

Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
 Data File : VV010646.D
 Acq On : 03 May 2019 03:51
 Operator : SY/MD
 Sample : K2603-20
 Misc : 5.00G/10ML/MSVOA V/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 GAJ72

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\SOM2VLM050219S.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	1.309	62	68	76	rVB	63801	63279	0.37%	0.073%
2	1.573	143	150	160	rVB	50211	59613	0.35%	0.068%
3	2.036	291	294	301	rVB4	1103	1427	0.01%	0.002%
4	2.119	307	320	331	rBV	124811	181962	1.06%	0.209%
5	2.235	354	356	361	rVB2	495	353	0.00%	0.000%
6	2.306	371	378	388	rBV4	2300	3537	0.02%	0.004%
7	2.531	438	448	459	rBV2	4394	6968	0.04%	0.008%
8	2.672	489	492	496	rBV2	223	155	0.00%	0.000%
9	2.798	526	531	534	rBV	413	368	0.00%	0.000%
10	2.852	545	548	550	rBV2	229	156	0.00%	0.000%
11	3.023	598	601	603	rBV2	239	183	0.00%	0.000%
12	3.068	608	615	627	rVB5	1461	2775	0.02%	0.003%
13	3.261	671	675	679	rBV2	298	262	0.00%	0.000%
14	3.524	752	757	761	rBV2	217	218	0.00%	0.000%
15	3.566	767	770	773	rVV2	280	250	0.00%	0.000%
16	3.682	803	806	810	rVB2	273	223	0.00%	0.000%
17	3.701	810	812	815	rBV2	202	152	0.00%	0.000%
18	3.926	871	882	902	rBV	14003	36075	0.21%	0.041%
19	4.164	953	956	961	rBV3	273	255	0.00%	0.000%
20	4.280	988	992	998	rVB2	251	255	0.00%	0.000%
21	4.393	1009	1027	1053	rBV2	104968	266208	1.55%	0.306%
22	4.823	1157	1161	1164	rVB2	192	199	0.00%	0.000%
23	4.865	1170	1174	1177	rBV2	234	168	0.00%	0.000%
24	4.949	1196	1200	1203	rBV2	192	175	0.00%	0.000%
25	5.007	1215	1218	1220	rBV2	273	220	0.00%	0.000%
26	5.090	1224	1244	1270	rBV2	244640	609002	3.54%	0.699%
27	5.402	1335	1341	1343	rBV2	227	244	0.00%	0.000%
28	5.659	1407	1421	1438	rBV	213829	422876	2.46%	0.486%
29	5.907	1494	1498	1501	rBV2	256	247	0.00%	0.000%
30	5.946	1501	1510	1525	rVB5	3083	6486	0.04%	0.007%
31	6.000	1525	1527	1530	rBV2	303	232	0.00%	0.000%
32	6.013	1530	1531	1537	rVB3	399	241	0.00%	0.000%
33	6.113	1548	1562	1588	rBV	143727	288593	1.68%	0.331%
34	6.318	1623	1626	1629	rBV2	207	157	0.00%	0.000%

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Integrator: RTE
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Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM050219S.M
 Title : VOC Analysis

35	6.363	1636	1640	1642	rBV2	300	260	0.00%	0.000%
36	6.495	1676	1681	1692	rVB2	315	522	0.00%	0.001%
37	6.544	1692	1696	1700	rBV2	280	260	0.00%	0.000%
38	6.585	1707	1709	1715	rBV	152	159	0.00%	0.000%
39	6.611	1715	1717	1721	rBV2	261	229	0.00%	0.000%
40	6.675	1732	1737	1739	rBV2	323	357	0.00%	0.000%
41	6.720	1748	1751	1753	rBV	232	152	0.00%	0.000%
42	6.820	1774	1782	1787	rVB6	1004	1357	0.01%	0.002%
43	6.888	1800	1803	1807	rVB2	224	169	0.00%	0.000%
44	6.910	1807	1810	1812	rBV2	226	157	0.00%	0.000%
45	7.026	1834	1846	1865	rBV	67921	119581	0.69%	0.137%
46	7.158	1884	1887	1889	rBV	267	153	0.00%	0.000%
47	7.219	1902	1906	1908	rVB2	324	175	0.00%	0.000%
48	7.357	1937	1949	1966	rBV	275500	488435	2.84%	0.561%
49	7.659	2033	2043	2059	rBV	38889	65685	0.38%	0.075%
50	7.759	2071	2074	2077	rBV2	287	189	0.00%	0.000%
51	8.058	2162	2167	2172	rBV2	345	376	0.00%	0.000%
52	8.132	2179	2190	2204	rBV	48524	87389	0.51%	0.100%
53	8.531	2310	2314	2320	rBV2	379	439	0.00%	0.001%
54	8.669	2351	2357	2360	rVB2	327	299	0.00%	0.000%
55	8.701	2360	2367	2370	rBV2	346	422	0.00%	0.000%
56	8.810	2394	2401	2403	rBV2	322	428	0.00%	0.000%
57	8.894	2414	2427	2465	rVB	286070	491917	2.86%	0.565%
58	9.058	2470	2478	2487	rVV3	2123	3016	0.02%	0.003%
59	9.180	2507	2516	2530	rVB	13385	21001	0.12%	0.024%
60	9.498	2611	2615	2616	rBV	261	174	0.00%	0.000%
61	9.598	2633	2646	2669	rVB4	23086	55362	0.32%	0.064%
62	9.920	2743	2746	2751	rBV2	301	275	0.00%	0.000%
63	9.945	2751	2754	2758	rVV	201	156	0.00%	0.000%
64	10.260	2840	2852	2866	rBV	94745	148256	0.86%	0.170%
65	10.492	2914	2924	2940	rBV2	25014	53218	0.31%	0.061%
66	10.582	2942	2952	2976	rVB	85702	131890	0.77%	0.151%
67	10.781	3005	3014	3019	rBV	27184	43333	0.25%	0.050%
68	10.958	3054	3069	3080	rBV	124932	191203	1.11%	0.220%
69	11.029	3080	3091	3099	rVV	592397	874765	5.08%	1.004%
70	11.077	3100	3106	3120	rVB	300199	438513	2.55%	0.503%
71	11.293	3163	3173	3189	rBV	252811	398311	2.31%	0.457%

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 Title : VOC Analysis

72	11.379	3190	3200	3211	rVV	381588	569962	3.31%	0.654%
73	11.440	3212	3219	3234	rVB	56110	84172	0.49%	0.097%
74	11.576	3248	3261	3270	rBV	247717	408185	2.37%	0.469%
75	11.672	3282	3291	3307	rVB2	273598	526721	3.06%	0.605%
76	11.743	3307	3313	3316	rBV	12583	14598	0.08%	0.017%
77	11.791	3317	3328	3344	rVB	3105549	4711082	27.37%	5.409%
78	11.984	3370	3388	3396	rBV4	101569	251002	1.46%	0.288%
79	12.035	3396	3404	3408	rBV	140338	203566	1.18%	0.234%
80	12.238	3459	3467	3478	rVB	873882	1283132	7.45%	1.473%
81	12.460	3527	3536	3543	rBV	250613	352535	2.05%	0.405%
82	12.514	3543	3553	3565	rBV2	711899	1535186	8.92%	1.763%
83	12.637	3582	3591	3601	rBV	722325	1188224	6.90%	1.364%
84	12.772	3626	3633	3648	rVB	237119	341526	1.98%	0.392%
85	12.926	3673	3681	3692	rVV2	1132048	1690222	9.82%	1.941%
86	12.993	3693	3702	3713	rVB	1521419	2237212	13.00%	2.569%
87	13.103	3724	3736	3751	rVB	2913541	4759993	27.65%	5.465%
88	13.193	3755	3764	3779	rVB4	335436	642528	3.73%	0.738%
89	13.556	3865	3877	3888	rBV	10642735	17213170	100.00%	19.763%
90	14.148	4051	4061	4070	rBV4	357465	661900	3.85%	0.760%
91	14.685	4217	4228	4243	rVB	7902050	12311705	71.52%	14.136%
92	14.871	4274	4286	4300	rBV	9411760	15004980	87.17%	17.228%
93	15.717	4538	4549	4572	rBV	2148698	3802086	22.09%	4.365%
94	15.865	4585	4595	4601	rBV	3256923	4953991	28.78%	5.688%
95	15.993	4627	4635	4645	rVB	543494	811395	4.71%	0.932%
96	16.061	4647	4656	4660	rBV	584458	879479	5.11%	1.010%
97	16.234	4701	4710	4726	rBV	344610	563211	3.27%	0.647%
98	16.325	4728	4738	4753	rVB	2289321	3592809	20.87%	4.125%
99	16.537	4796	4804	4820	rVB2	369928	618978	3.60%	0.711%
100	17.447	5077	5087	5105	rBV2	167827	311473	1.81%	0.358%

