

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU040719\
 Data File : VU030840.D
 Acq On : 07 Apr 2019 06:51
 Operator : JC/SP
 Sample : K2249-14
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 E3YE6

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\SOMULM040519WMA.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	1.180	23	28	35	rVB2	15220	14660	0.05%	0.016%
2	1.399	87	96	110	rBV2	1593595	1684640	5.35%	1.872%
3	1.695	175	188	206	rVB2	791345	1277758	4.06%	1.420%
4	1.830	224	230	239	rBV	16222	21719	0.07%	0.024%
5	2.273	356	368	394	rVB3	721602	1363812	4.33%	1.516%
6	2.441	413	420	436	rVB2	13693	24914	0.08%	0.028%
7	2.617	467	475	488	rBV2	6251	10661	0.03%	0.012%
8	2.698	492	500	506	rVV5	4353	5992	0.02%	0.007%
9	2.823	533	539	541	rBV3	2115	1902	0.01%	0.002%
10	2.891	556	560	563	rBV	644	476	0.00%	0.001%
11	2.913	564	567	572	rVB3	906	561	0.00%	0.001%
12	2.981	575	588	610	rBV	112920	216326	0.69%	0.240%
13	3.077	615	618	622	rVV2	1213	893	0.00%	0.001%
14	3.225	662	664	667	rVB3	799	493	0.00%	0.001%
15	3.296	674	686	688	rBV4	1102	1639	0.01%	0.002%
16	3.440	711	731	765	rBV	6504512	14384259	45.71%	15.986%
17	4.219	948	973	1014	rBV	12766465	31471296	100.00%	34.977%
18	4.643	1093	1105	1130	rVB	324790	812118	2.58%	0.903%
19	4.910	1169	1188	1226	rBV	3880599	9293454	29.53%	10.329%
20	5.334	1299	1320	1359	rVB2	689499	2120025	6.74%	2.356%
21	5.723	1437	1441	1444	rBV4	632	639	0.00%	0.001%
22	5.794	1459	1463	1469	rBV4	1506	1810	0.01%	0.002%
23	5.881	1477	1490	1512	rBV	592124	1176348	3.74%	1.307%
24	6.186	1574	1585	1605	rVB7	19878	41596	0.13%	0.046%
25	6.325	1614	1628	1648	rBV2	453782	931276	2.96%	1.035%
26	6.569	1702	1704	1708	rVB4	1260	722	0.00%	0.001%
27	6.768	1762	1766	1773	rVB5	683	813	0.00%	0.001%
28	6.826	1778	1784	1786	rBV5	889	741	0.00%	0.001%
29	6.910	1805	1810	1816	rVB5	1301	1524	0.00%	0.002%
30	6.955	1821	1824	1827	rBV2	876	579	0.00%	0.001%
31	7.116	1867	1874	1875	rBV3	614	634	0.00%	0.001%
32	7.225	1893	1908	1929	rBV	213797	395060	1.26%	0.439%
33	7.402	1960	1963	1967	rBV4	1105	872	0.00%	0.001%
34	7.466	1973	1983	1993	rBV4	11860	20352	0.06%	0.023%

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Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM040519WMA.M
 Title : VOC Analysis

35	7.563	1999	2013	2025	rBV	865364	1501138	4.77%	1.668%
36	7.633	2025	2035	2054	rVB	726284	1260194	4.00%	1.401%
37	7.845	2086	2101	2119	rBV3	161928	349111	1.11%	0.388%
38	8.019	2151	2155	2159	rBV6	820	716	0.00%	0.001%
39	8.071	2167	2171	2182	rVB9	4049	6338	0.02%	0.007%
40	8.161	2195	2199	2202	rBV2	1759	1644	0.01%	0.002%
41	8.225	2208	2219	2234	rBV3	74451	130467	0.41%	0.145%
42	8.308	2234	2245	2283	rBV	615405	1185276	3.77%	1.317%
43	8.633	2341	2346	2348	rBV5	761	730	0.00%	0.001%
44	8.778	2389	2391	2394	rBV3	797	539	0.00%	0.001%
45	8.839	2407	2410	2414	rVB3	953	657	0.00%	0.001%
46	8.881	2420	2423	2427	rVB3	716	539	0.00%	0.001%
47	8.955	2441	2446	2450	rBV5	1417	1413	0.00%	0.002%
48	8.984	2450	2455	2456	rBV3	2829	1816	0.01%	0.002%
49	9.019	2464	2466	2470	rVB2	885	595	0.00%	0.001%
50	9.087	2474	2487	2515	rBV	884944	1492150	4.74%	1.658%
51	9.247	2526	2537	2559	rBV	516721	844824	2.68%	0.939%
52	9.370	2563	2575	2611	rVB	1621274	2677180	8.51%	2.975%
53	9.775	2687	2701	2731	rBV	1008273	1624273	5.16%	1.805%
54	9.923	2744	2747	2748	rBV2	1212	717	0.00%	0.001%
55	9.945	2751	2754	2755	rBV3	926	543	0.00%	0.001%
56	10.164	2811	2822	2834	rBV	63467	104222	0.33%	0.116%
57	10.318	2860	2870	2878	rVV9	5257	8493	0.03%	0.009%
58	10.369	2882	2886	2889	rVV2	1266	1124	0.00%	0.001%
59	10.427	2892	2904	2928	rBV	606880	985787	3.13%	1.096%
60	10.585	2942	2953	2969	rVV	164536	258651	0.82%	0.287%
61	10.678	2970	2982	3000	rVV	756120	1612676	5.12%	1.792%
62	10.768	3001	3010	3029	rVB	454261	704501	2.24%	0.783%
63	10.861	3037	3039	3041	rBV2	854	492	0.00%	0.001%
64	10.913	3049	3055	3061	rBV4	10440	15212	0.05%	0.017%
65	10.968	3061	3072	3097	rVB	562440	887064	2.82%	0.986%
66	11.064	3099	3102	3103	rBV3	1050	606	0.00%	0.001%
67	11.093	3106	3111	3115	rBV8	2643	3397	0.01%	0.004%
68	11.144	3115	3127	3148	rBV	2234362	3386216	10.76%	3.763%
69	11.424	3202	3214	3221	rBV2	31719	51221	0.16%	0.057%
70	11.479	3221	3231	3247	rVV	884921	1409919	4.48%	1.567%
71	11.566	3247	3258	3286	rVB	910886	1428410	4.54%	1.588%

