

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV041819\
 Data File : VV010377.D
 Acq On : 19 Apr 2019 09:39
 Operator : SY/MD
 Sample : K2402-07
 Misc : 5.07G/10mL/MSVOA V/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 GAHH9

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_V\METHOD\SOM2VLM041819S.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	2.135	317	325	337	rVB2	17444	24174	14.60%	2.951%
2	2.238	355	357	360	rVB2	661	369	0.22%	0.045%
3	2.351	389	392	393	rBV	462	207	0.13%	0.025%
4	2.380	399	401	405	rVB	926	394	0.24%	0.048%
5	2.460	424	426	428	rBV	378	221	0.13%	0.027%
6	2.537	444	450	451	rBV3	2400	2330	1.41%	0.284%
7	2.592	464	467	471	rBV2	452	355	0.21%	0.043%
8	2.827	537	540	542	rBV2	451	333	0.20%	0.041%
9	2.952	577	579	582	rVB2	539	231	0.14%	0.028%
10	2.997	590	593	595	rVB2	623	337	0.20%	0.041%
11	3.029	600	603	606	rBV2	460	334	0.20%	0.041%
12	3.048	606	609	610	rBV	541	297	0.18%	0.036%
13	3.180	648	650	654	rBV	341	257	0.16%	0.031%
14	3.209	657	659	664	rVB	576	483	0.29%	0.059%
15	3.238	664	668	669	rBV	525	405	0.24%	0.049%
16	3.386	711	714	716	rBV	360	196	0.12%	0.024%
17	3.415	719	723	724	rBV	585	352	0.21%	0.043%
18	3.447	730	733	734	rBV2	573	294	0.18%	0.036%
19	3.505	747	751	753	rBV	560	357	0.22%	0.044%
20	3.560	765	768	772	rBV2	531	395	0.24%	0.048%
21	3.605	780	782	786	rBV3	507	443	0.27%	0.054%
22	3.939	881	886	888	rBV2	1089	1143	0.69%	0.140%
23	4.061	922	924	926	rBV2	486	258	0.16%	0.031%
24	4.183	960	962	965	rBV2	436	259	0.16%	0.032%
25	4.306	992	1000	1003	rBV2	639	784	0.47%	0.096%
26	4.341	1008	1011	1012	rVV2	492	234	0.14%	0.029%
27	4.405	1016	1031	1045	rBV2	11091	27965	16.89%	3.414%
28	4.521	1064	1067	1071	rBV	633	490	0.30%	0.060%
29	4.733	1131	1133	1137	rVB2	454	239	0.14%	0.029%
30	4.823	1158	1161	1162	rBV	501	288	0.17%	0.035%
31	4.981	1207	1210	1212	rBV	562	287	0.17%	0.035%
32	5.032	1223	1226	1230	rVB2	296	197	0.12%	0.024%
33	5.097	1230	1246	1266	rBV3	23116	65190	39.37%	7.959%
34	5.322	1312	1316	1318	rBV2	442	342	0.21%	0.042%

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Integration Parameters: LSCINT.P

Integrator: RTE
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 Title : VOC Analysis

35	5.444	1352	1354	1357	rBV	338	246	0.15%	0.030%
36	5.563	1387	1391	1392	rBV	362	248	0.15%	0.030%
37	5.611	1404	1406	1409	rVB	487	212	0.13%	0.026%
38	5.666	1413	1423	1439	rBV2	21706	43762	26.43%	5.343%
39	5.849	1478	1480	1483	rVB	549	237	0.14%	0.029%
40	5.865	1483	1485	1489	rVB2	691	332	0.20%	0.041%
41	5.884	1489	1491	1495	rVB2	683	397	0.24%	0.048%
42	5.907	1495	1498	1501	rBV	499	364	0.22%	0.044%
43	6.039	1536	1539	1541	rBV	436	207	0.13%	0.025%
44	6.052	1541	1543	1549	rVB2	498	303	0.18%	0.037%
45	6.080	1549	1552	1553	rBV2	559	363	0.22%	0.044%
46	6.119	1554	1564	1577	rVV5	13333	27992	16.91%	3.417%
47	6.341	1630	1633	1635	rBV3	399	277	0.17%	0.034%
48	6.521	1687	1689	1691	rVB	571	241	0.15%	0.029%
49	6.798	1772	1775	1777	rBV	389	282	0.17%	0.034%
50	6.920	1809	1813	1817	rBV3	616	485	0.29%	0.059%
51	7.032	1835	1848	1859	rBV8	6341	12861	7.77%	1.570%
52	7.164	1886	1889	1890	rBV2	415	195	0.12%	0.024%
53	7.232	1907	1910	1913	rBV	385	237	0.14%	0.029%
54	7.312	1932	1935	1938	rBV2	861	641	0.39%	0.078%
55	7.363	1938	1951	1969	rVV3	25171	49238	29.74%	6.011%
56	7.560	2009	2012	2013	rBV	378	229	0.14%	0.028%
57	7.630	2030	2034	2035	rBV2	460	270	0.16%	0.033%
58	7.736	2064	2067	2072	rVB	712	517	0.31%	0.063%
59	8.058	2164	2167	2174	rBV	481	470	0.28%	0.057%
60	8.238	2220	2223	2225	rBV2	349	196	0.12%	0.024%
61	8.399	2269	2273	2278	rBV4	589	587	0.35%	0.072%
62	8.460	2290	2292	2294	rBV	438	218	0.13%	0.027%
63	8.608	2336	2338	2342	rVB2	427	251	0.15%	0.031%
64	8.637	2344	2347	2350	rVB	500	275	0.17%	0.034%
65	8.839	2408	2410	2414	rVB2	497	243	0.15%	0.030%
66	8.900	2418	2429	2441	rBV2	25727	45937	27.74%	5.608%
67	9.029	2467	2469	2471	rBV	538	284	0.17%	0.035%
68	9.135	2501	2502	2505	rVV2	385	197	0.12%	0.024%
69	9.495	2610	2614	2615	rBV3	504	413	0.25%	0.050%
70	9.810	2710	2712	2718	rBV2	391	374	0.23%	0.046%
71	10.003	2770	2772	2776	rBV2	483	199	0.12%	0.024%

