4		3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91
5		2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	90



Data Path : W:\HPCHEM1\MSVOA U\DATA\VU050818\
Data File : VU023741.D
Acq On : 08 May 2018 16:58
Operator : MD/SY
Sample : J2725-10 50X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

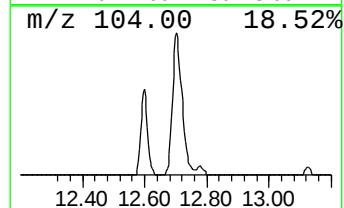
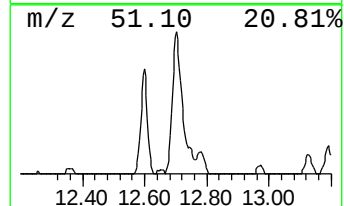
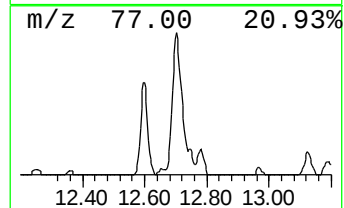
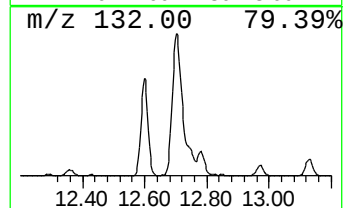
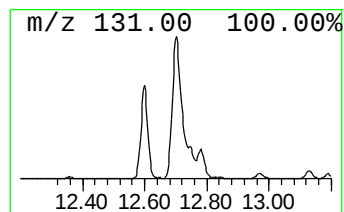
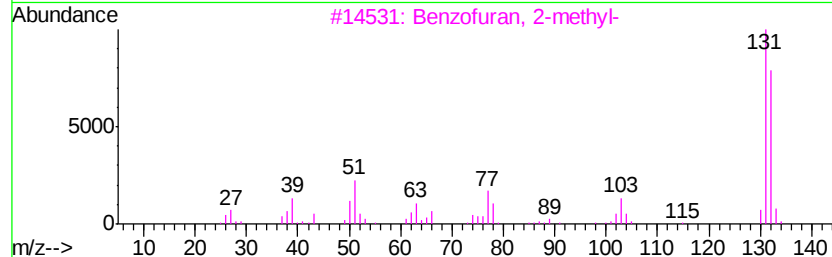
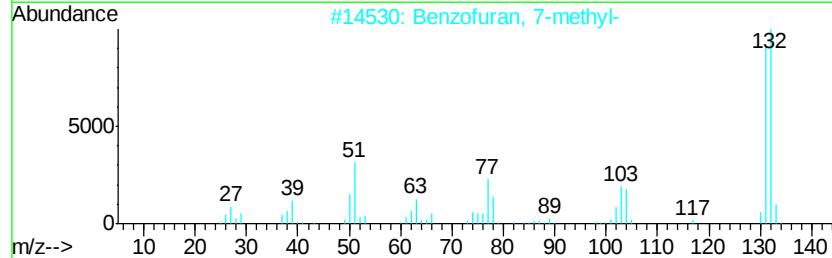
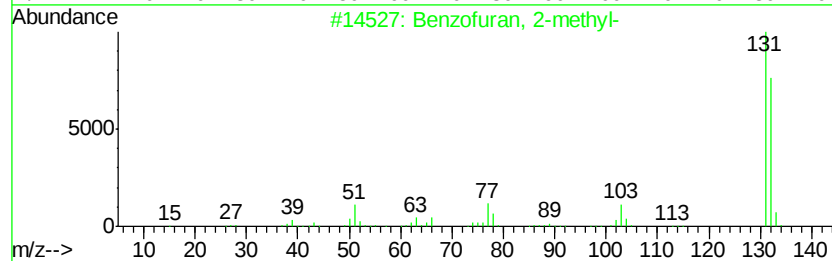
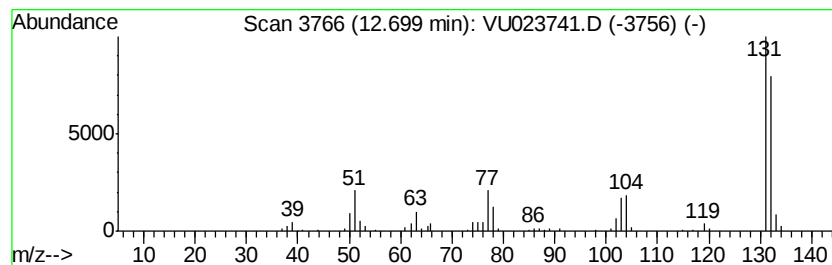
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Quant Title : VOC Analysis