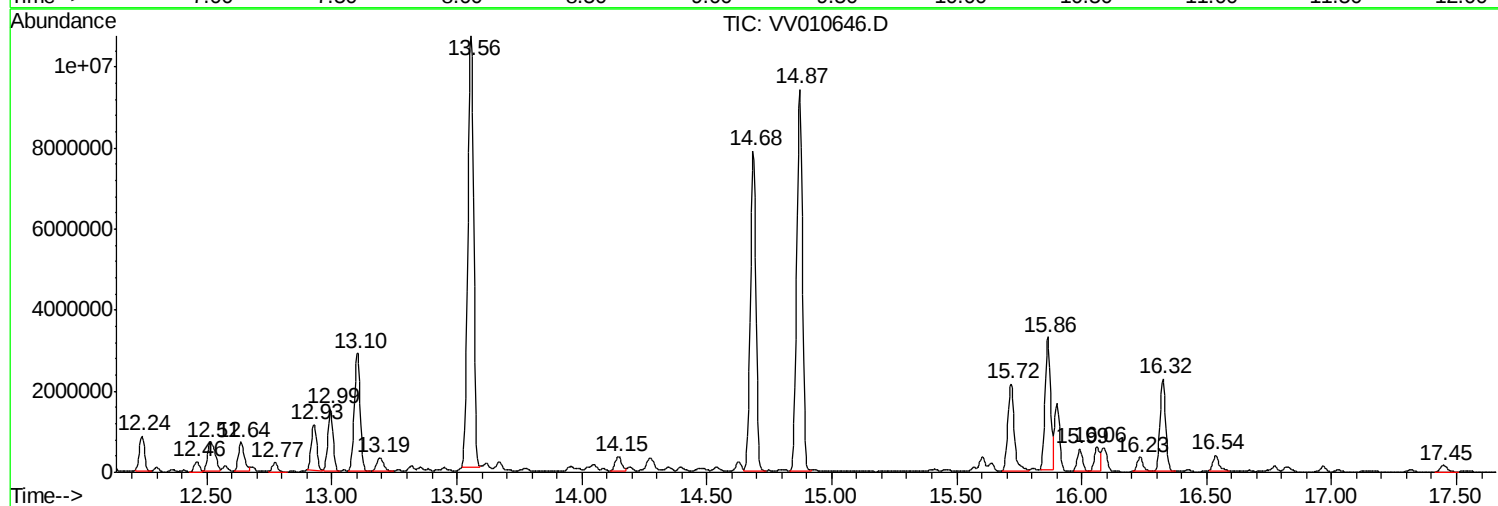
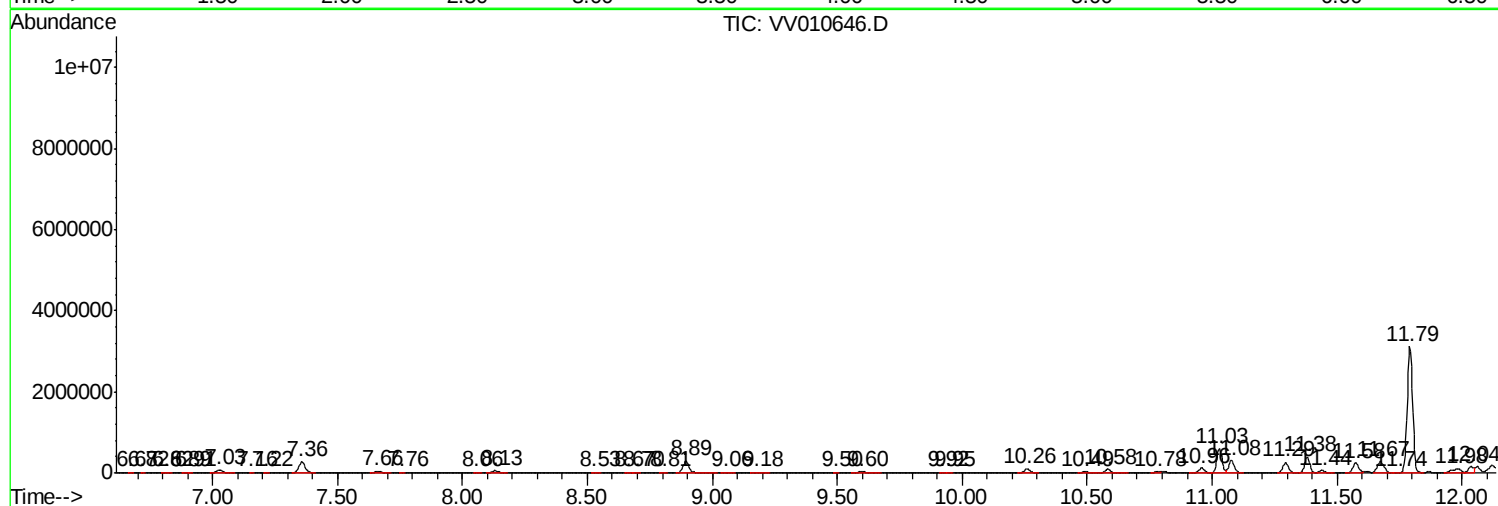
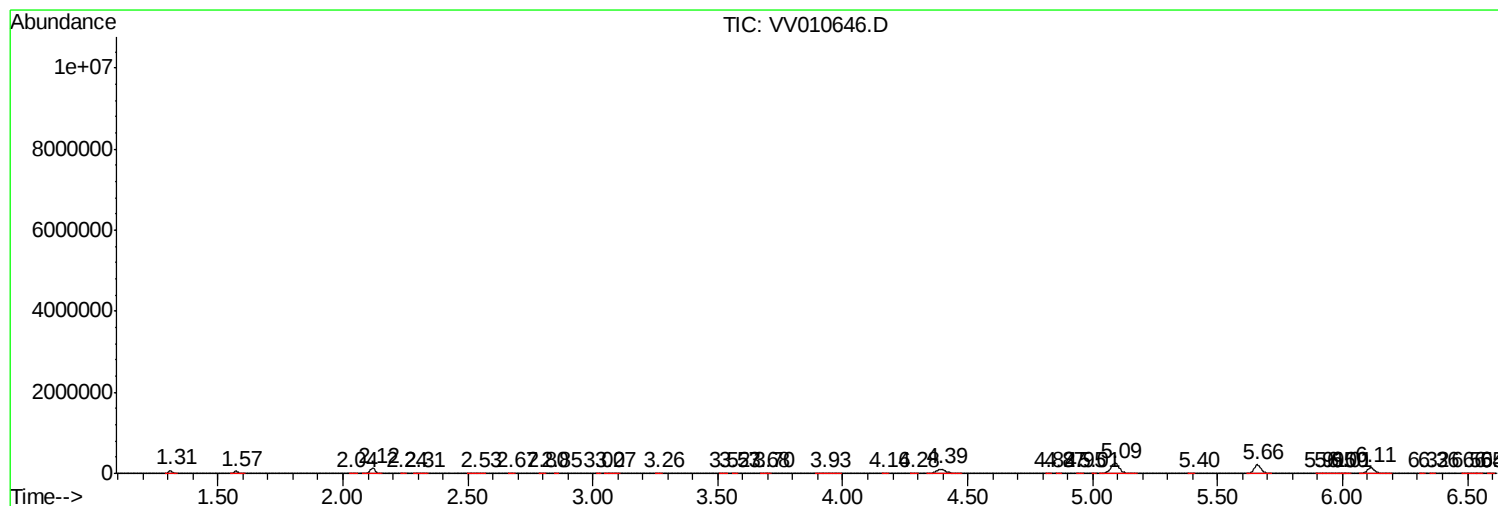
Sum of corrected areas: 87097120

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Instrument :
MSVOA_V
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM050219S.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
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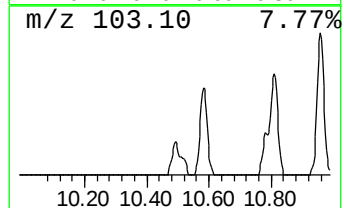
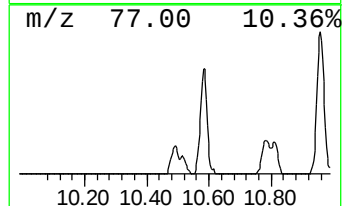
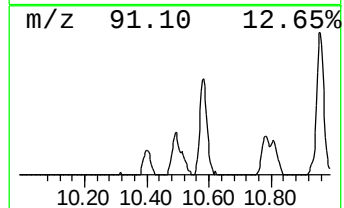
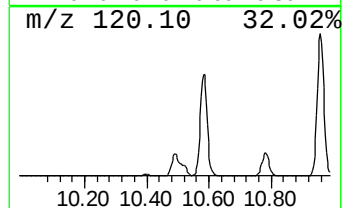
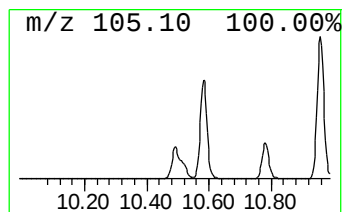
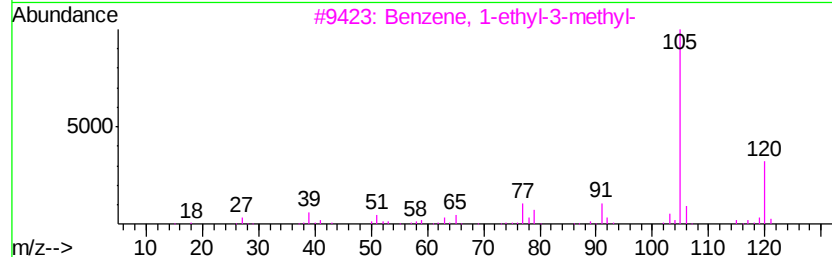
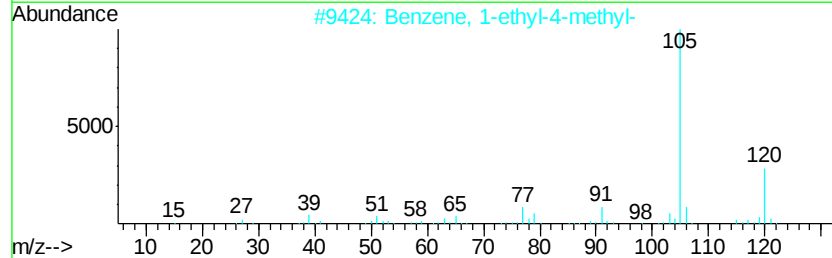
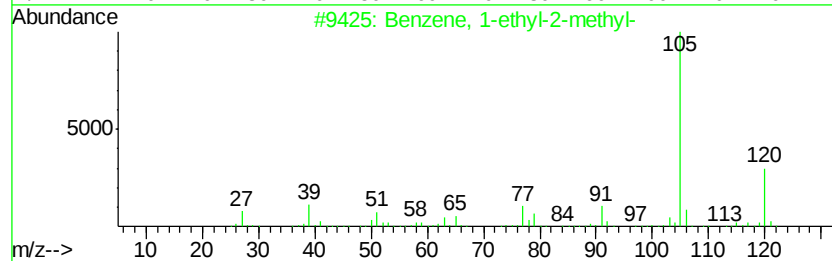
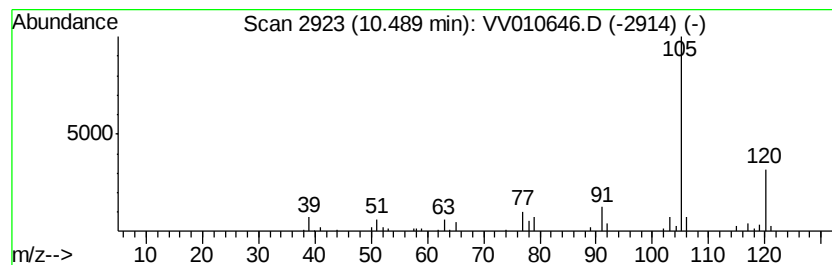
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM050219S.M
Quant Title : VOC Analysis