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72	11.762	3308	3319	3330	rBV	293087	497642	1.58%	0.553%
73	11.855	3338	3348	3375	rVB	830163	1453630	4.62%	1.616%
74	11.980	3381	3387	3399	rVB9	10365	17777	0.06%	0.020%
75	12.051	3403	3409	3419	rVB	16717	26422	0.08%	0.029%
76	12.144	3427	3438	3461	rBV3	142884	292565	0.93%	0.325%
77	12.250	3463	3471	3478	rBV2	34592	52100	0.17%	0.058%
78	12.353	3493	3503	3521	rBV3	34356	77907	0.25%	0.087%
79	12.479	3540	3542	3543	rBV	1294	619	0.00%	0.001%
80	12.549	3552	3564	3573	rBV5	19274	32909	0.10%	0.037%
81	12.649	3588	3595	3603	rVB3	16449	23474	0.07%	0.026%
82	12.701	3603	3611	3627	rVB2	30469	49254	0.16%	0.055%
83	12.784	3635	3637	3641	rVB4	1549	912	0.00%	0.001%
84	12.919	3677	3679	3681	rBV3	955	589	0.00%	0.001%
85	12.961	3685	3692	3704	rVB3	7663	13314	0.04%	0.015%
86	13.122	3731	3742	3758	rVV3	47340	81192	0.26%	0.090%
87	13.286	3784	3793	3809	rBV4	17548	30486	0.10%	0.034%
88	13.382	3819	3823	3826	rBV4	1880	1681	0.01%	0.002%
89	13.498	3856	3859	3860	rBV2	943	560	0.00%	0.001%
90	13.520	3860	3866	3870	rVB6	2051	2407	0.01%	0.003%
91	13.588	3884	3887	3890	rBV	1085	610	0.00%	0.001%
92	13.607	3891	3893	3897	rBV4	923	724	0.00%	0.001%
93	13.749	3926	3937	3961	rBV	36940	94485	0.30%	0.105%
94	14.147	4058	4061	4065	rVB4	1202	964	0.00%	0.001%
95	14.176	4065	4070	4073	rBV5	1672	1822	0.01%	0.002%
96	14.241	4088	4090	4094	rBV3	1193	954	0.00%	0.001%
97	14.347	4121	4123	4126	rBV	1119	627	0.00%	0.001%
98	15.016	4328	4331	4335	rVB3	1006	686	0.00%	0.001%
99	15.289	4414	4416	4417	rBV2	1178	506	0.00%	0.001%
100	15.324	4423	4427	4431	rBV5	1596	1488	0.00%	0.002%

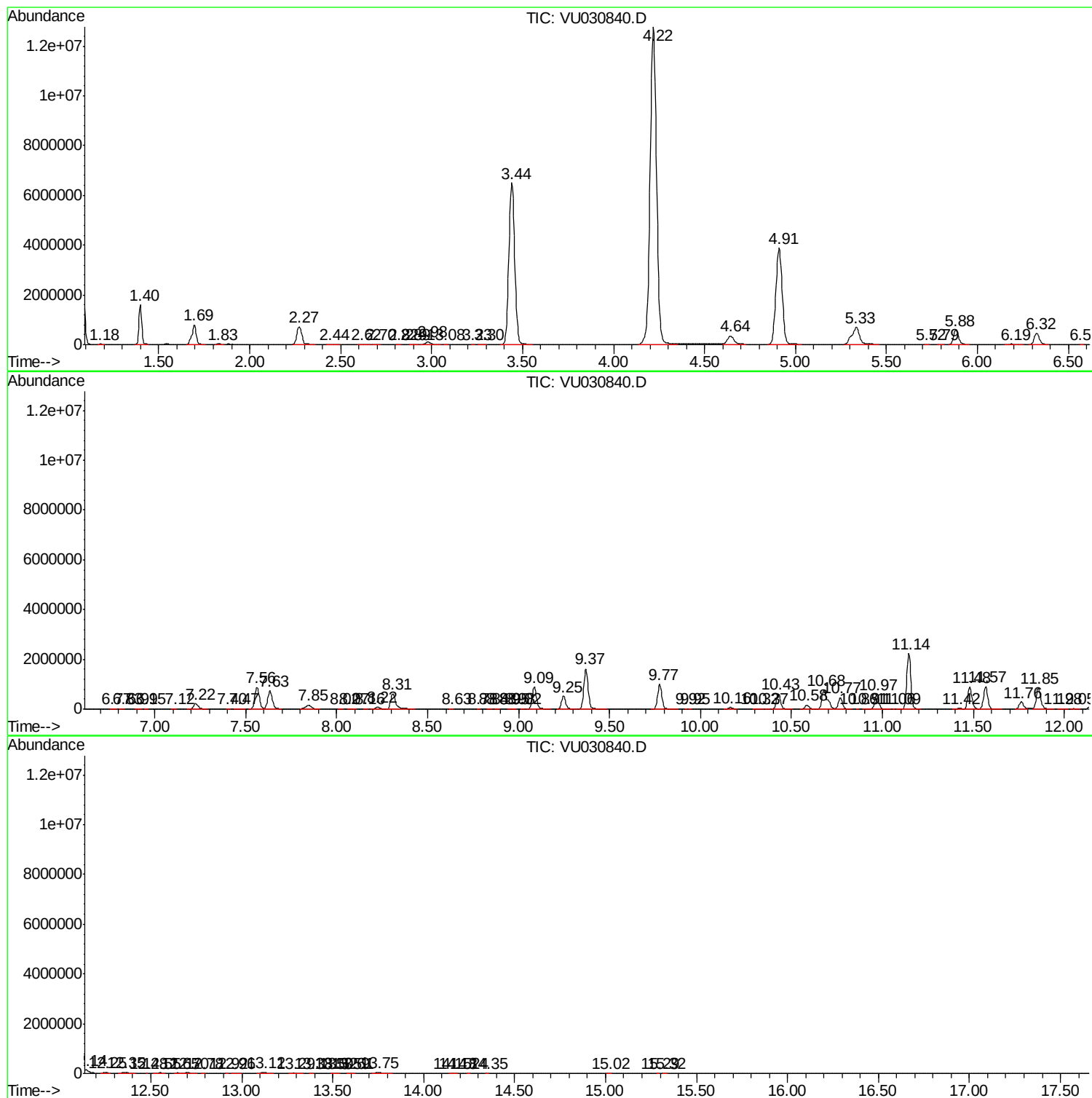
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM040519WMA.M
Quant Title : VOC Analysis