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

Peak Number 7 Benzofuran, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	8.00 ug/L	164557	1,4-Dichlorobenzene-d4	11.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2			Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	93
3			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
4			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
5			3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	90



Data Path : W:\HPCHEM1\MSVOA U\DATA\VU050818\
Data File : VU023741.D
Acq On : 08 May 2018 16:58
Operator : MD/SY
Sample : J2725-10 50X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

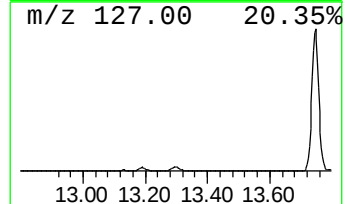
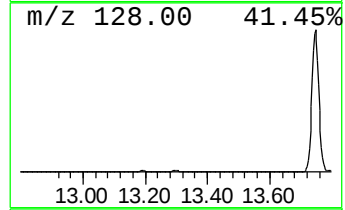
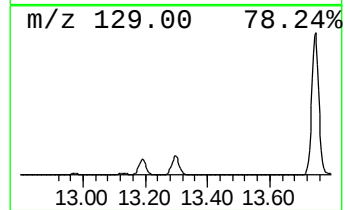
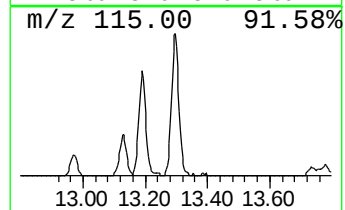
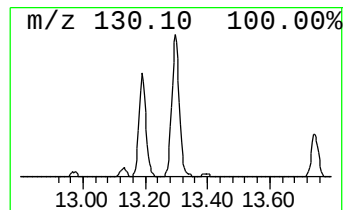
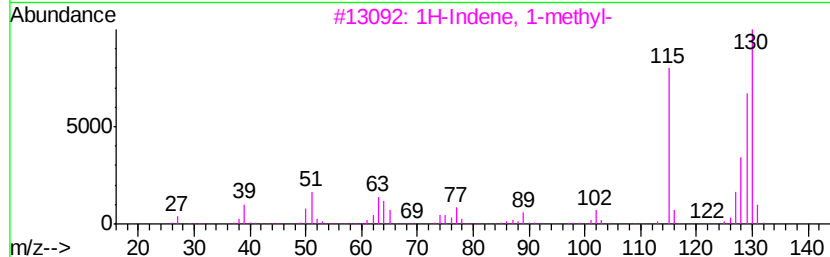
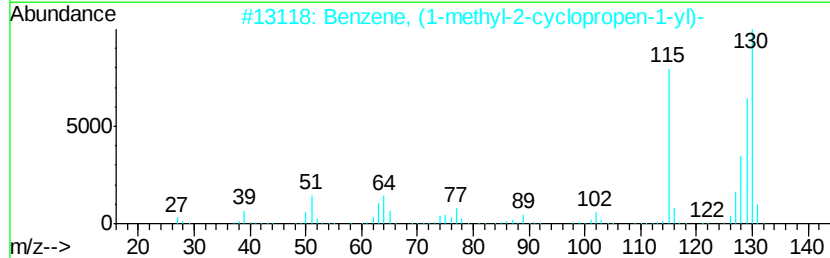
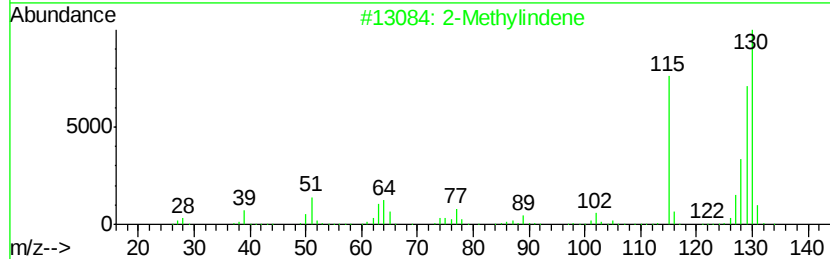
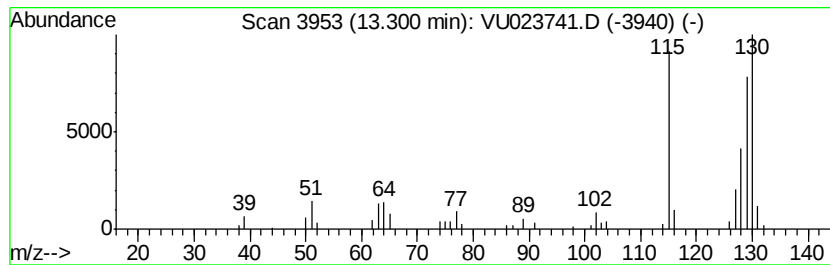
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Quant Title : VOC Analysis