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

Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.49	3.34 ug/L	53218	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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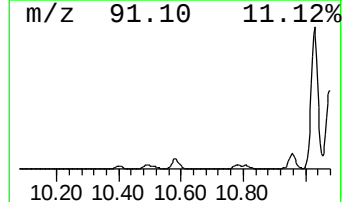
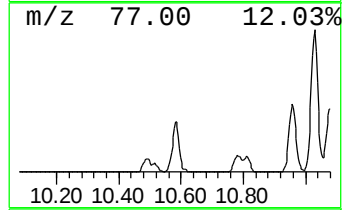
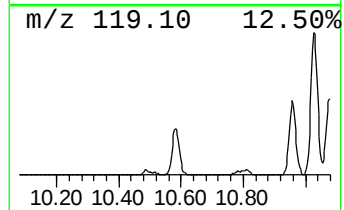
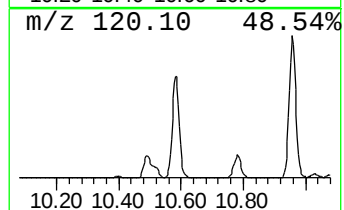
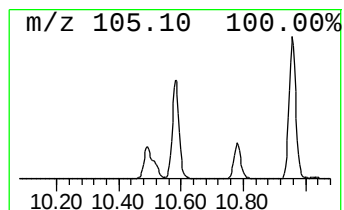
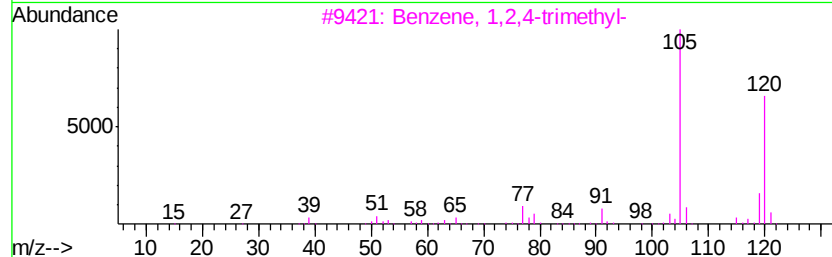
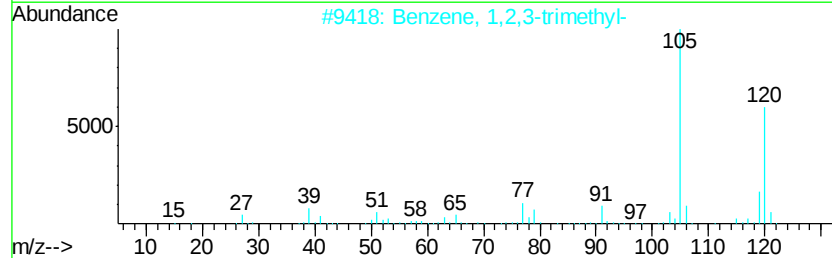
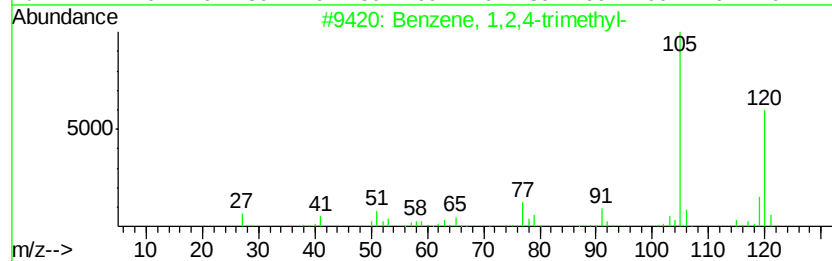
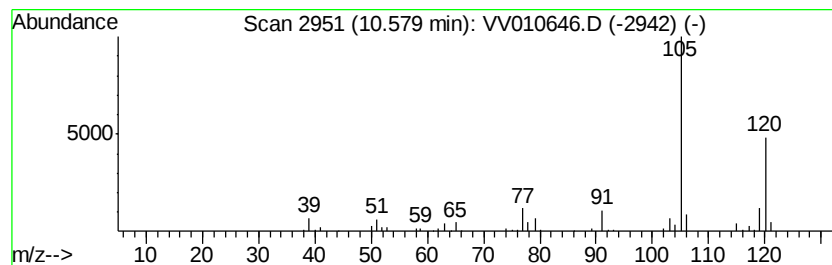
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Quant Title : VOC Analysis

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

Peak Number 3 Benzene, 1,2,4-trimethyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.58	8.28 ug/L	131890	1,4-Dichlorobenzene-d4	11.29

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



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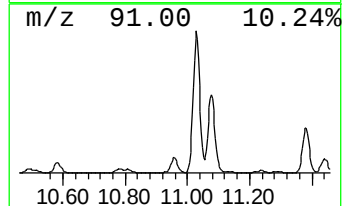
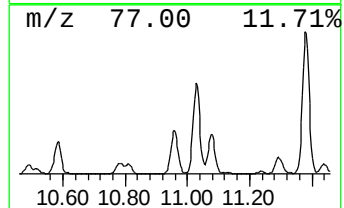
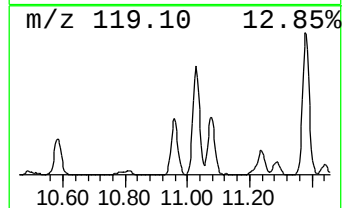
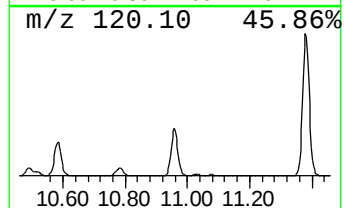
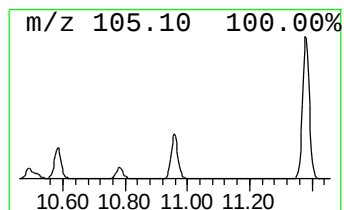
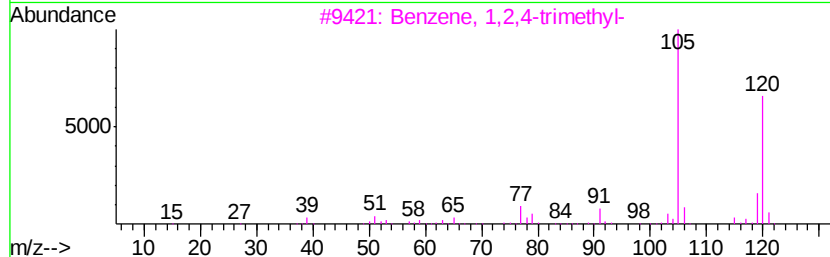
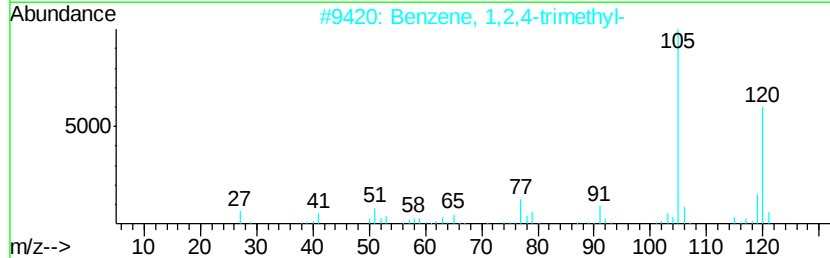
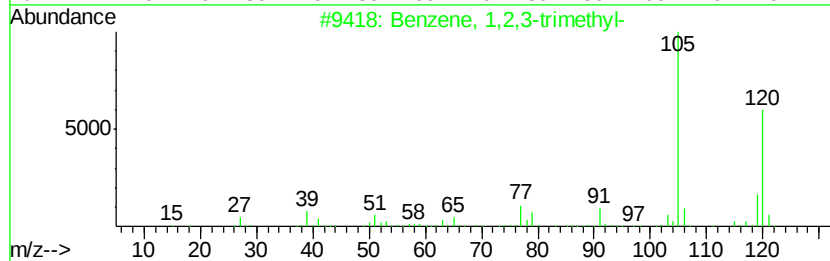
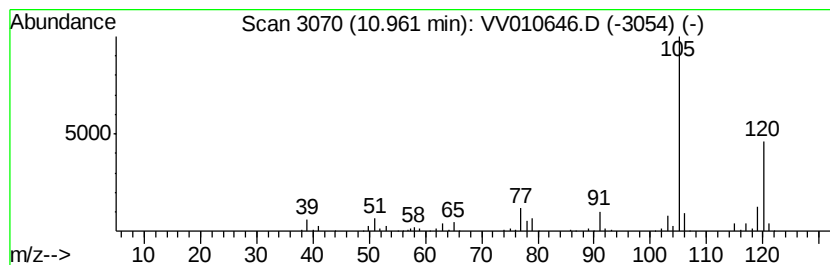
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Quant Title : VOC Analysis

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

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.96	12.00 ug/L	191203	1,4-Dichlorobenzene-d4	11.29

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		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4		Mesitylene	120	C9H12	000108-67-8	94
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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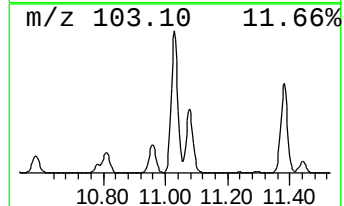
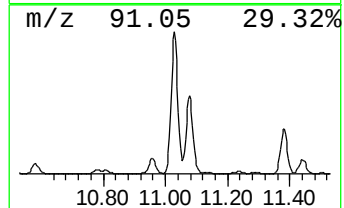
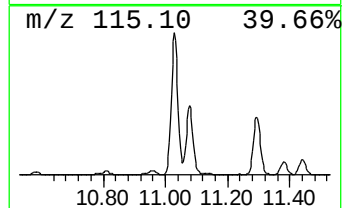
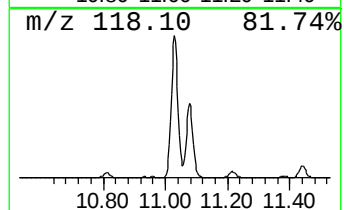
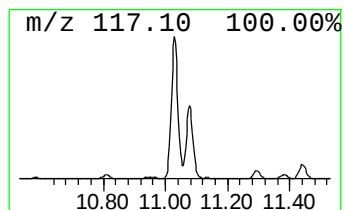
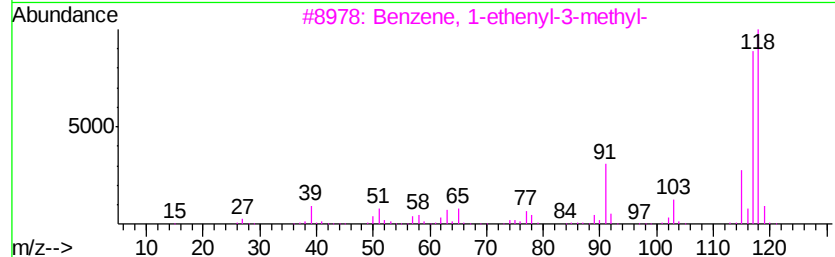
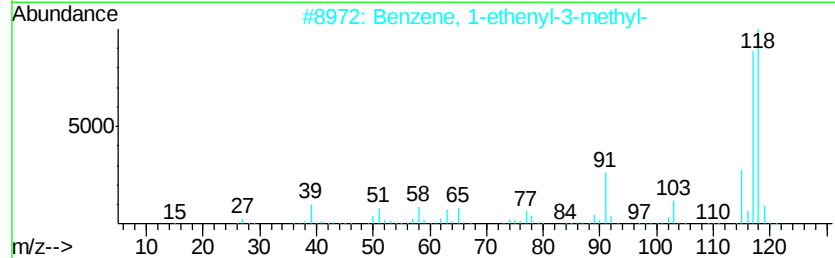
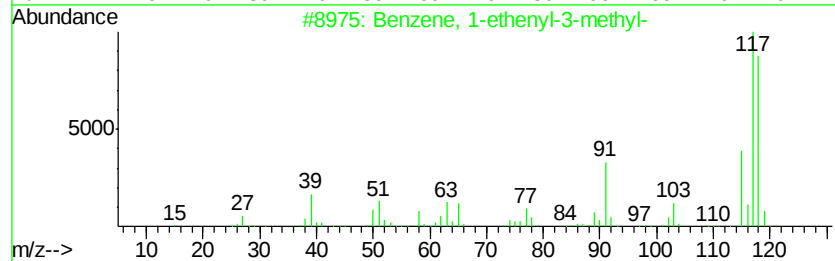
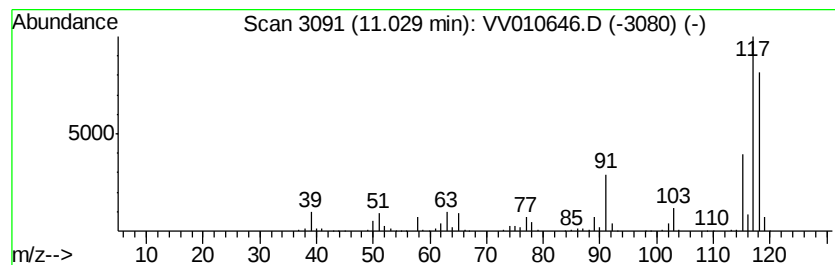
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Quant Title : VOC Analysis