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



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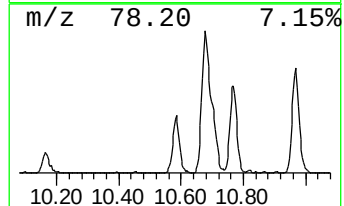
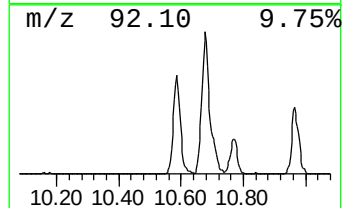
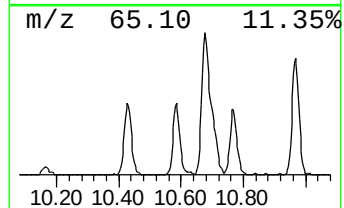
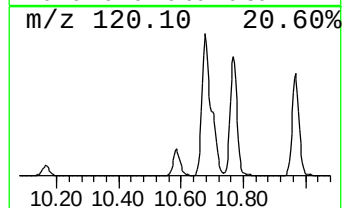
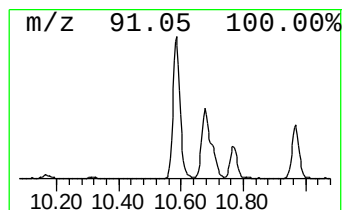
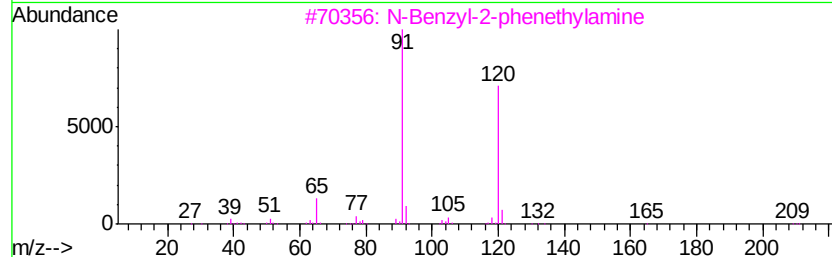
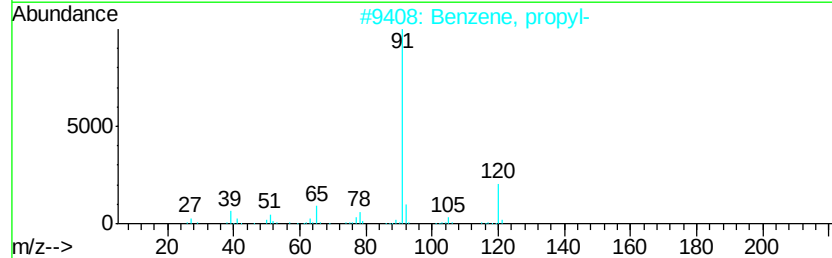
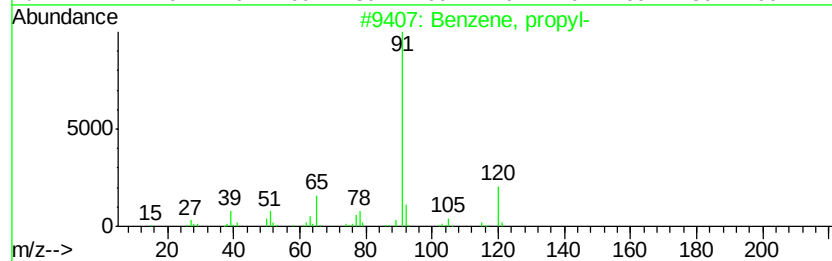
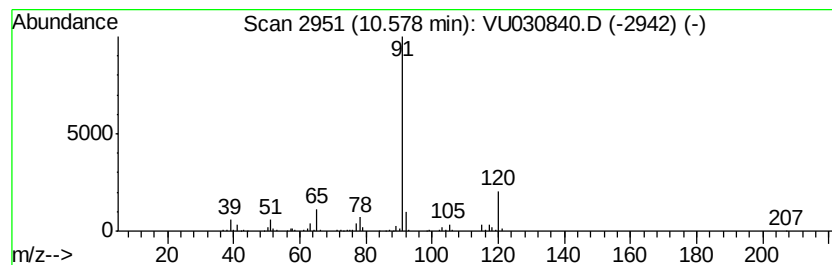
Instrument :
MSVOA_U
ClientSampled :
E3YE6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM040519WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 Benzene, propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.58	9.17 ug/L	258651	1,4-Dichlorobenzene-d4	11.48
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, propyl-	120 C9H12	000103-65-1 90
2		Benzene, propyl-	120 C9H12	000103-65-1 87
3		N-Benzyl-2-phenethylamine	211 C15H17N	003647-71-0 78
4		N-Benzyl-2-phenethylamine	211 C15H17N	003647-71-0 78
5		Benzene, propyl-	120 C9H12	000103-65-1 72



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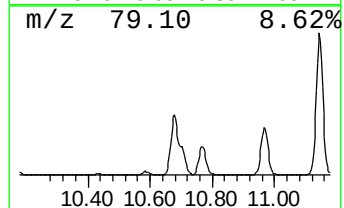
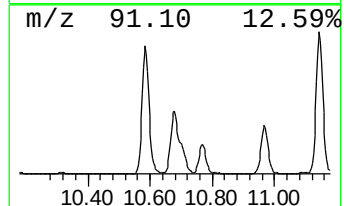
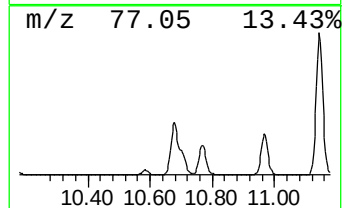
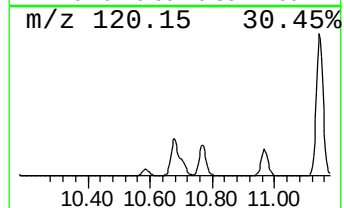
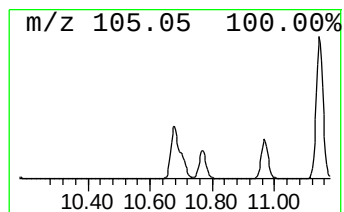
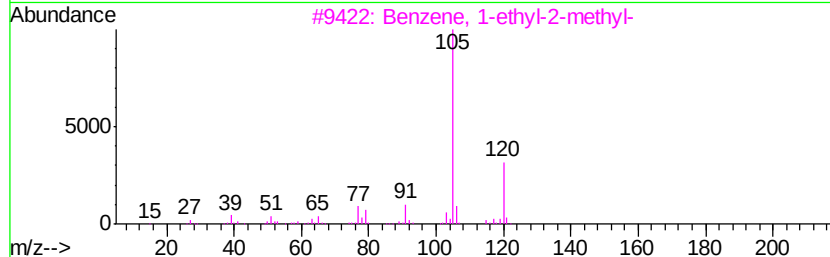
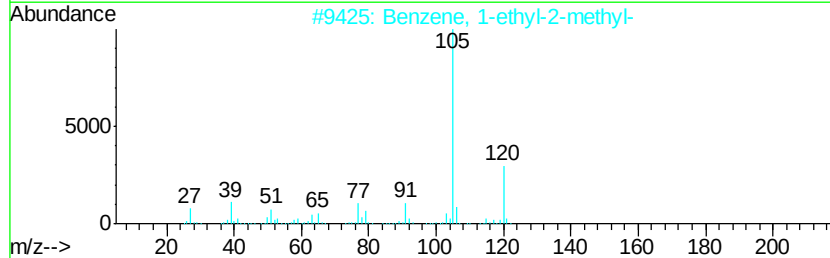
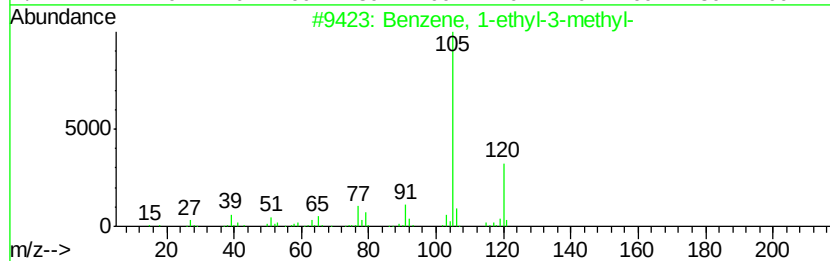
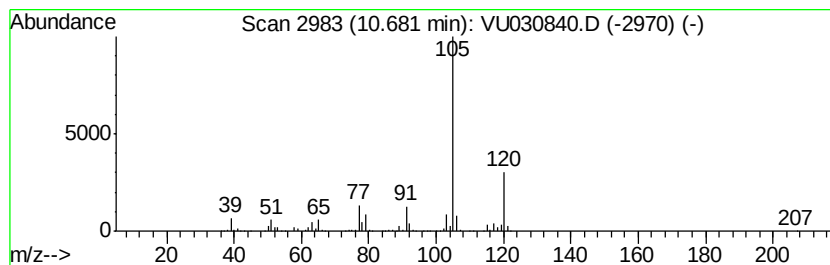
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Quant Title : VOC Analysis