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72	10.260	2843	2852	2865	rVB3	10940	17288	10.44%	2.111%
73	10.360	2881	2883	2887	rBV	538	315	0.19%	0.038%
74	10.476	2915	2919	2921	rBV2	552	445	0.27%	0.054%
75	10.621	2960	2964	2968	rBV2	454	387	0.23%	0.047%
76	10.714	2989	2993	2996	rBV2	381	273	0.16%	0.033%
77	10.945	3063	3065	3067	rBV2	556	312	0.19%	0.038%
78	11.299	3165	3175	3189	rVB3	13816	23711	14.32%	2.895%
79	11.379	3197	3200	3201	rBV2	465	271	0.16%	0.033%
80	11.437	3216	3218	3221	rVB2	1056	470	0.28%	0.057%
81	11.582	3262	3263	3266	rBV2	433	237	0.14%	0.029%
82	11.675	3284	3292	3305	rVB4	17529	28089	16.96%	3.429%
83	11.981	3386	3387	3390	rBV2	637	305	0.18%	0.037%
84	12.013	3393	3397	3399	rBV2	396	286	0.17%	0.035%
85	12.026	3399	3401	3402	rBV2	524	196	0.12%	0.024%
86	12.058	3409	3411	3417	rVB4	1487	970	0.59%	0.118%
87	12.241	3461	3468	3480	rBV9	4090	6874	4.15%	0.839%
88	12.360	3503	3505	3510	rBV2	475	327	0.20%	0.040%
89	12.495	3545	3547	3548	rBV2	587	241	0.15%	0.029%
90	12.633	3586	3590	3592	rBV4	2594	2168	1.31%	0.265%
91	12.717	3615	3616	3619	rBV	323	208	0.13%	0.025%
92	12.775	3630	3634	3641	rBV4	1058	1275	0.77%	0.156%
93	12.932	3675	3683	3693	rVB7	4675	7280	4.40%	0.889%
94	12.997	3694	3703	3711	rVV3	8265	12405	7.49%	1.514%
95	13.103	3730	3736	3748	rVB6	12616	20987	12.68%	2.562%
96	13.244	3778	3780	3781	rBV6	800	383	0.23%	0.047%
97	13.418	3832	3834	3839	rBV5	1113	893	0.54%	0.109%
98	13.553	3866	3876	3889	rVB	101623	165572	100.00%	20.214%
99	14.685	4216	4228	4247	rBV	85353	145527	87.89%	17.767%
100	14.868	4277	4285	4299	rVB	37949	61188	36.96%	7.470%

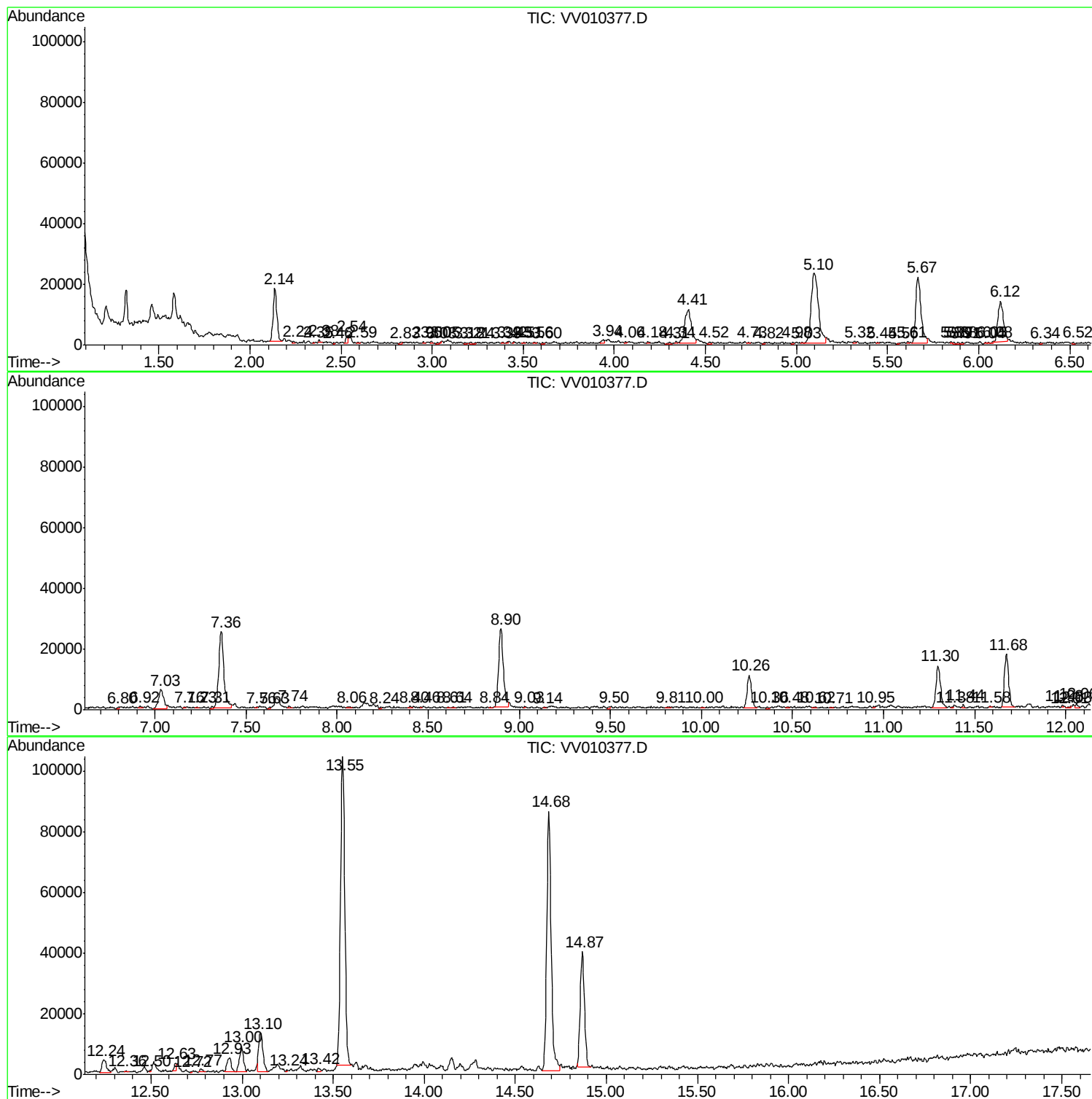
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM041819S.M
Quant Title : VOC Analysis