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

Peak Number 8 2-Methylindene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.30	2.98 ug/L	61286	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Methylindene	130	C10H10	002177-47-1	97
2		Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	96
3		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	95
4		Cycloprop[alindene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	95
5		1,4-Dihydronaphthalene	130	C10H10	000612-17-9	93



Data Path : W:\HPCHEM1\MSVOA U\DATA\VU050818\
Data File : VU023741.D
Acq On : 08 May 2018 16:58
Operator : MD/SY
Sample : J2725-10 50X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

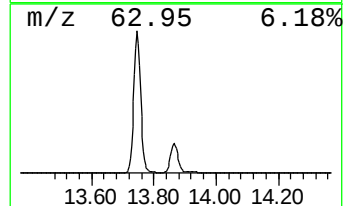
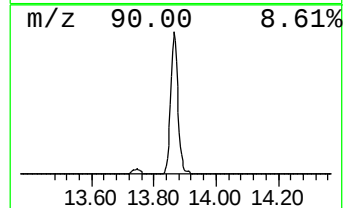
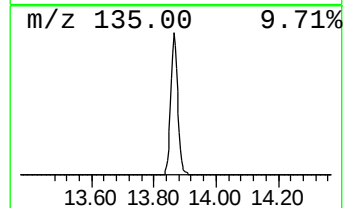
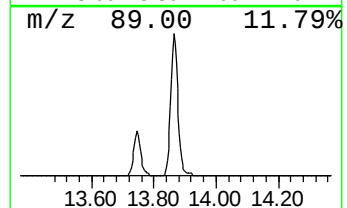
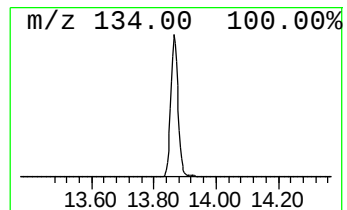
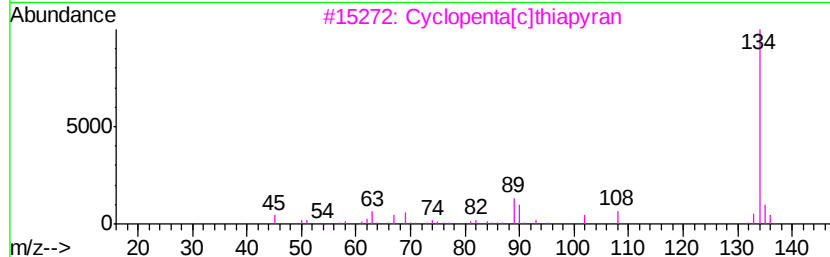
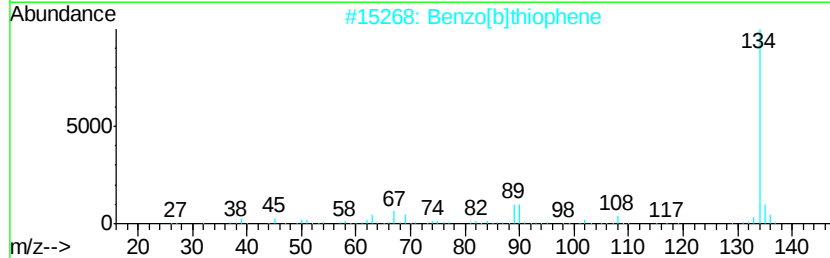
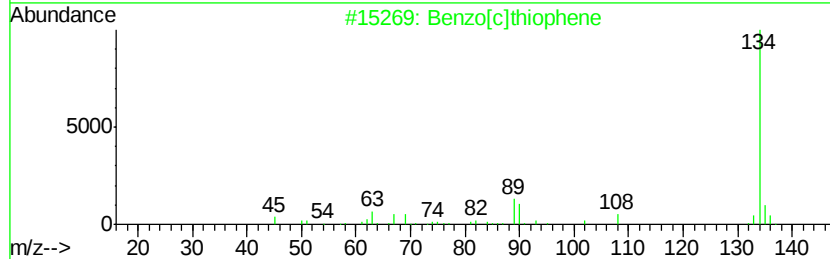
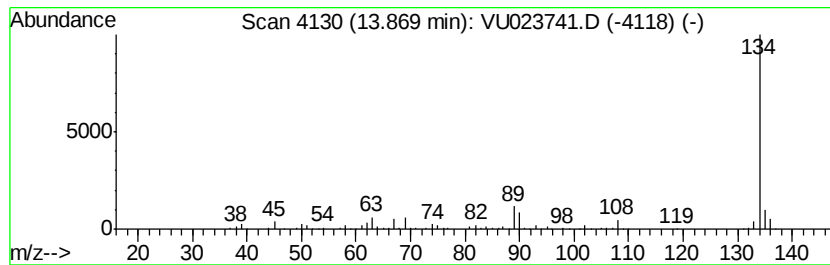
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Quant Title : VOC Analysis