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.68	57.19 ug/L	1612680	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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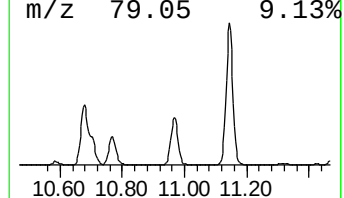
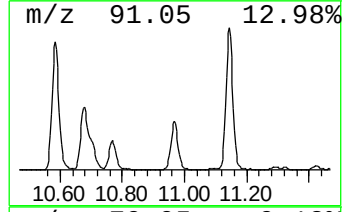
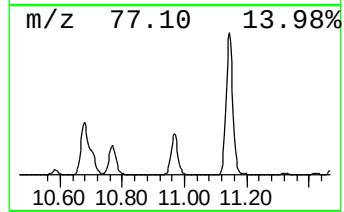
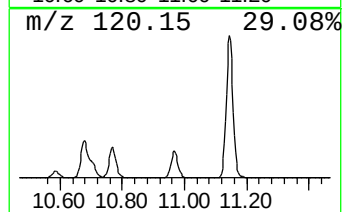
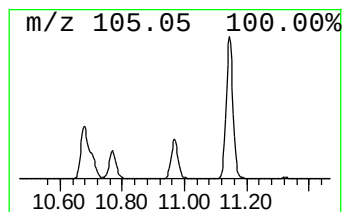
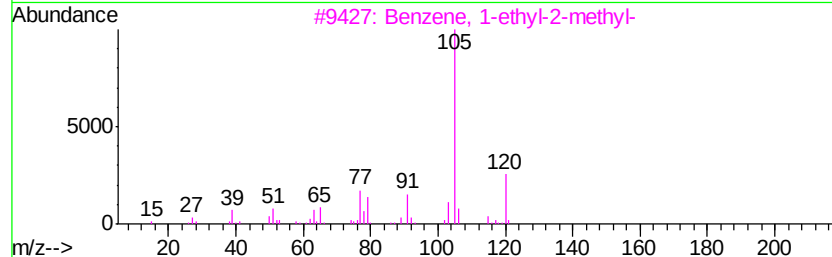
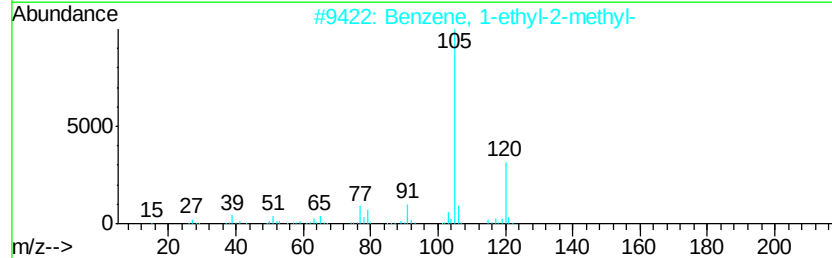
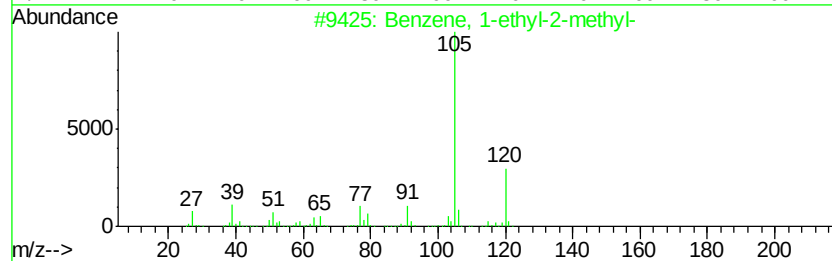
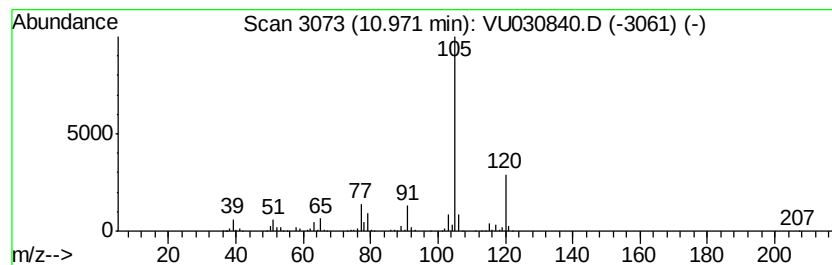
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Quant Title : VOC Analysis

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Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.97	31.46 ug/L	887064	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU040719\
Data File : VU030840.D
Acq On : 07 Apr 2019 06:51
Operator : JC/SP
Sample : K2249-14
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E3YE6