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



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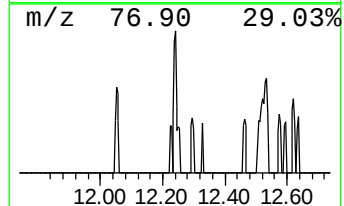
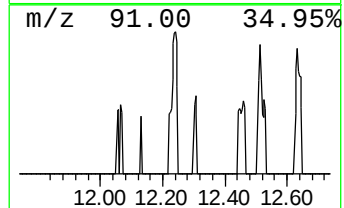
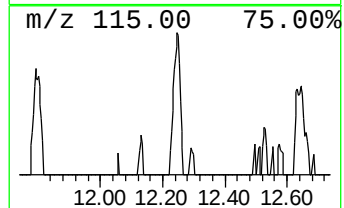
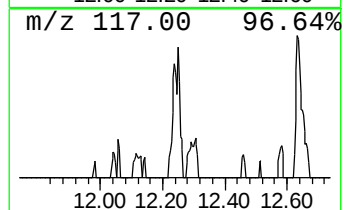
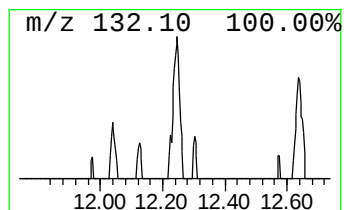
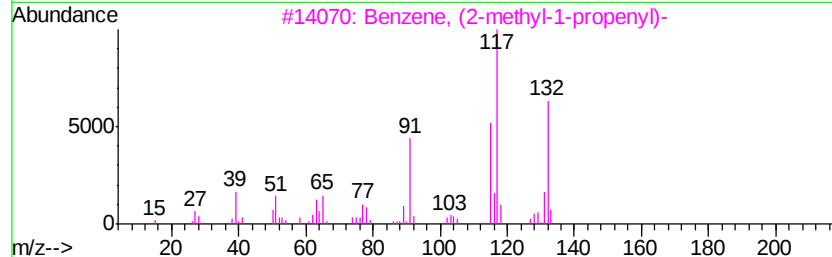
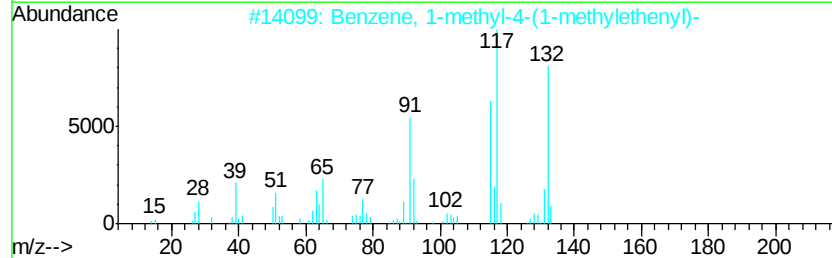
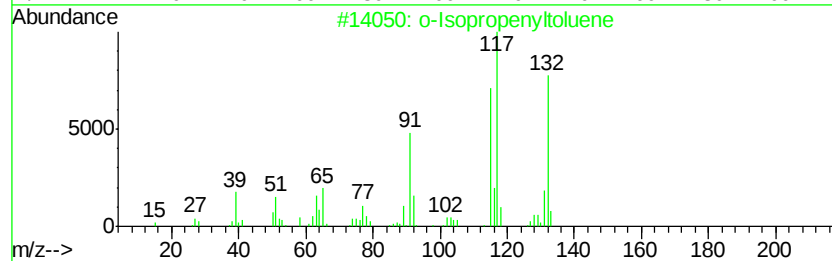
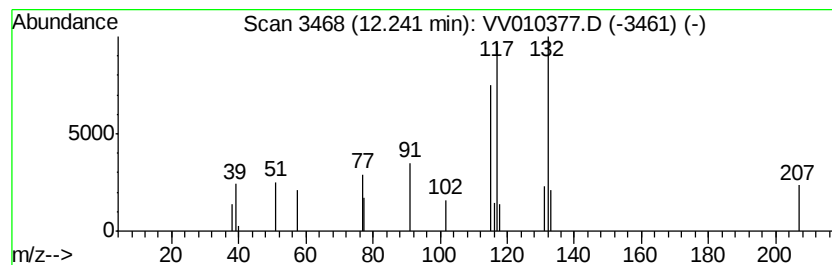
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM041819S.M
Quant Title : VOC Analysis

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

Peak Number 3 o-Isopropenyltoluene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.24	7.25 ug/L	6874	1,4-Dichlorobenzene-d4	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Isopropenyltoluene	132	C10H12	007399-49-7	64
2		Benzene, 1-methyl-4-(1-methyleth...	132	C10H12	001195-32-0	59
3		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	58
4		Benzene, 4-ethenyl-1,2-dimethyl-	132	C10H12	027831-13-6	58
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	53



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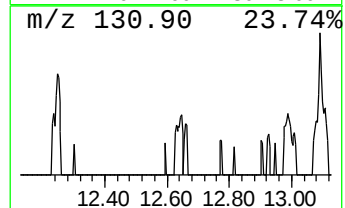
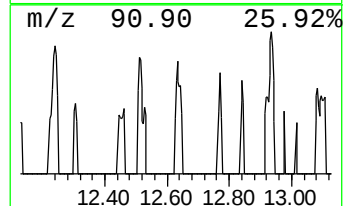
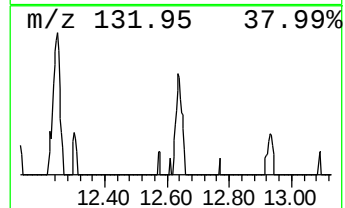
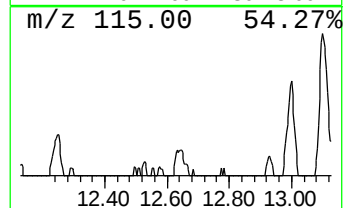
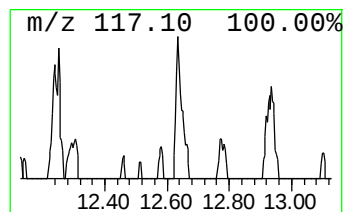
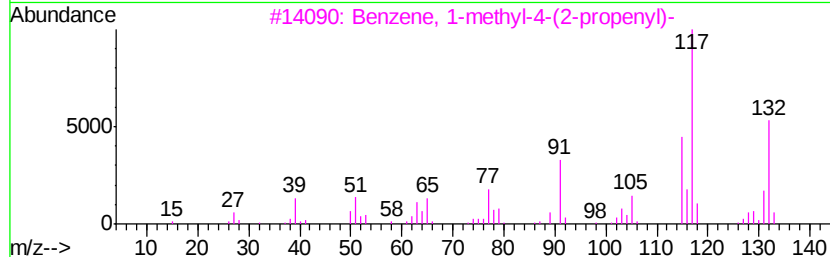
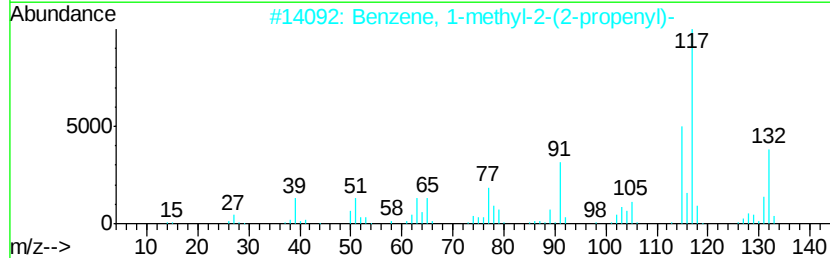
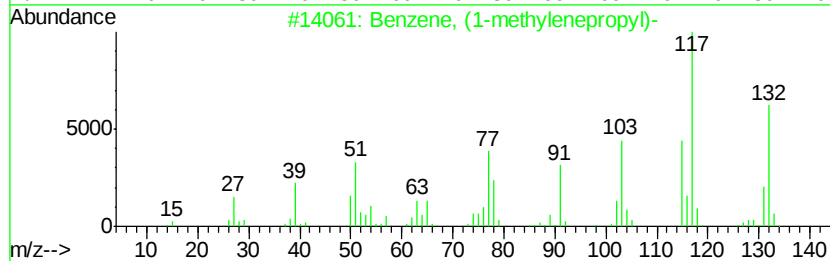
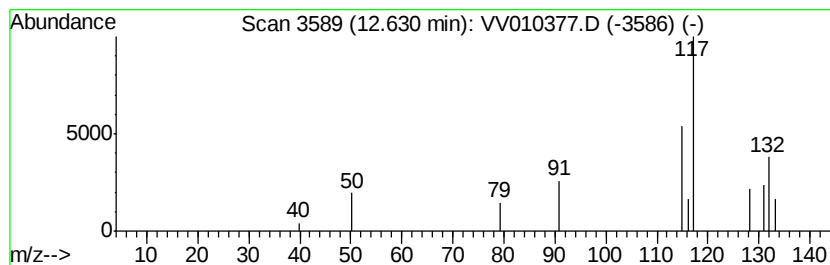
Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM041819S.M
Quant Title : VOC Analysis