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

Peak Number 6 Benzene, 1-ethenyl-3-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.03	54.90 ug/L	874765	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	96
2		Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	96
3		Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	96
4		Benzene, 1-propenyl-	118	C9H10	000637-50-3	95
5		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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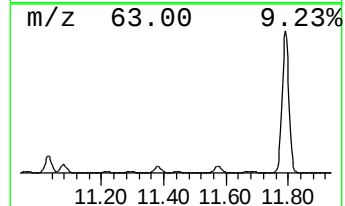
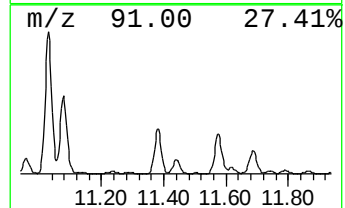
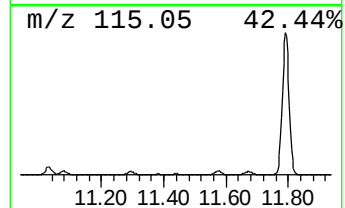
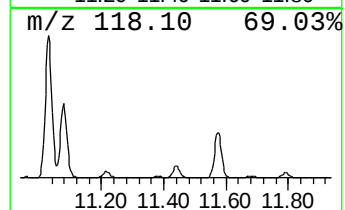
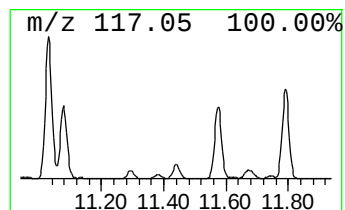
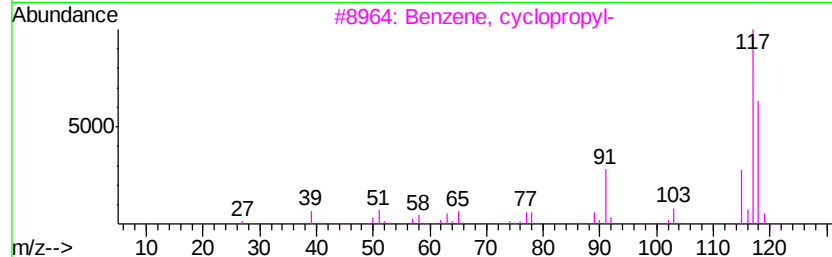
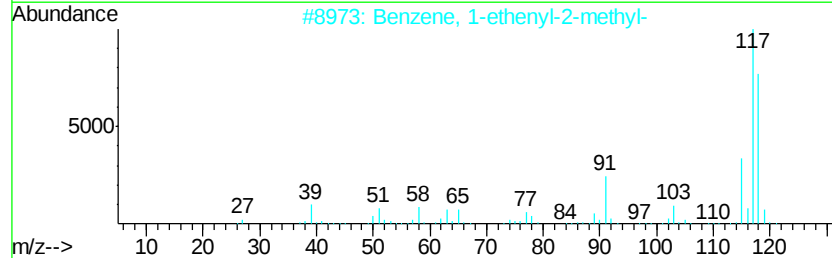
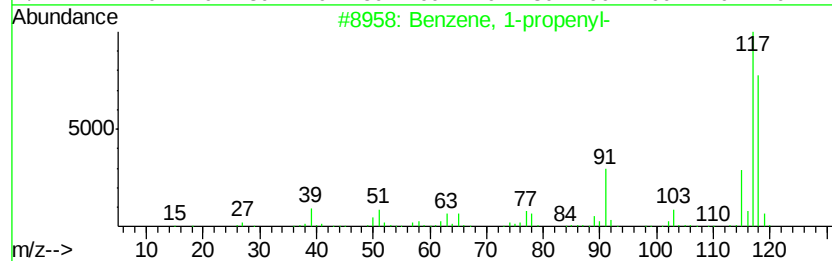
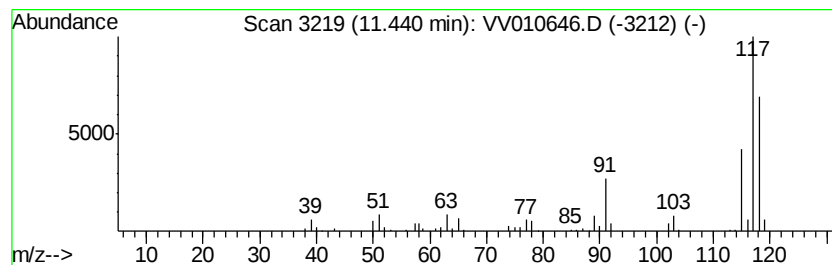
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Quant Title : VOC Analysis

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

Peak Number 9 Benzene, 1-propenyl- Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	5.28 ug/L	84172	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	95
2		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	93
3		Benzene, cyclopropyl-	118	C9H10	000873-49-4	92
4		Benzene, 1-propenyl-	118	C9H10	000637-50-3	89
5		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	89



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

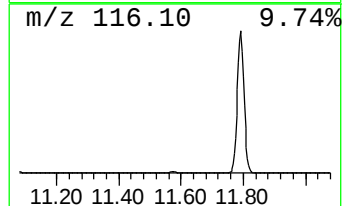
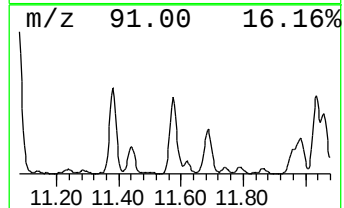
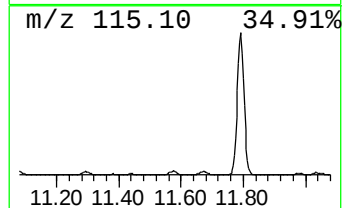
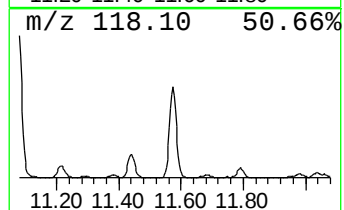
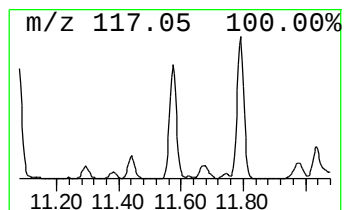
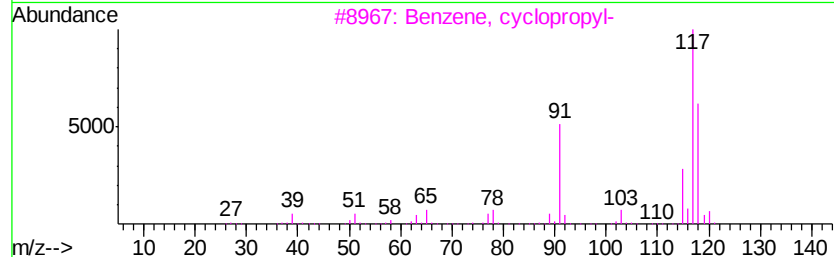
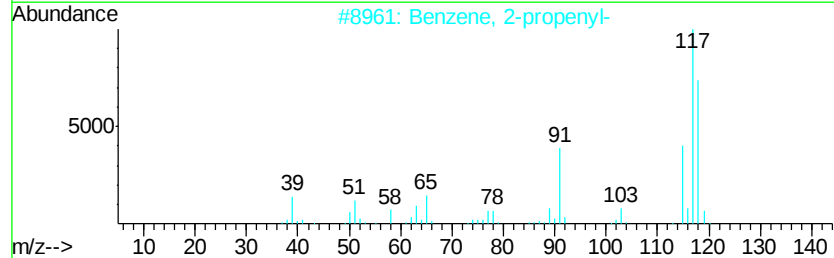
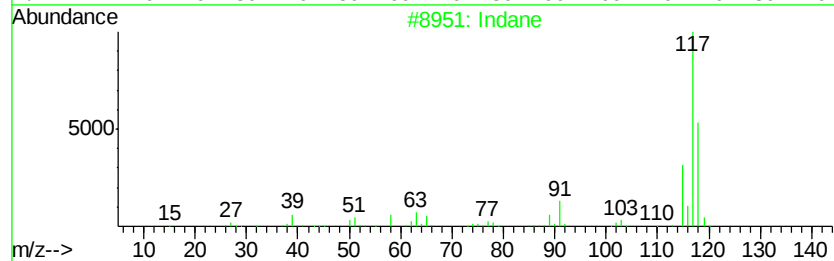
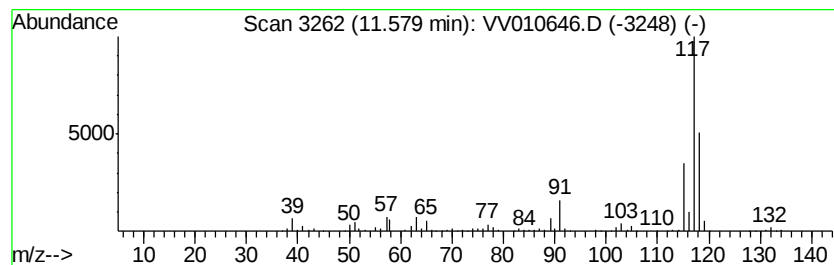
Instrument :
MSVOA_V
ClientSampleId :
GAJ72

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM050219S.M
Quant Title : VOC Analysis