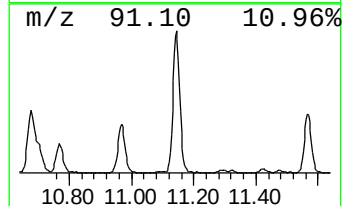
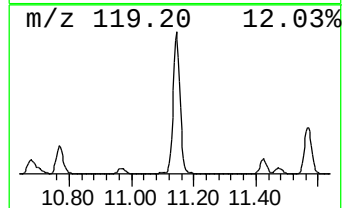
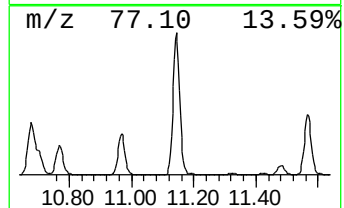
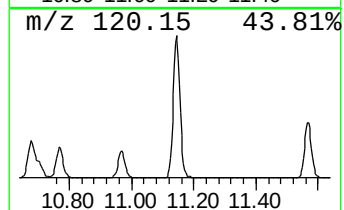
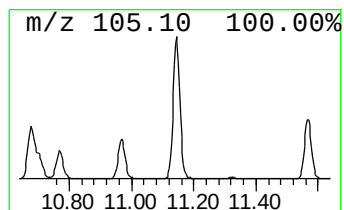
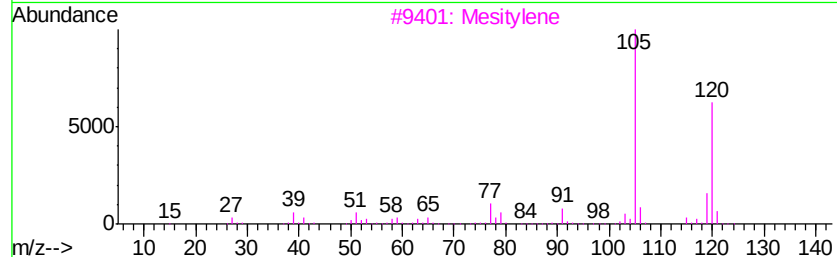
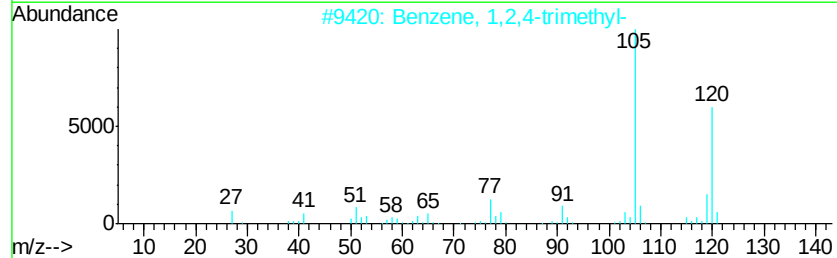
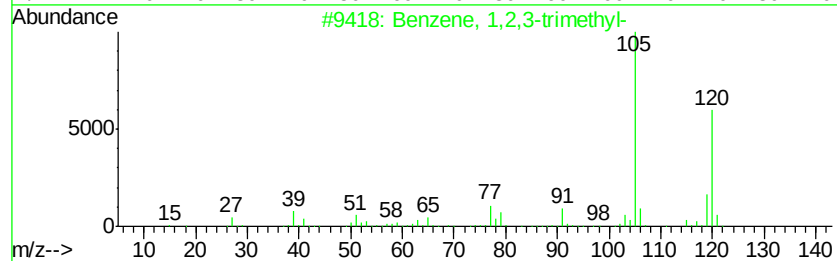
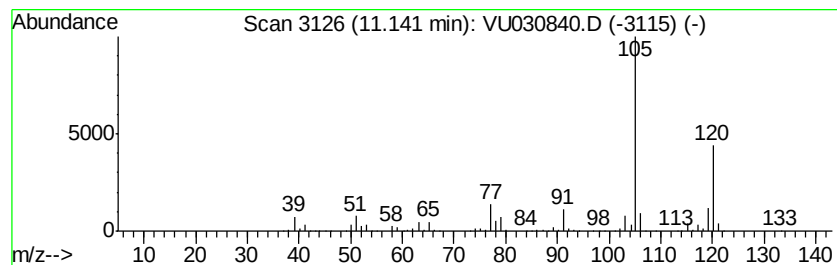
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Quant Title : VOC Analysis

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

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.14	120.09 ug/L	3386220	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		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3		Mesitylene	120	C9H12	000108-67-8	94
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU040719\
Data File : VU030840.D
Acq On : 07 Apr 2019 06:51
Operator : JC/SP
Sample : K2249-14
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E3YE6

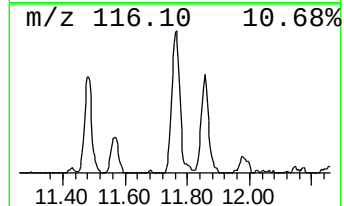
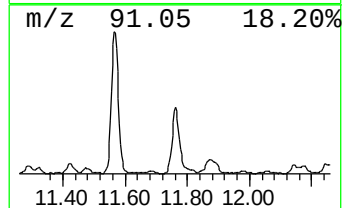
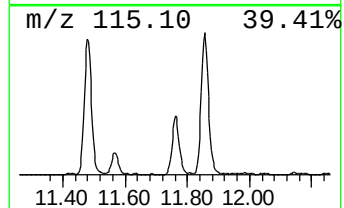
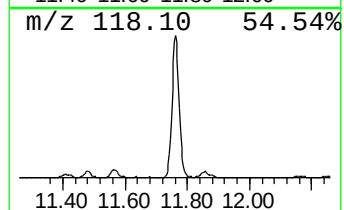
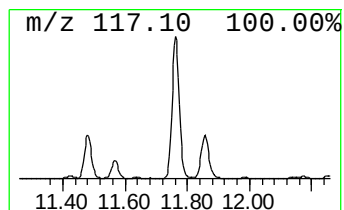
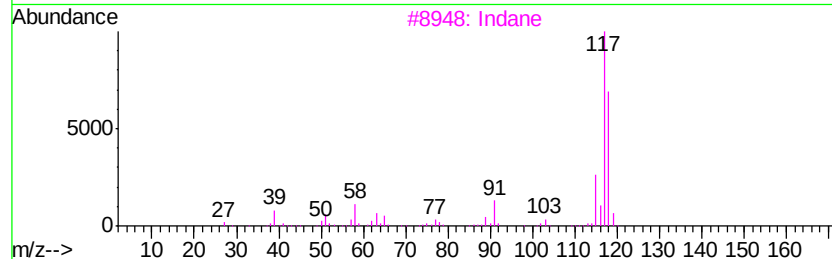
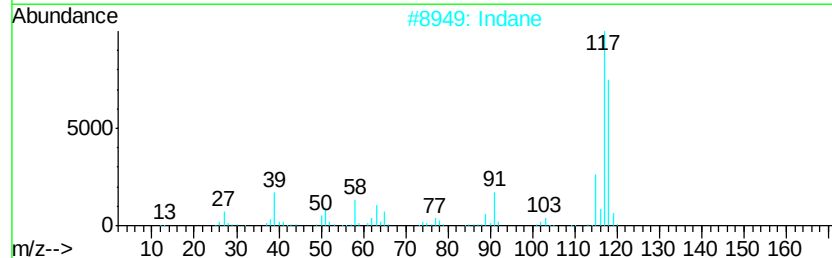
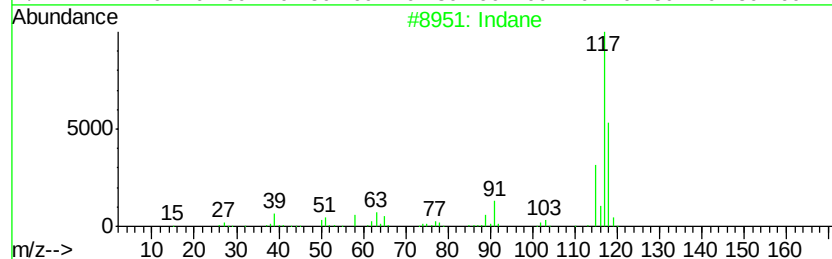
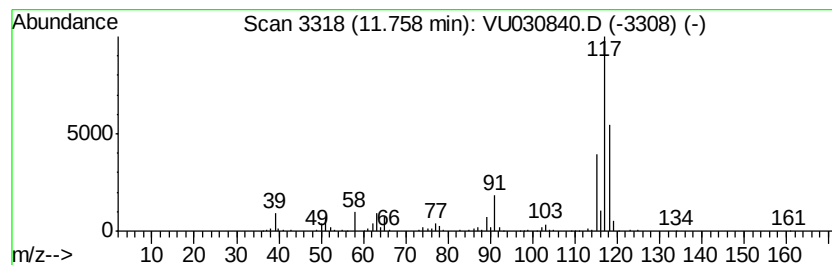
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Quant Title : VOC Analysis