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

Peak Number 4 Benzene, (1-methylenepropyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.63	2.29 ug/L	2168	1,4-Dichlorobenzene-d4	11.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1-methylenepropyl)-	132 C10H12	002039-93-2 86
2		Benzene, 1-methyl-2-(2-propenyl)-	132 C10H12	001587-04-8 86
3		Benzene, 1-methyl-4-(2-propenyl)-	132 C10H12	003333-13-9 86
4		Benzene, (2-methyl-2-propenyl)-	132 C10H12	003290-53-7 78
5		(E)-1-Phenyl-1-butene	132 C10H12	001005-64-7 78



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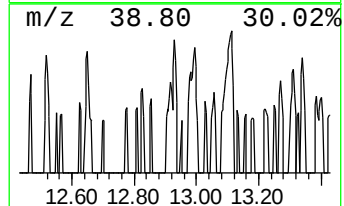
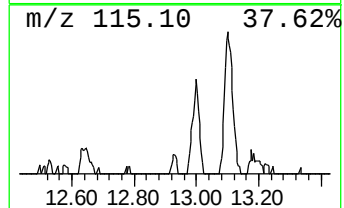
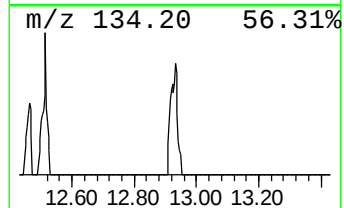
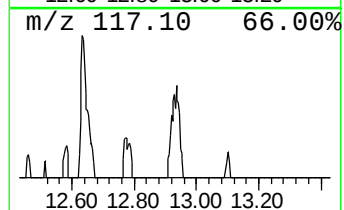
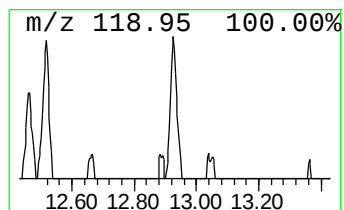
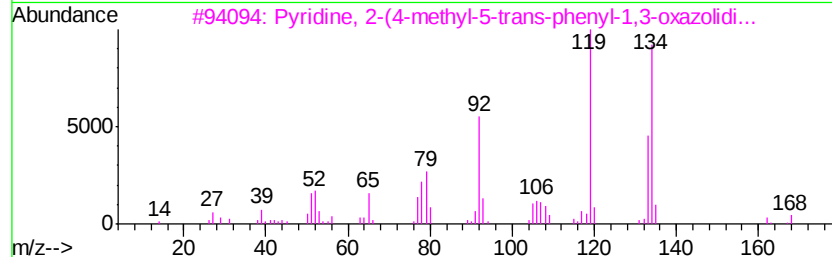
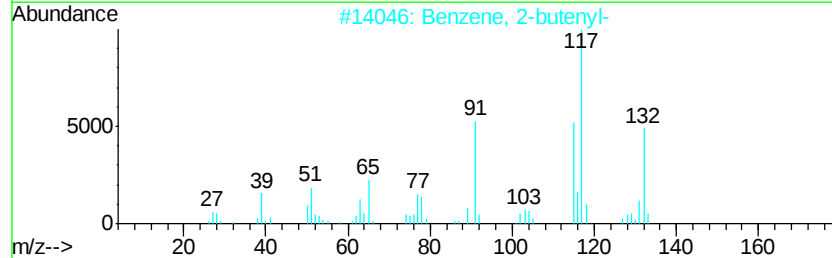
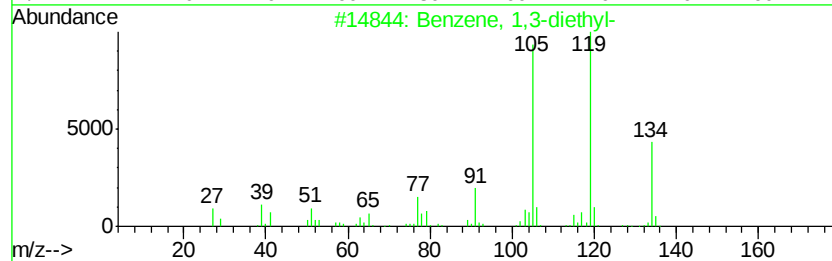
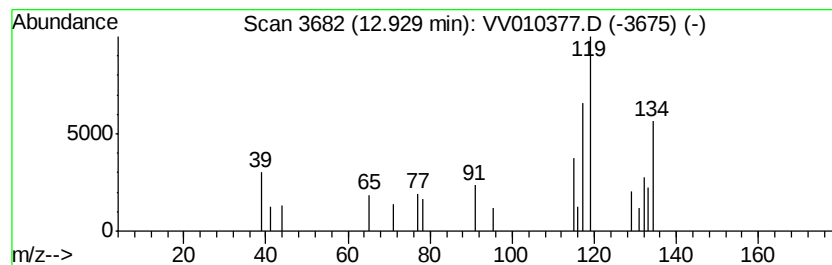
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Quant Title : VOC Analysis