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

Peak Number 10 Benzo[c]thiophene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	13.42 ug/L	275982	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[c]thiophene	134	C8H6S	000270-82-6	97
2		Benzo[b]thiophene	134	C8H6S	000095-15-8	95
3		Cyclopenta[c]thiopyran	134	C8H6S	000270-63-3	94
4		Benzo[b]thiophene	134	C8H6S	000095-15-8	93
5		Benzo[b]thiophene	134	C8H6S	000095-15-8	91



Data Path : W:\HPCHEM1\MSVOA U\DATA\VU050818\
Data File : VU023741.D
Acq On : 08 May 2018 16:58
Operator : MD/SY
Sample : J2725-10 50X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

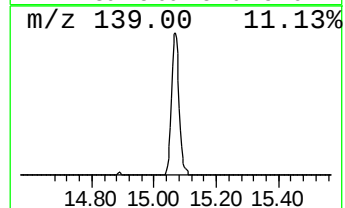
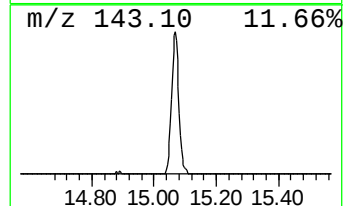
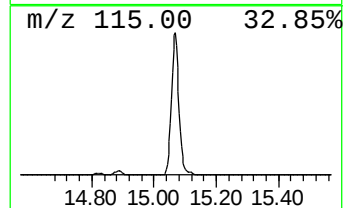
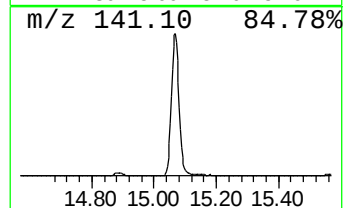
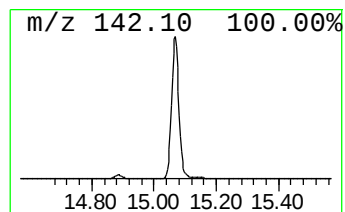
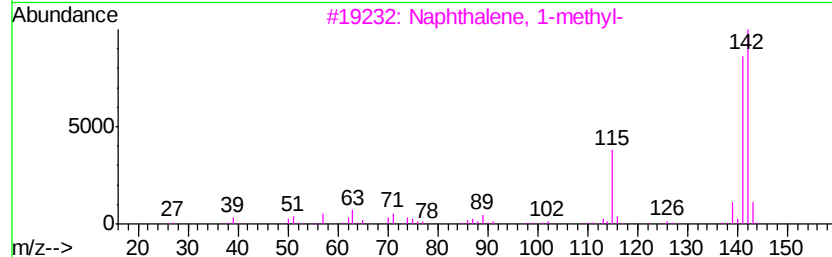
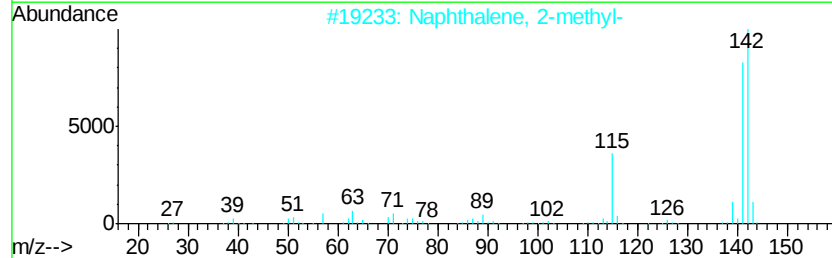
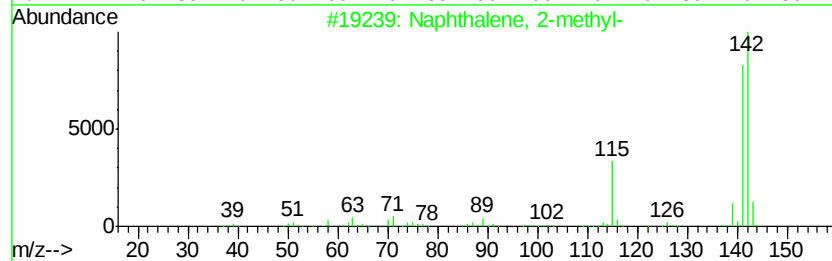
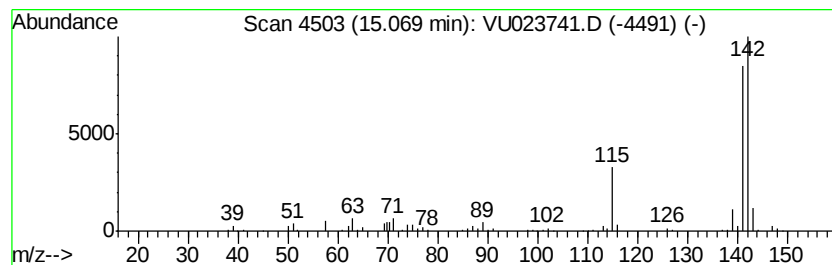
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Quant Title : VOC Analysis