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

Peak Number 10 Indane Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.58	25.62 ug/L	408185	1,4-Dichlorobenzene-d4	11.29
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane		118 C9H10	000496-11-7 95
2	Benzene, 2-propenyl-		118 C9H10	000300-57-2 83
3	Benzene, cyclopropyl-		118 C9H10	000873-49-4 83
4	Benzene, cyclopropyl-		118 C9H10	000873-49-4 81
5	Deltacyclene		118 C9H10	007785-10-6 72



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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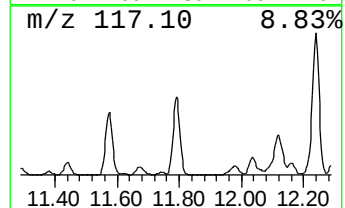
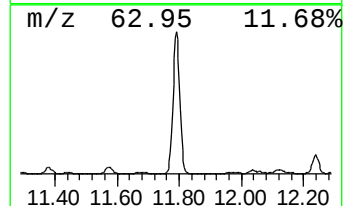
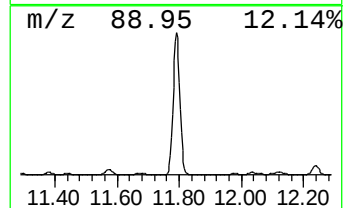
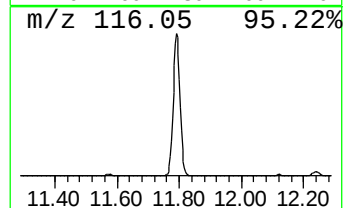
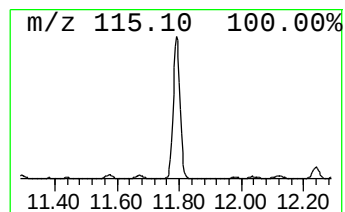
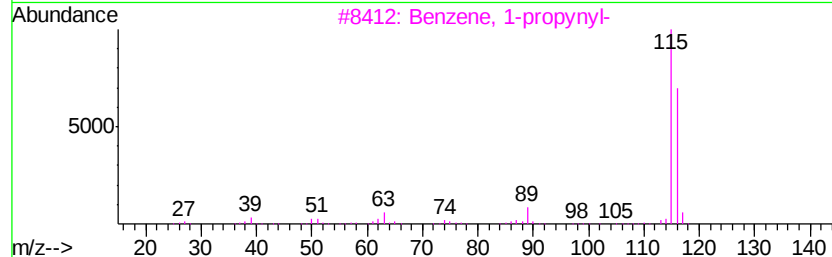
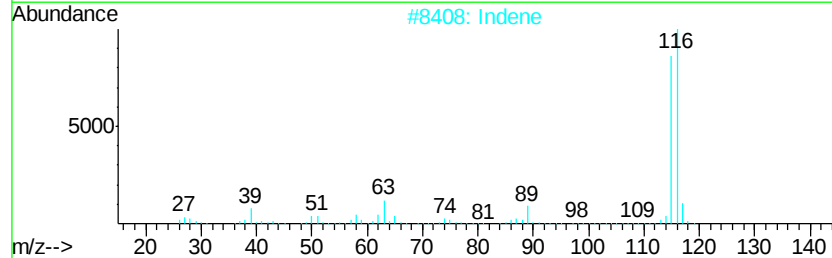
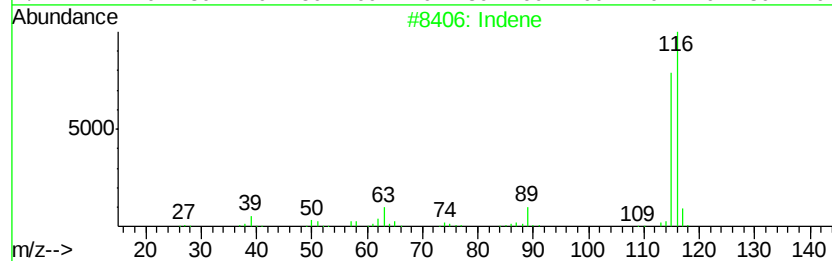
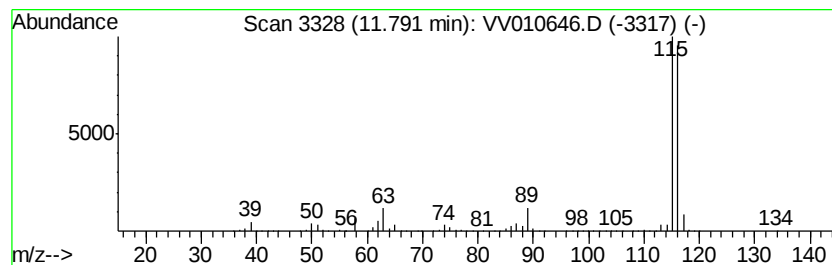
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Quant Title : VOC Analysis

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

Peak Number 11 Indene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	295.69 ug/L	4711080	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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		Benzene, 1-propynyl-	116	C9H8	000673-32-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
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Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAJ72

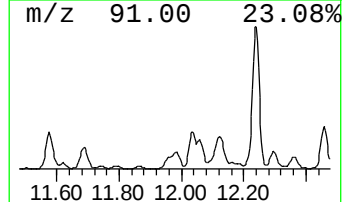
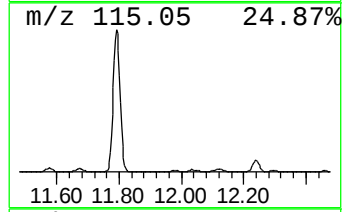
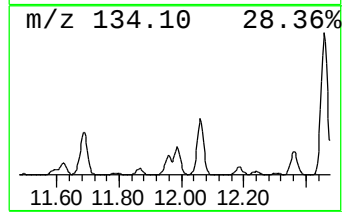
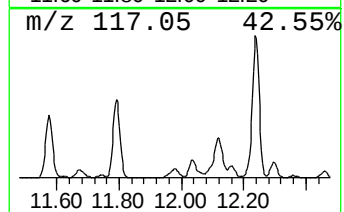
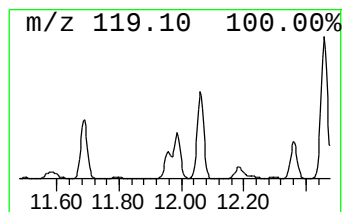
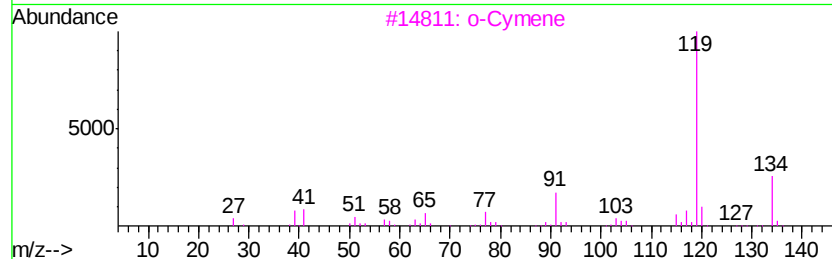
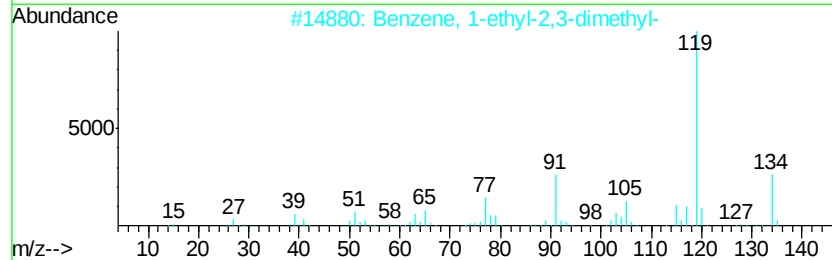
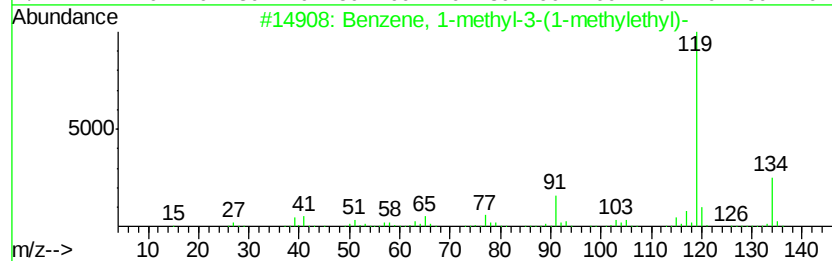
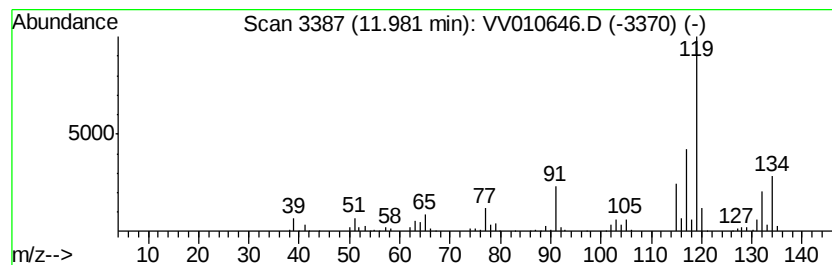
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Quant Title : VOC Analysis