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

Peak Number 6 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.93	7.68 ug/L	7280	1,4-Dichlorobenzene-d4	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	32
2		Benzene, 2-butenyl-	132	C10H12	001560-06-1	25
3		Pvridine, 2-(4-methyl-5-trans-ph...	240	C15H16N2O	1000142-75-0	25
4		Cyclopropa[1,2:1,3]dicyclopenten...	176	C12H16O	091531-53-2	25
5		2-(1-Hydroxyethyl)hydroxymethylb...	152	C9H12O2	057259-71-9	23



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV041819\
Data File : VV010377.D
Acq On : 19 Apr 2019 09:39
Operator : SY/MD
Sample : K2402-07
Misc : 5.07G/10mL/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

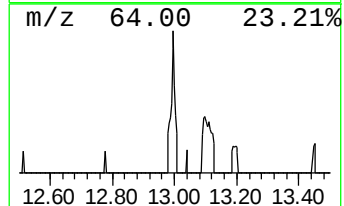
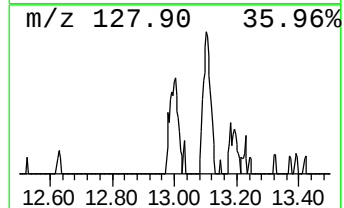
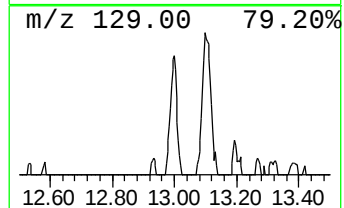
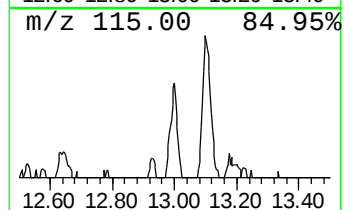
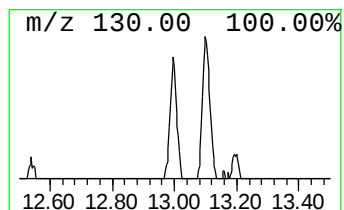
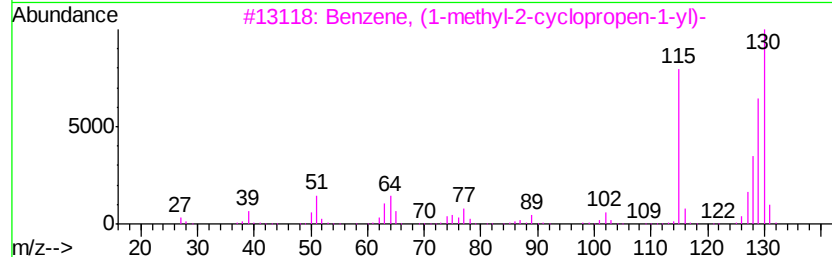
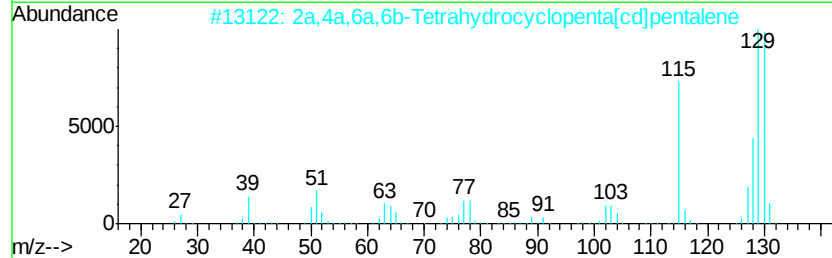
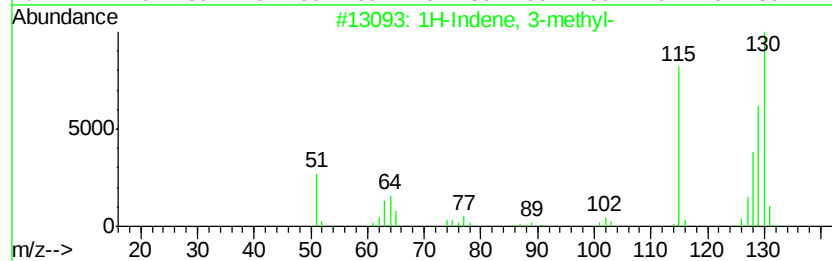
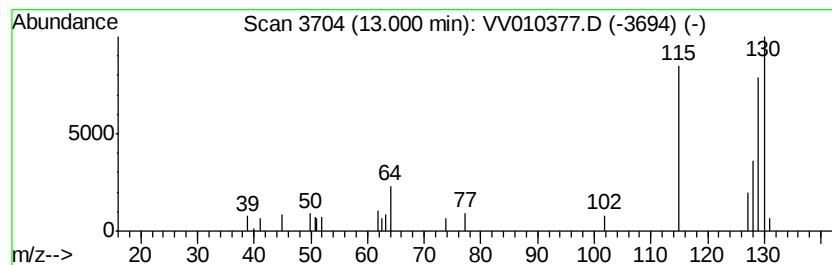
Instrument :
MSVOA_V
ClientSampleId :
GAHH9

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM041819S.M
Quant Title : VOC Analysis

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

Peak Number 7 1H-Indene, 3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.00	13.08 ug/L	12405	1,4-Dichlorobenzene-d4	11.30	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 3-methyl-	130	C10H10	000767-60-2	80
2	2a,4a,6a,6b-Tetrahydrocyclopenta...	130	C10H10	006053-74-3	80
3	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	80
4	2-Methylindene	130	C10H10	002177-47-1	80
5	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV041819\
Data File : VV010377.D
Acq On : 19 Apr 2019 09:39
Operator : SY/MD
Sample : K2402-07
Misc : 5.07G/10mL/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