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

Peak Number 8 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.76	17.65 ug/L	497642	1,4-Dichlorobenzene-d4	11.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	95
2	Indane		118	C9H10	000496-11-7	94
3	Indane		118	C9H10	000496-11-7	92
4	Indane		118	C9H10	000496-11-7	87
5	Benzene, 1-ethenyl-2-methyl-		118	C9H10	000611-15-4	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU040719\
Data File : VU030840.D
Acq On : 07 Apr 2019 06:51
Operator : JC/SP
Sample : K2249-14
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 50 Sample Multiplier: 1

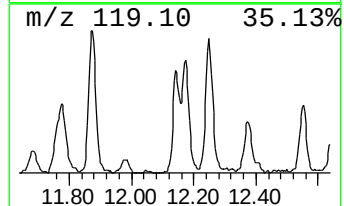
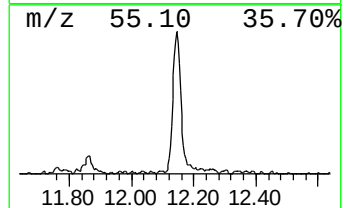
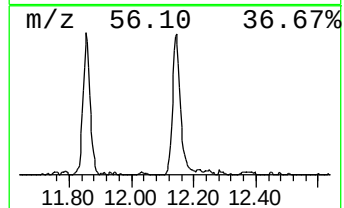
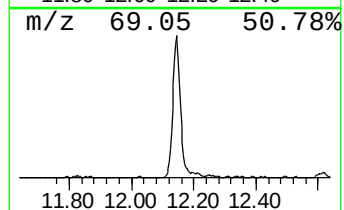
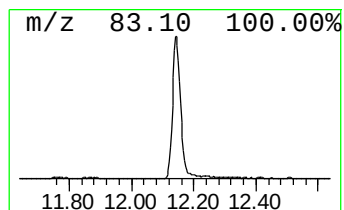
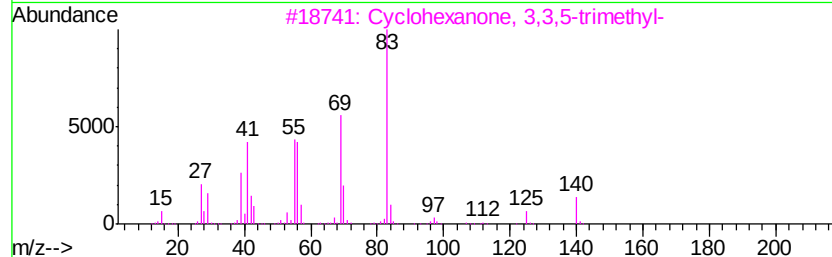
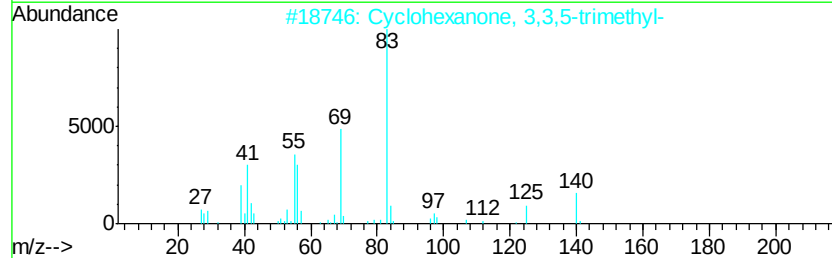
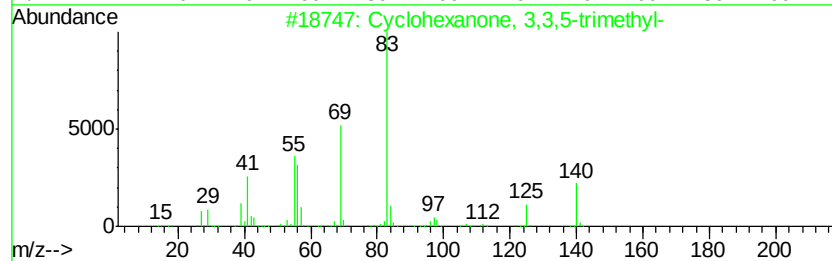
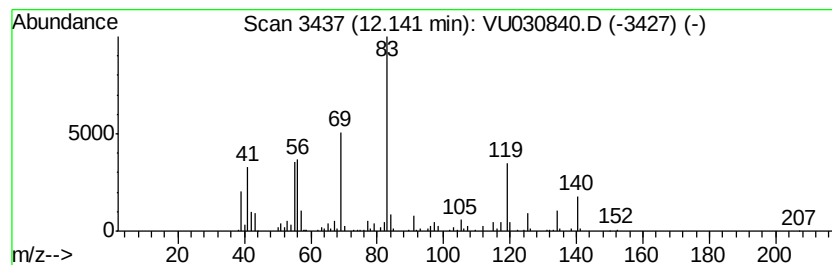
Instrument :
MSVOA_U
ClientSampleId :
E3YE6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM040519WMA.M
Quant Title : VOC Analysis