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

Peak Number 12 Benzene, 1-methyl-3-(1-meth... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	15.75 ug/L	251002	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	70
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	70
3		o-Cymene	134	C10H14	000527-84-4	70
4		o-Cymene	134	C10H14	000527-84-4	70
5		p-Cymene	134	C10H14	000099-87-6	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
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ALS Vial : 1 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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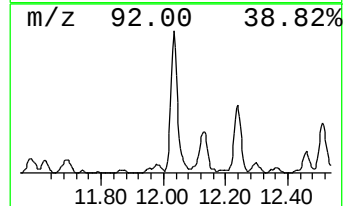
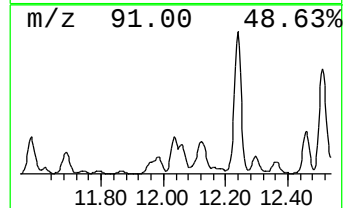
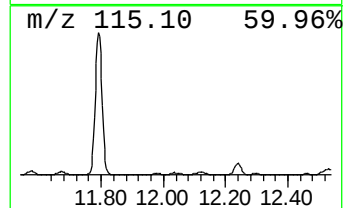
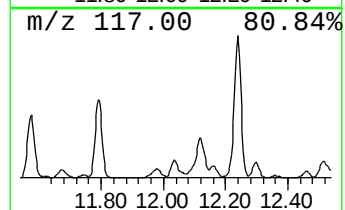
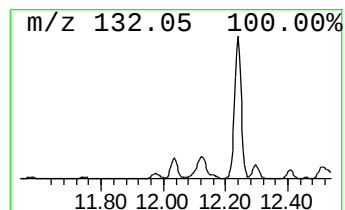
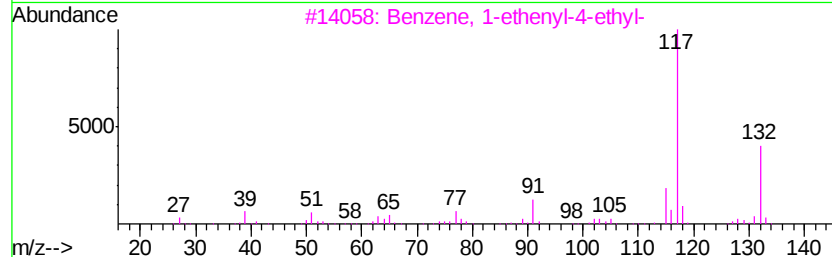
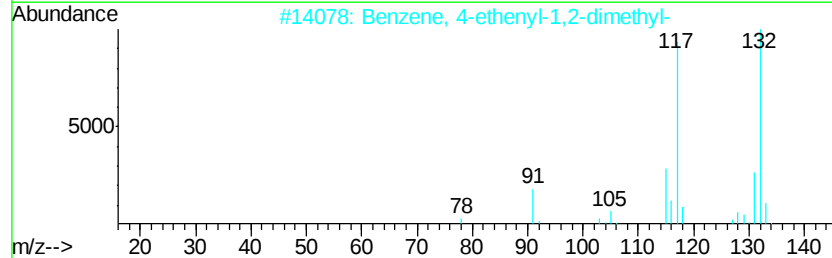
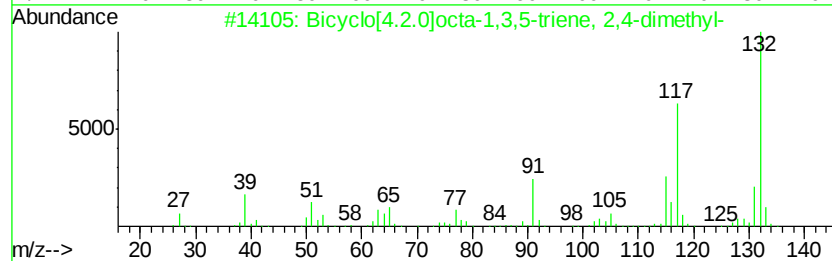
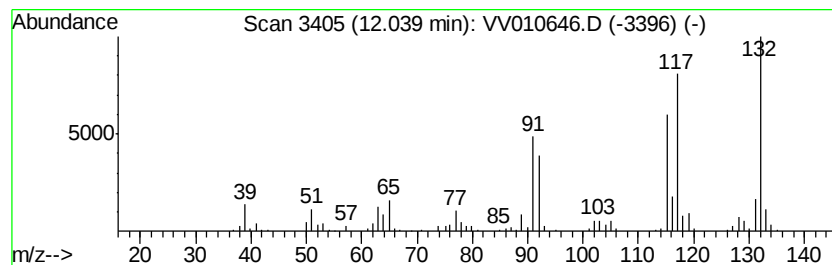
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Quant Title : VOC Analysis

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

Peak Number 13 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.04	12.78 ug/L	203566	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	94
2		Benzene, 4-ethenyl-1,2-dimethyl-	132	C10H12	027831-13-6	89
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	70
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	70
5		Benzene, 1-methyl-4-(1-methyleth...	132	C10H12	001195-32-0	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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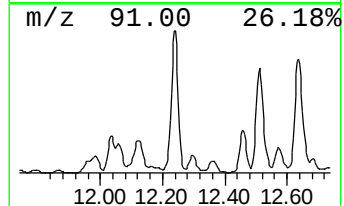
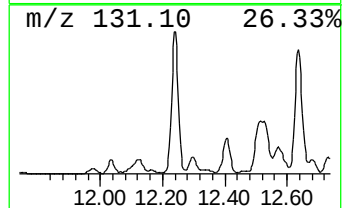
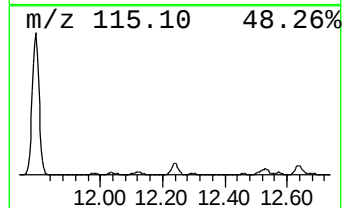
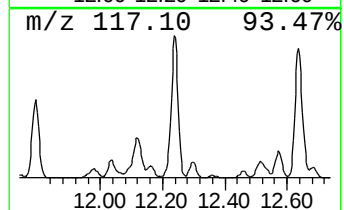
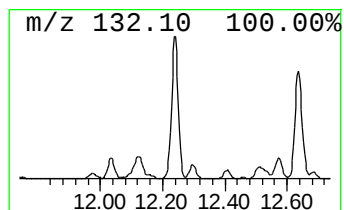
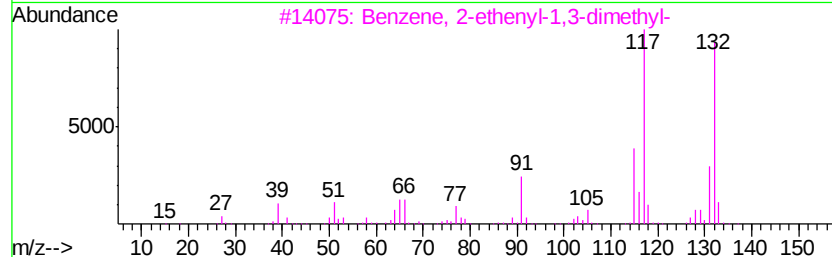
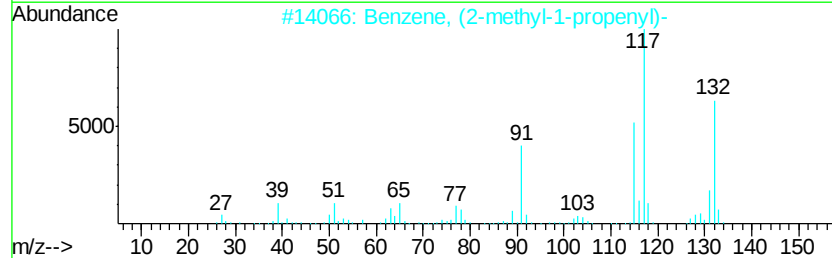
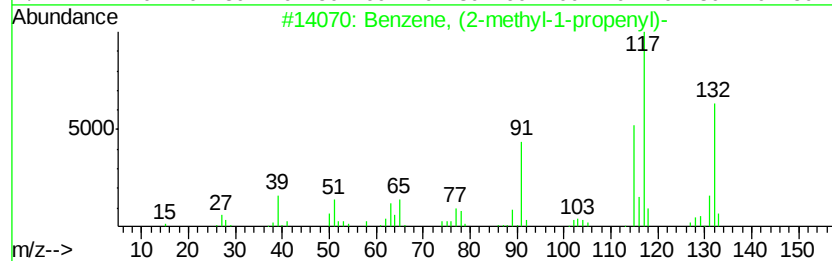
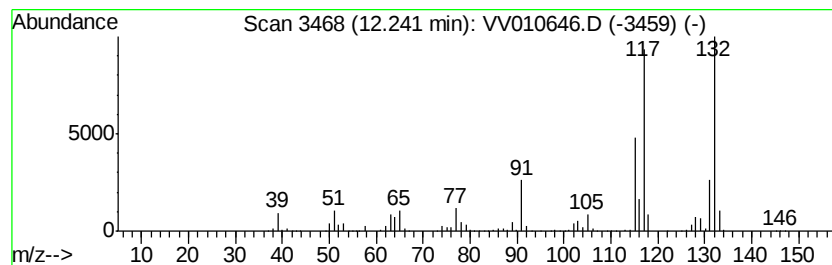
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Quant Title : VOC Analysis

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

Peak Number 14 Benzene, (2-methyl-1-propen... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.24	80.54 ug/L	1283130	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	95
2		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	95
3		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	95
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	95
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampled :
GAJ72

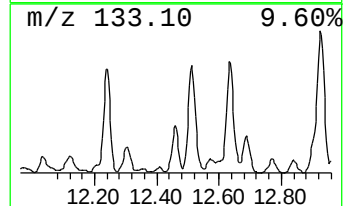
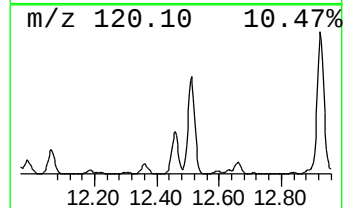
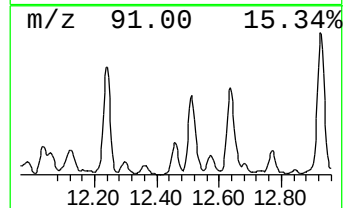
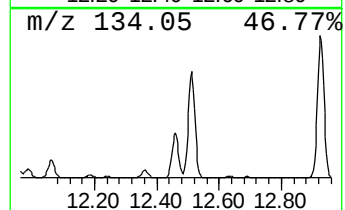
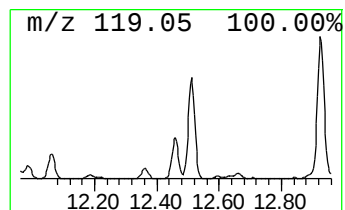
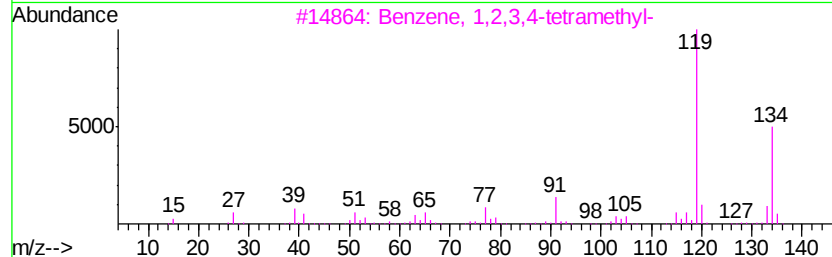
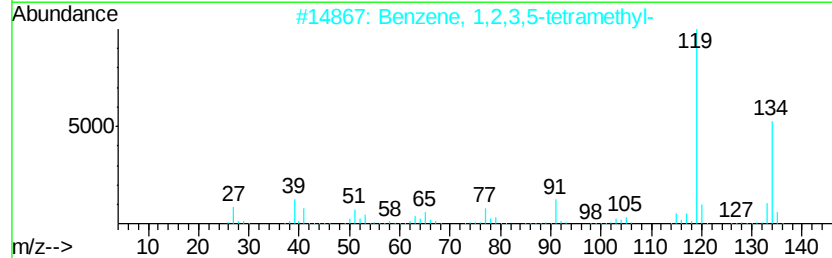
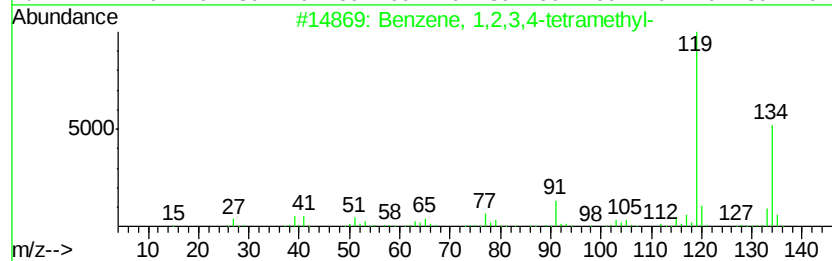
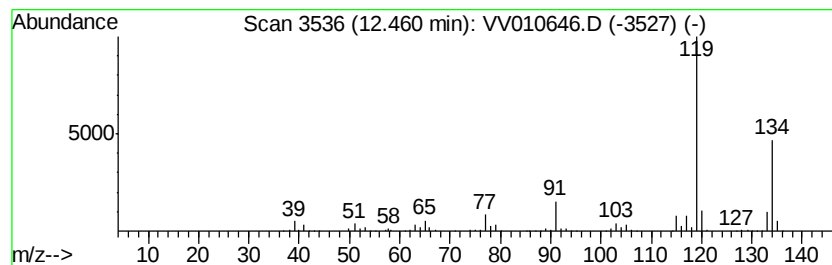
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Quant Title : VOC Analysis