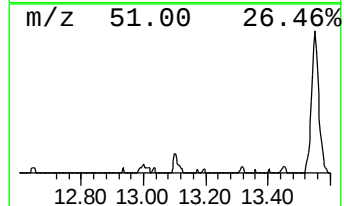
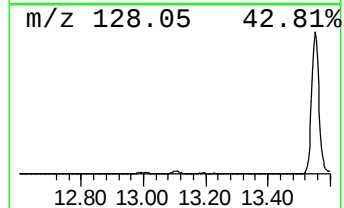
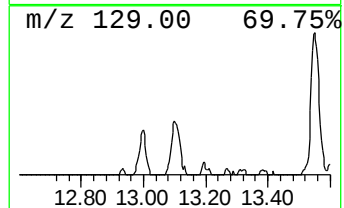
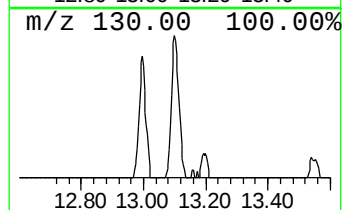
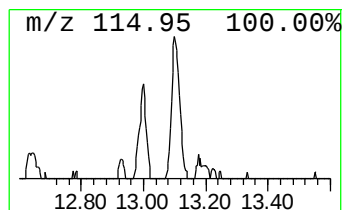
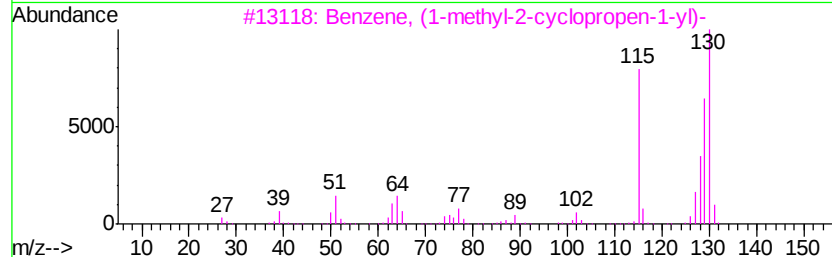
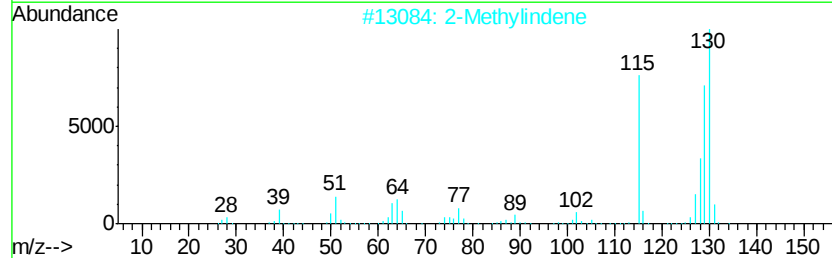
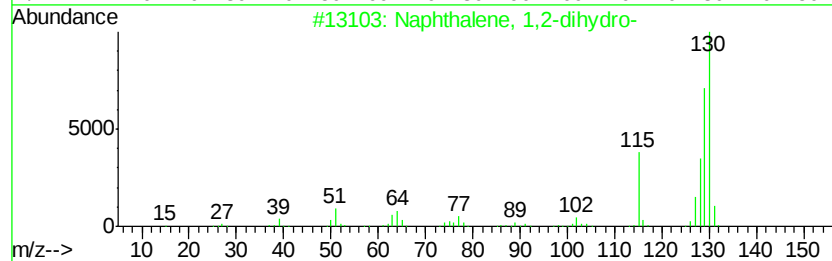
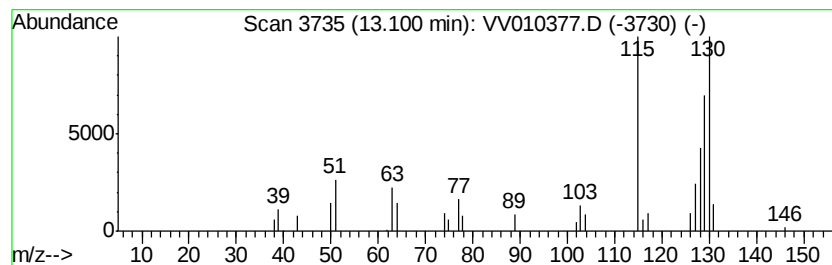
Instrument :
MSVOA_V
ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM041819S.M
Quant Title : VOC Analysis

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

Peak Number 8 Naphthalene, 1,2-dihydro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.10	22.13 ug/L	20987	1,4-Dichlorobenzene-d4	11.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2-dihydro-	130 C10H10	000447-53-0 68
2		2-Methylindene	130 C10H10	002177-47-1 64
3		Benzene, (1-methyl-2-cyclopropen...	130 C10H10	065051-83-4 64
4		1H-Indene, 1-methyl-	130 C10H10	000767-59-9 64
5		1H-Indene, 3-methyl-	130 C10H10	000767-60-2 59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV041819\
Data File : VV010377.D
Acq On : 19 Apr 2019 09:39
Operator : SY/MD
Sample : K2402-07
Misc : 5.07G/10mL/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