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

Peak Number 9 Cyclohexanone, 3,3,5-trimet... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	10.38 ug/L	292565	1,4-Dichlorobenzene-d4	11.48
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexanone, 3,3,5-trimethyl-	140 C9H16O	000873-94-9 76
2		Cyclohexanone, 3,3,5-trimethyl-	140 C9H16O	000873-94-9 76
3		Cyclohexanone, 3,3,5-trimethyl-	140 C9H16O	000873-94-9 68
4		1H-1,2,4-Triazole, 3-methyl-	83 C3H5N3	007170-01-6 47
5		Cyclopentanone, 2,4,4-trimethyl-	126 C8H14O	004694-12-6 47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU040719\
Data File : VU030840.D
Acq On : 07 Apr 2019 06:51
Operator : JC/SP
Sample : K2249-14
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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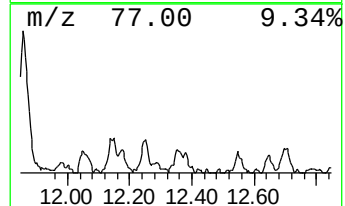
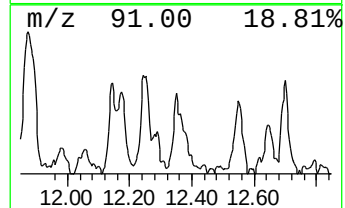
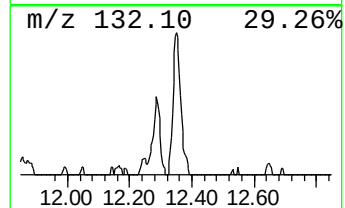
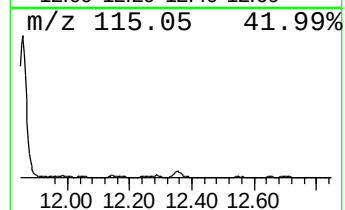
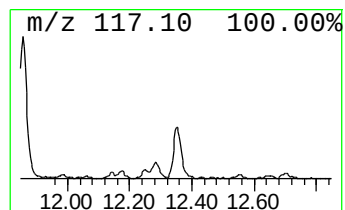
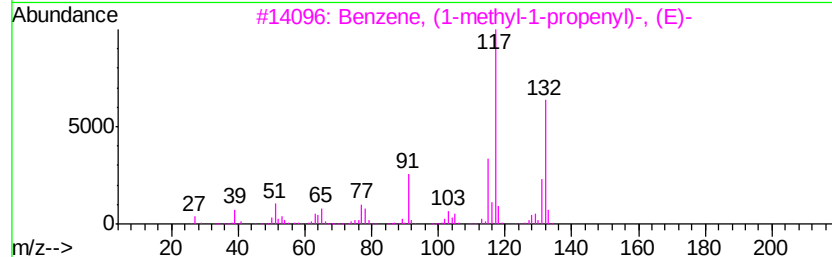
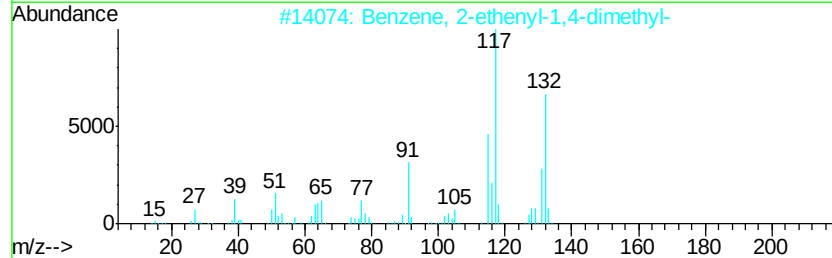
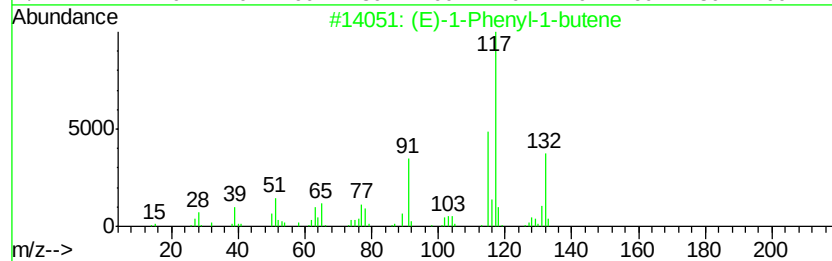
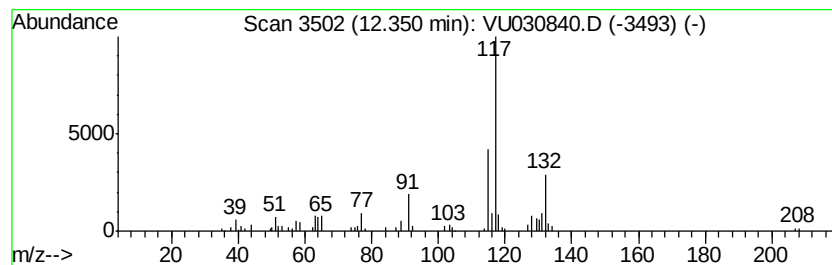
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Quant Title : VOC Analysis

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

Peak Number 10 (E)-1-Phenyl-1-butene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.35	2.76 ug/L	77907	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	94
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	91
5		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU040719\
Data File : VU030840.D
Acq On : 07 Apr 2019 06:51
Operator : JC/SP
Sample : K2249-14
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E3YE6

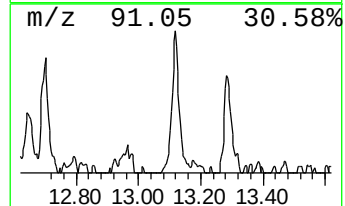
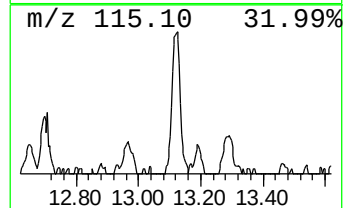
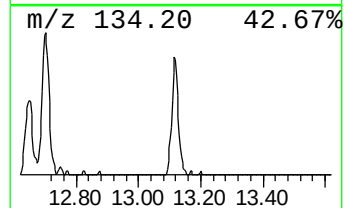
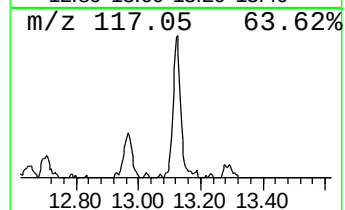
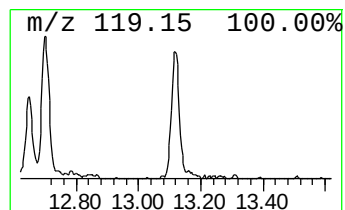
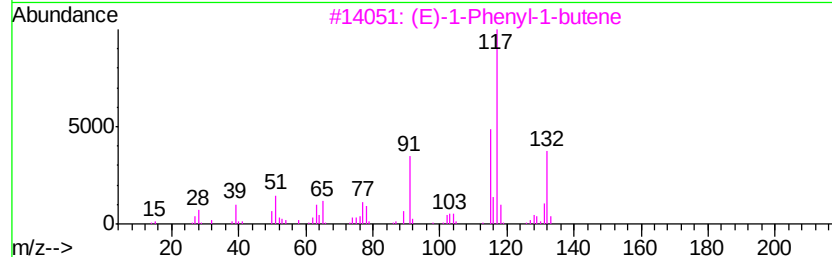
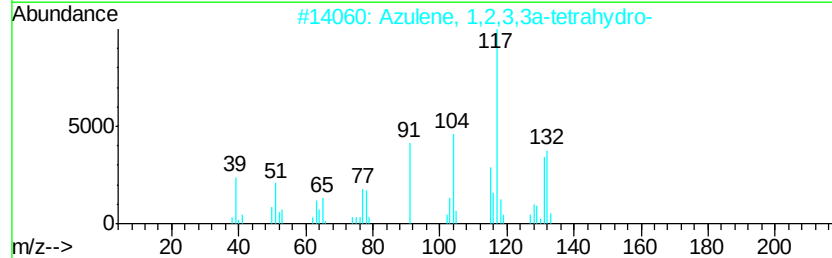
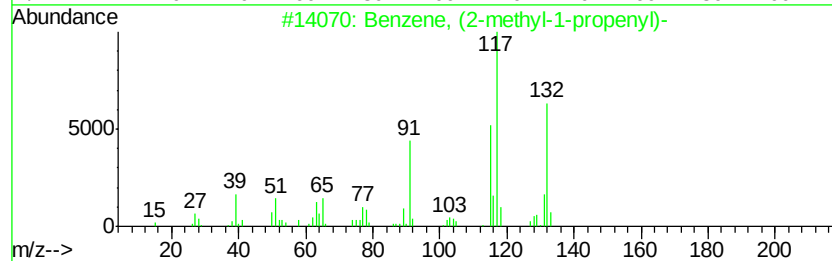
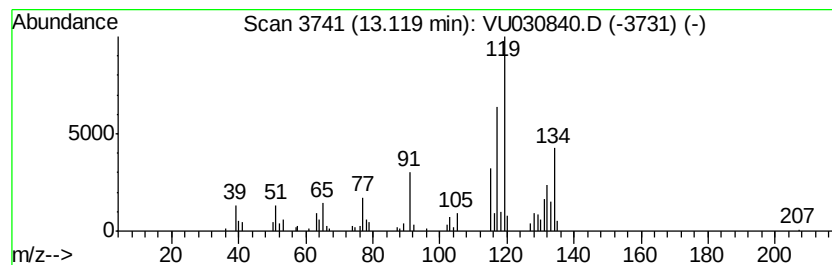
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Quant Title : VOC Analysis