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

Peak Number 11 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	10.96 ug/L	225348	1,4-Dichlorobenzene-d4	11.49

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, 1-methyl-	142	C11H10	000090-12-0	94
5			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94



Data Path : W:\HPCHEM1\MSVOA_U\DATA\VU050818\
Data File : VU023741.D
Acq On : 08 May 2018 16:58
Operator : MD/SY
Sample : J2725-10 50X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Quant Method : W:\HPCHEM1\MSVOA_U\METHOD\SOMULM050818WMA.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
Mesitylene	11.15	4.2	ug/L	86963	3	11.49	1028190	50.0
Indane	11.77	26.3	ug/L	540389	3	11.49	1028190	50.0
Indene	11.99	54.1	ug/L	1111590	3	11.49	1028190	50.0
1H-Indenol	12.60	3.8	ug/L	77084	3	11.49	1028190	50.0
Benzofuran, 2-met...	12.70	8.0	ug/L	164557	3	11.49	1028190	50.0
2-Methylindene	13.30	3.0	ug/L	61286	3	11.49	1028190	50.0
Benzo[c]thiophene	13.87	13.4	ug/L	275982	3	11.49	1028190	50.0
Naphthalene, 1-me...	15.07	11.0	ug/L	225348	3	11.49	1028190	50.0