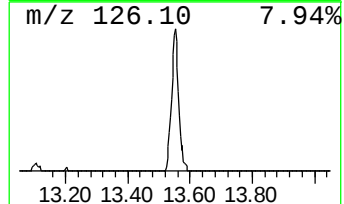
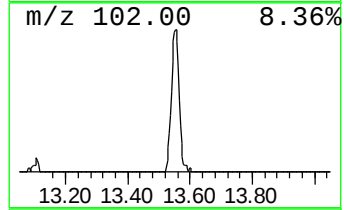
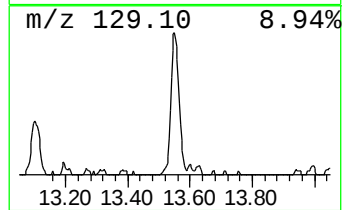
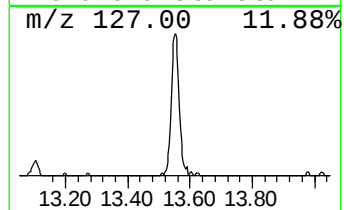
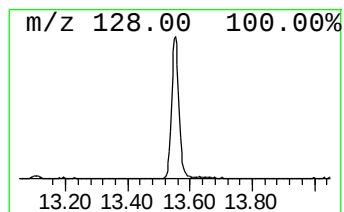
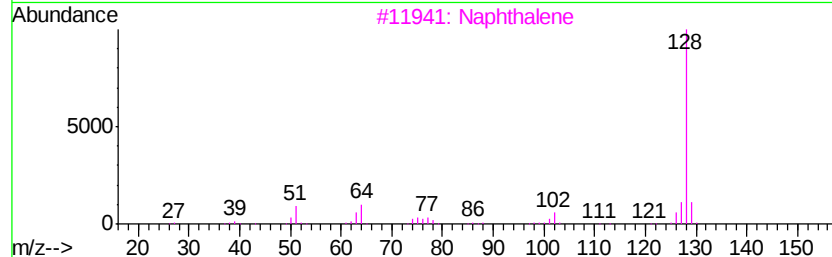
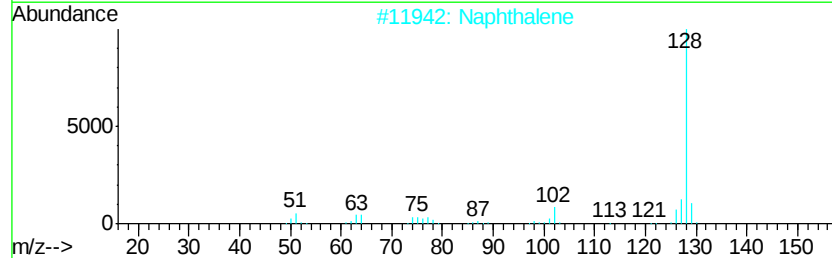
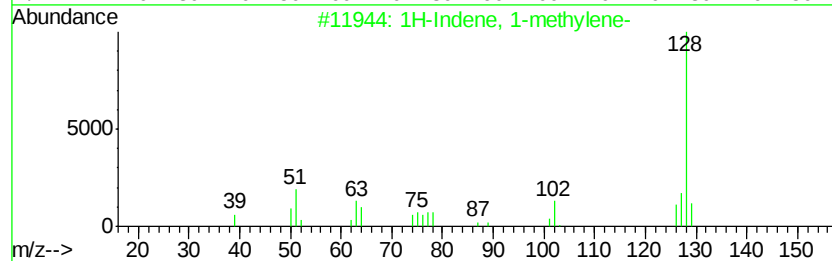
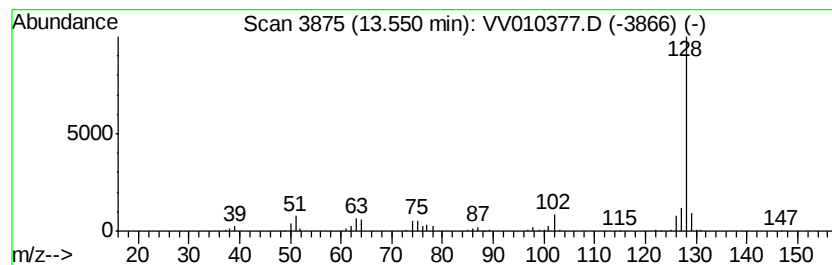
Instrument :
MSVOA_V
ClientSampleId :
GAHH9

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM041819S.M
Quant Title : VOC Analysis

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

Peak Number 9 1H-Indene, 1-methylene- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.55	174.57 ug/L	165572	1,4-Dichlorobenzene-d4	11.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Indene, 1-methylene-	128 C10H8	002471-84-3 94
2		Naphthalene	128 C10H8	000091-20-3 93
3		Naphthalene	128 C10H8	000091-20-3 90
4		Naphthalene	128 C10H8	000091-20-3 90
5		Azulene	128 C10H8	000275-51-4 90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV041819\
Data File : VV010377.D
Acq On : 19 Apr 2019 09:39
Operator : SY/MD
Sample : K2402-07
Misc : 5.07G/10mL/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAHH9

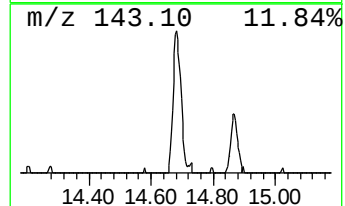
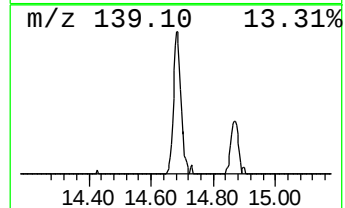
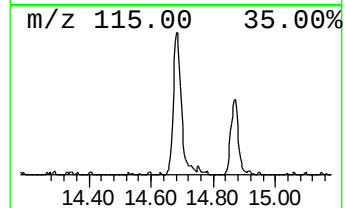
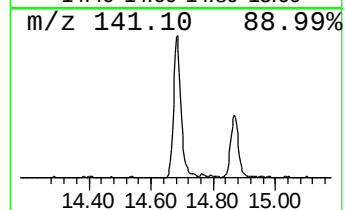
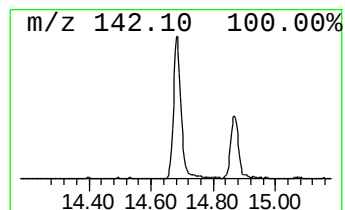
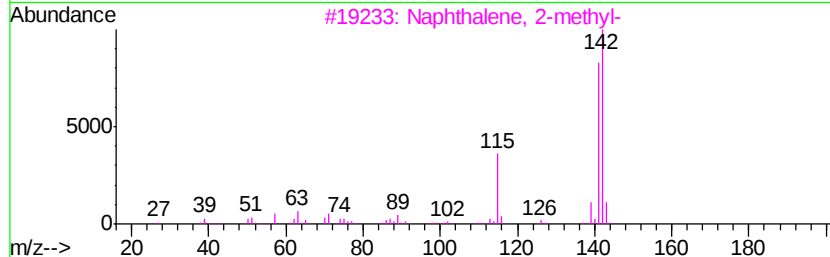
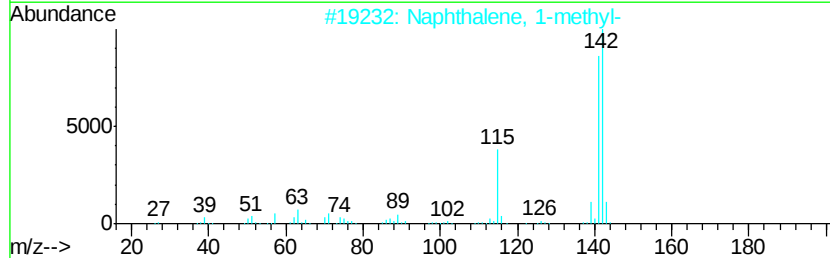
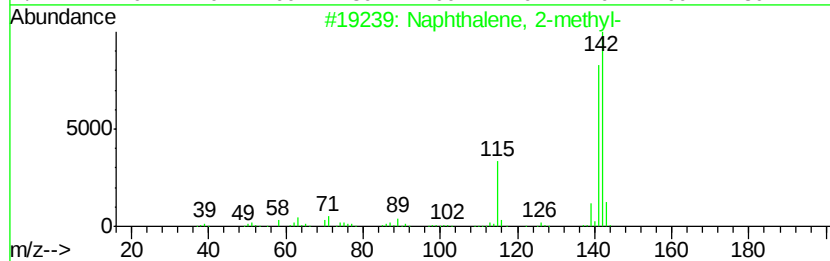
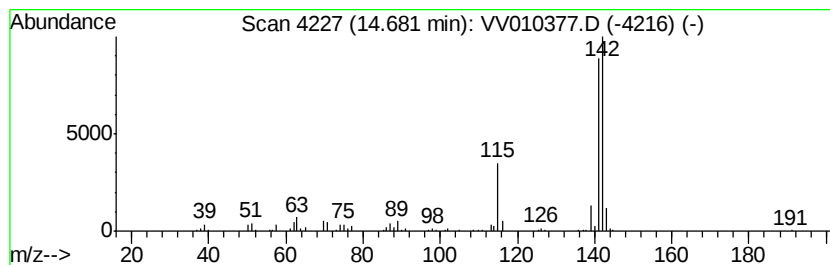
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Quant Title : VOC Analysis