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

Peak Number 15 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.46	22.13 ug/L	352535	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
5		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAJ72

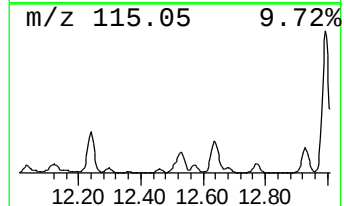
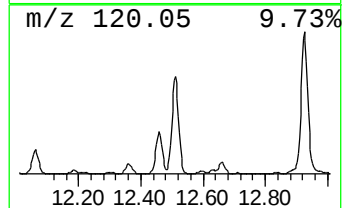
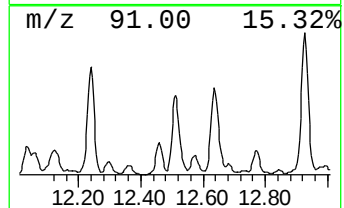
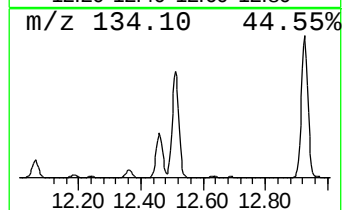
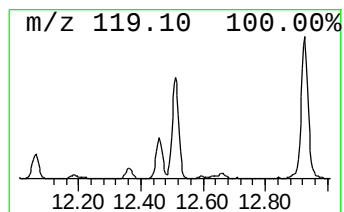
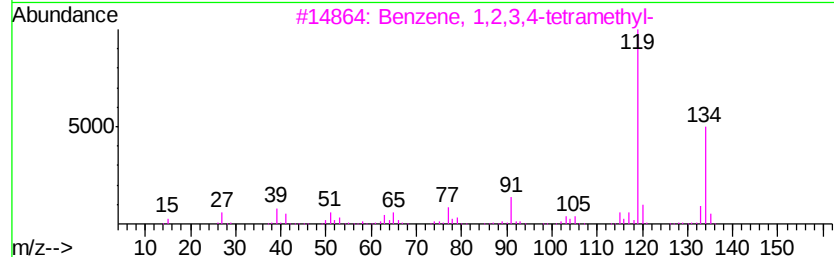
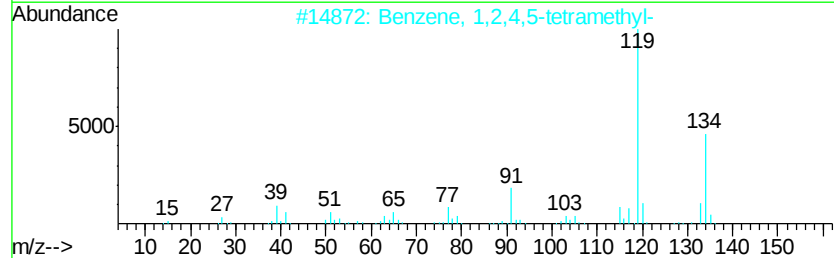
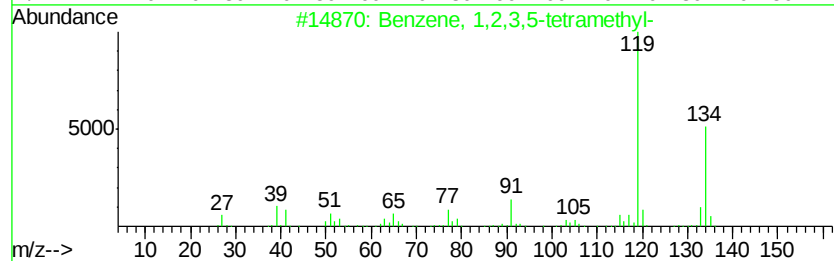
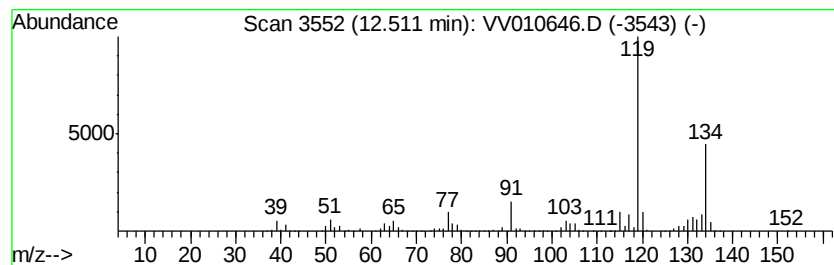
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Quant Title : VOC Analysis

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

Peak Number 16 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.51	96.36 ug/L	1535190	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	93
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	93
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAJ72

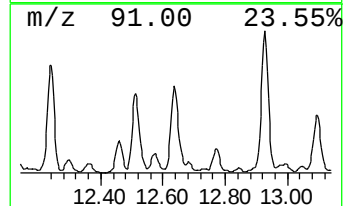
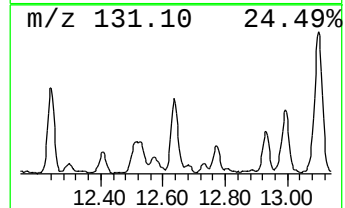
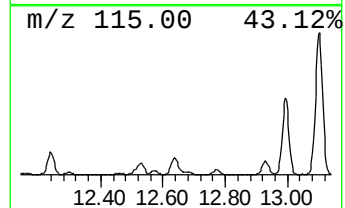
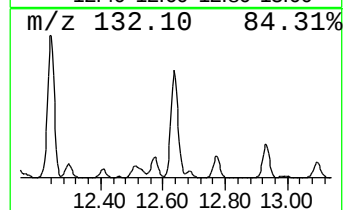
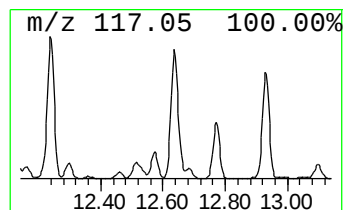
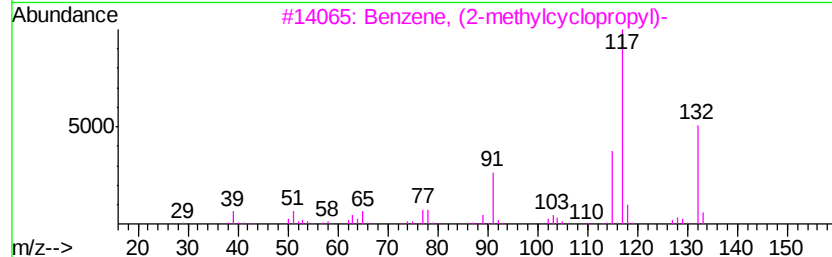
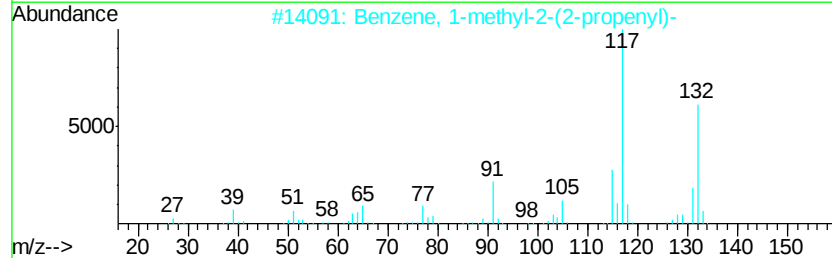
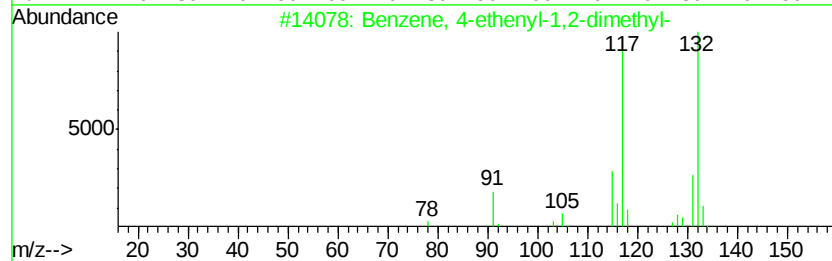
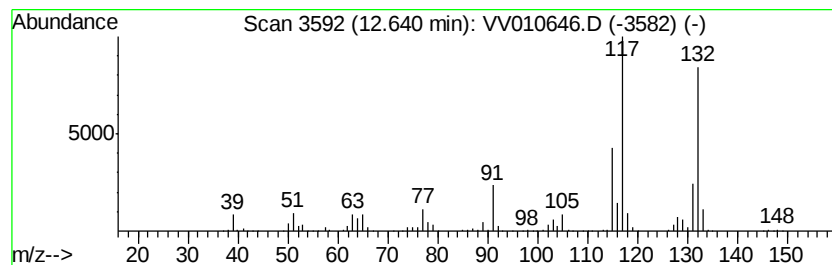
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Quant Title : VOC Analysis