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

Peak Number 11 Benzene, (2-methyl-1-propen... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	2.88 ug/L	81192	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	50
2		Azulene, 1,2,3,3a-tetrahydro-	132	C10H12	033877-87-1	50
3		(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	50
4		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	50
5		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU040719\
Data File : VU030840.D
Acq On : 07 Apr 2019 06:51
Operator : JC/SP
Sample : K2249-14
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E3YE6

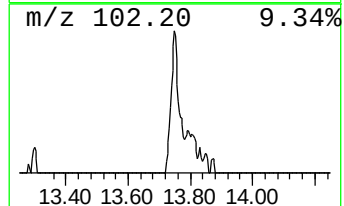
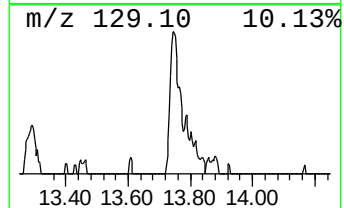
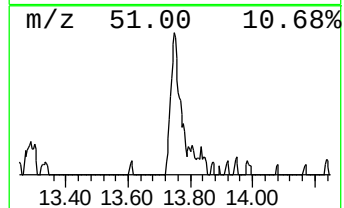
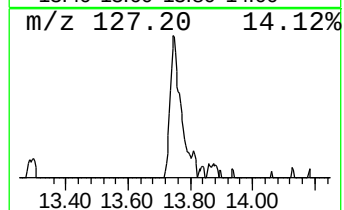
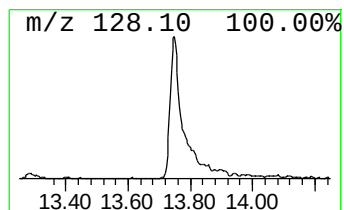
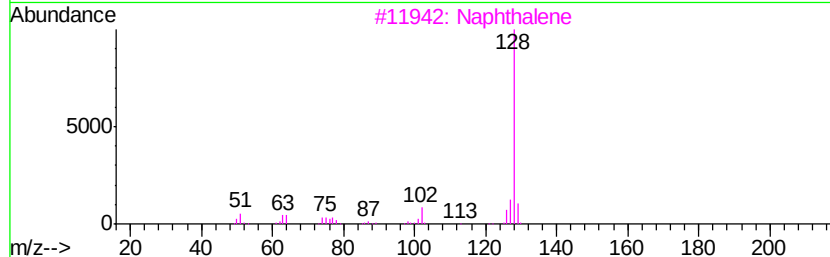
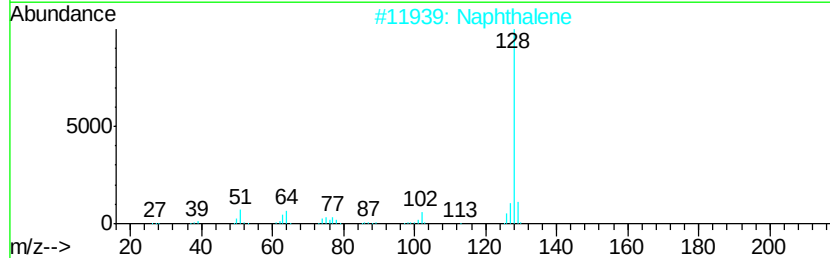
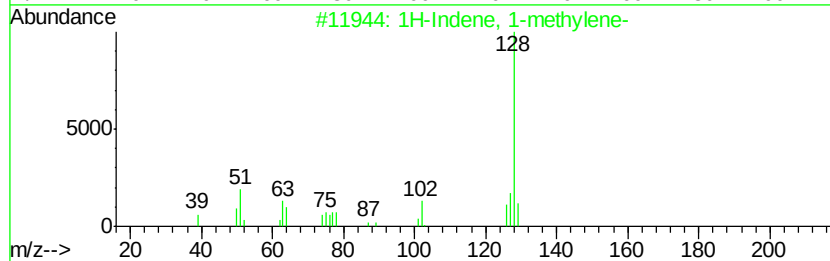
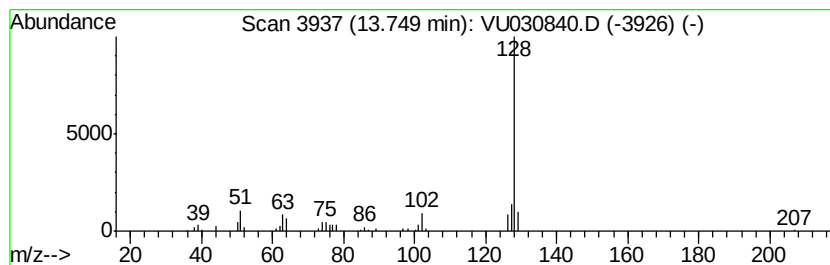
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM040519WMA.M
Quant Title : VOC Analysis

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

Peak Number 12 1H-Indene, 1-methylene- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.75	3.35 ug/L	94485	1,4-Dichlorobenzene-d4	11.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	95
2			Naphthalene	128	C10H8	000091-20-3	94
3			Naphthalene	128	C10H8	000091-20-3	94
4			Naphthalene	128	C10H8	000091-20-3	93
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU040719\
Data File : VU030840.D
Acq On : 07 Apr 2019 06:51
Operator : JC/SP
Sample : K2249-14
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E3YE6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM040519WMA.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, propyl-	10.58	9.2	ug/L	258651	3	11.48	1409920	50.0
Benzene, 1-ethyl-...	10.68	57.2	ug/L	1612680	3	11.48	1409920	50.0
Benzene, 1-ethyl-...	10.97	31.5	ug/L	887064	3	11.48	1409920	50.0
Benzene, 1,2,3-tr...	11.14	120.1	ug/L	3386220	3	11.48	1409920	50.0
Indane	11.76	17.6	ug/L	497642	3	11.48	1409920	50.0
Cyclohexanone, 3,...	12.14	10.4	ug/L	292565	3	11.48	1409920	50.0
(E)-1-Phenyl-1-bu...	12.35	2.8	ug/L	77907	3	11.48	1409920	50.0
Benzene, (2-methy...	13.12	2.9	ug/L	81192	3	11.48	1409920	50.0
1H-Indene, 1-meth...	13.75	3.4	ug/L	94485	3	11.48	1409920	50.0