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.68	153.44 ug/L	145527	1,4-Dichlorobenzene-d4	11.30

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV041819\
Data File : VV010377.D
Acq On : 19 Apr 2019 09:39
Operator : SY/MD
Sample : K2402-07
Misc : 5.07G/10mL/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAHH9

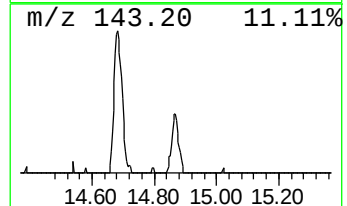
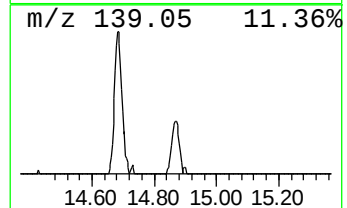
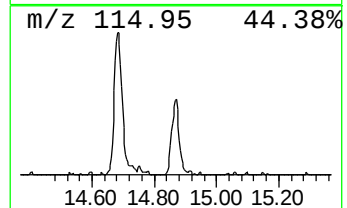
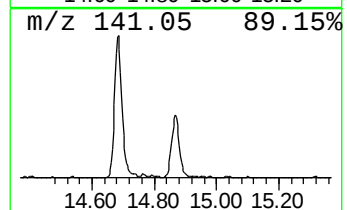
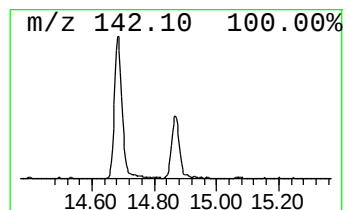
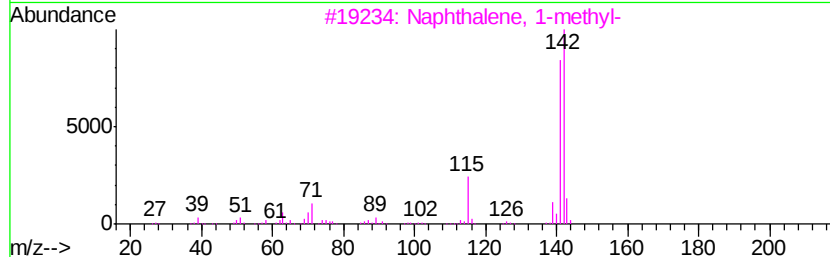
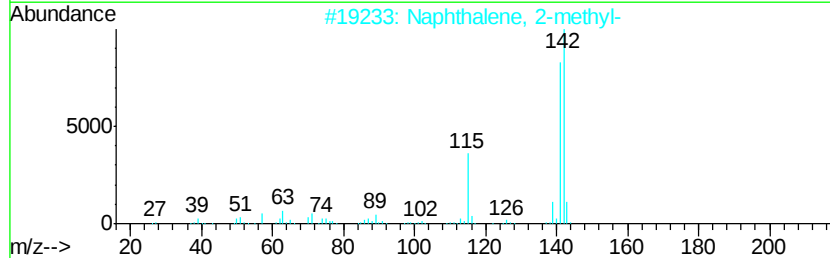
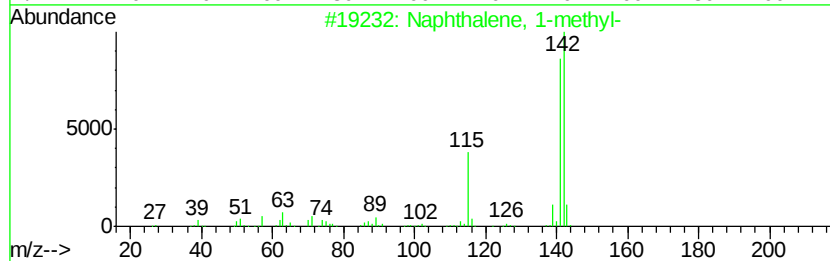
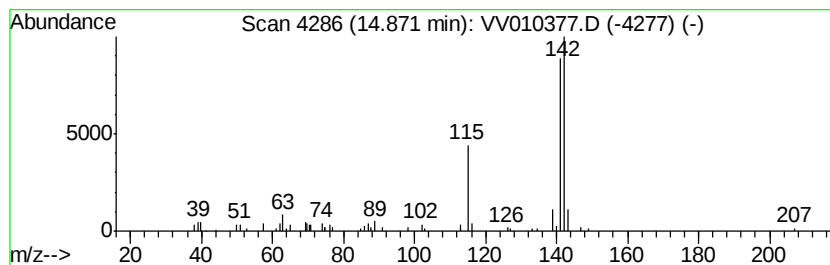
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Quant Title : VOC Analysis

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

Peak Number 11 Naphthalene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	64.51 ug/L	61188	1,4-Dichlorobenzene-d4	11.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
3			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV041819\
Data File : VV010377.D
Acq On : 19 Apr 2019 09:39
Operator : SY/MD
Sample : K2402-07
Misc : 5.07G/10mL/MSVOA_V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAHH9

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM041819S.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
o-Isopropenyltoluene	12.24	7.3	ug/L	6874	3	11.30	23711	25.0
Benzene, (1-methy...	12.63	2.3	ug/L	2168	3	11.30	23711	25.0
unknown-01	12.93	7.7	ug/L	7280	3	11.30	23711	25.0
1H-Indene, 3-methyl-	13.00	13.1	ug/L	12405	3	11.30	23711	25.0
Naphthalene, 1,2-...	13.10	22.1	ug/L	20987	3	11.30	23711	25.0
1H-Indene, 1-meth...	13.55	174.6	ug/L	165572	3	11.30	23711	25.0
Naphthalene, 1-me...	14.68	153.4	ug/L	145527	3	11.30	23711	25.0
Naphthalene, 2-me...	14.87	64.5	ug/L	61188	3	11.30	23711	25.0