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

Peak Number 17 Benzene, 4-ethenyl-1,2-dime... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.64	74.58 ug/L	1188220	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethenyl-1,2-dimethyl-	132	C10H12	027831-13-6	95
2		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	95
3		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	95
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	94
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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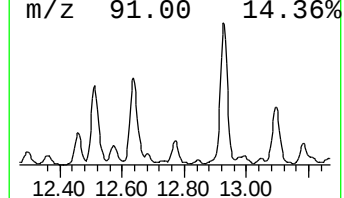
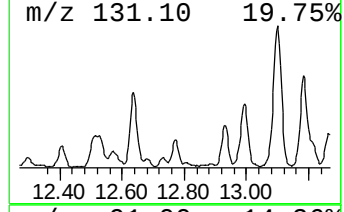
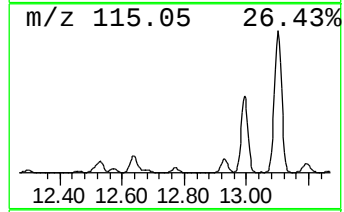
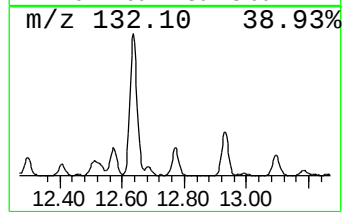
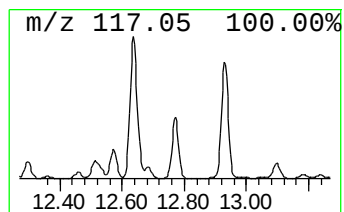
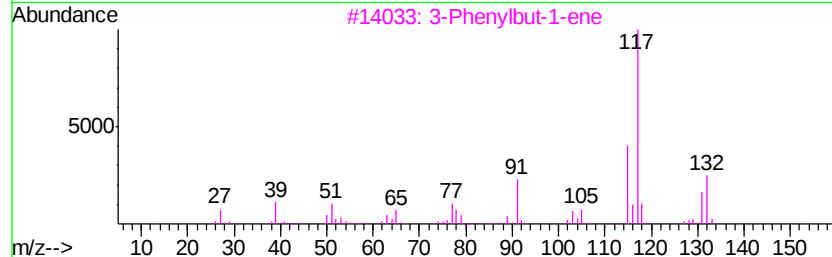
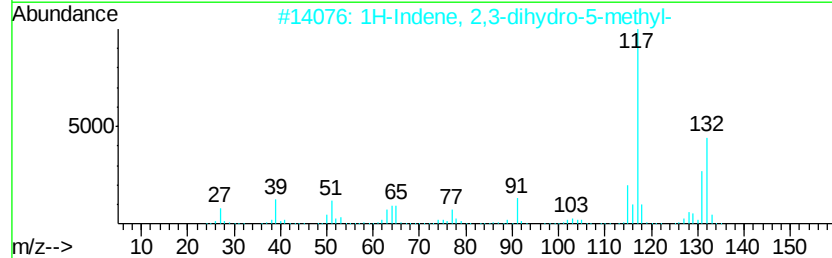
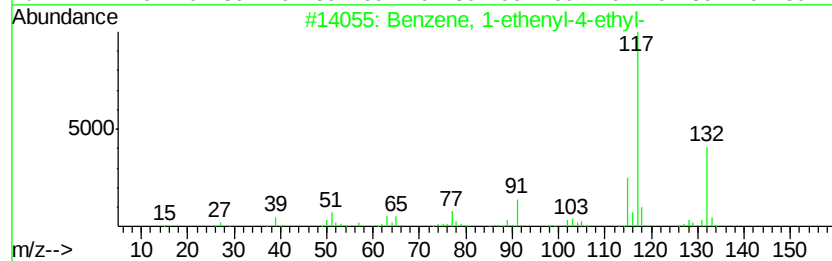
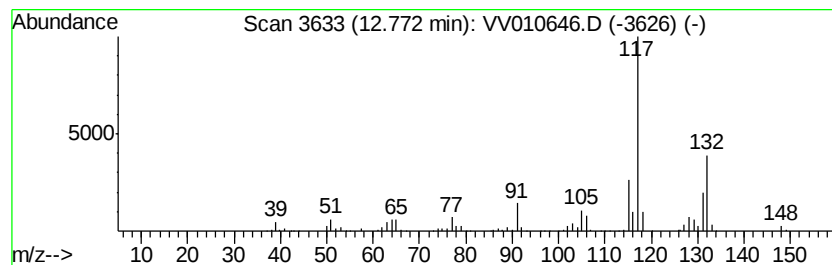
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Quant Title : VOC Analysis

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

Peak Number 18 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	21.44 ug/L	341526	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
5		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAJ72

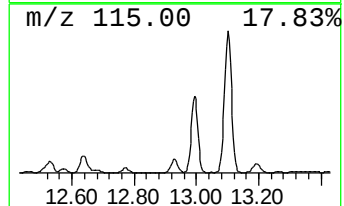
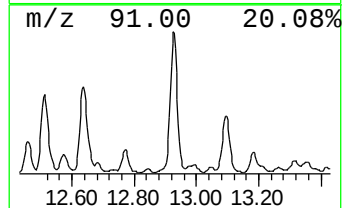
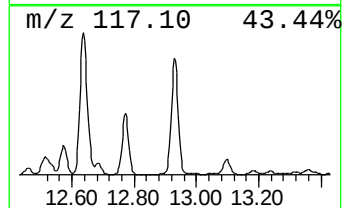
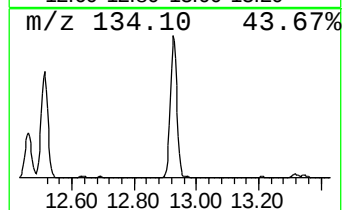
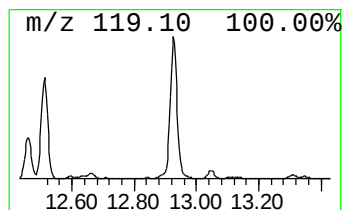
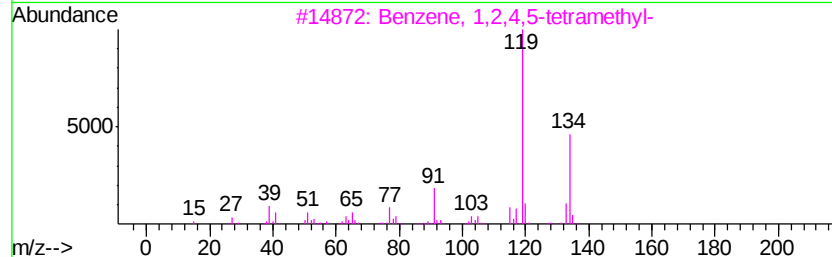
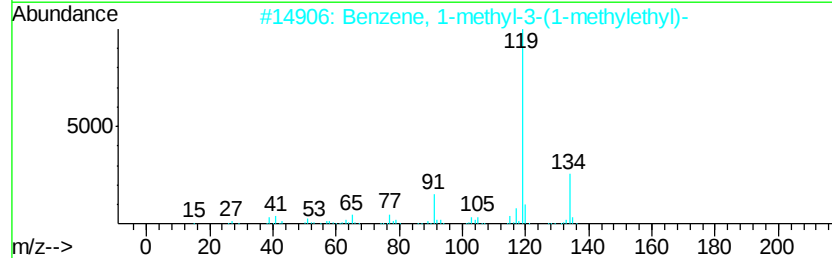
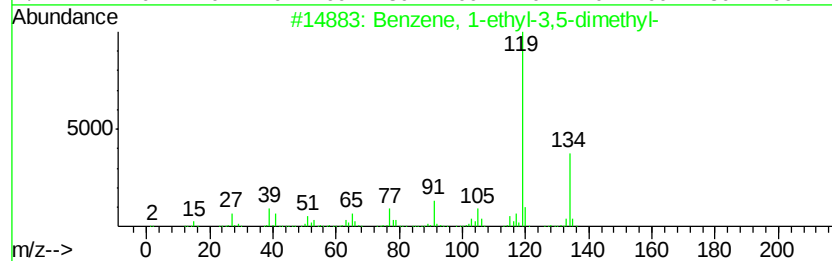
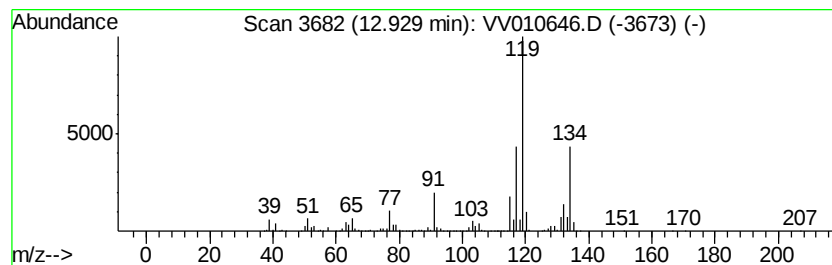
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Quant Title : VOC Analysis

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

Peak Number 19 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.93	106.09 ug/L	1690220	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	76
2		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	70
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	70
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	70
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAJ72

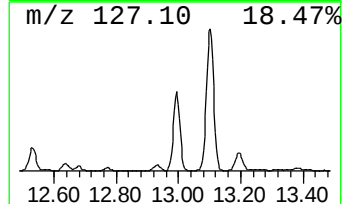
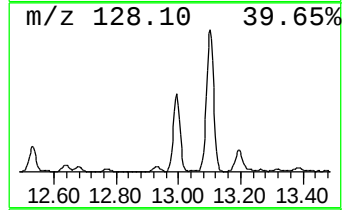
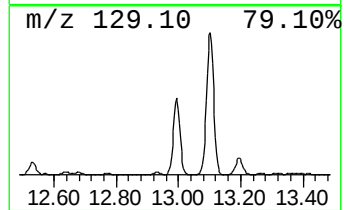
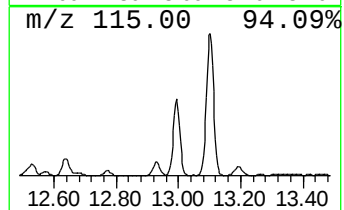
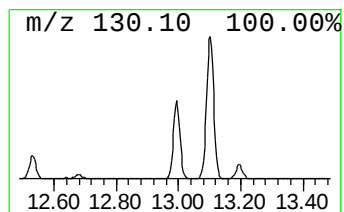
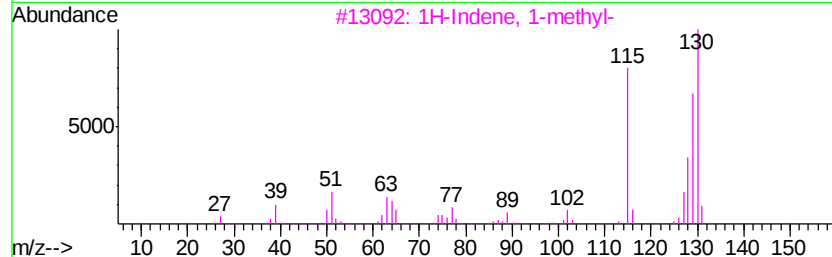
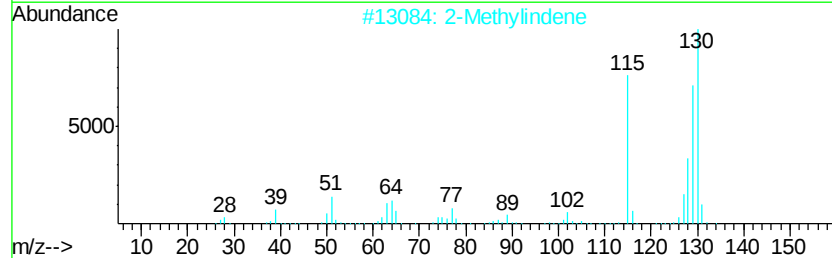
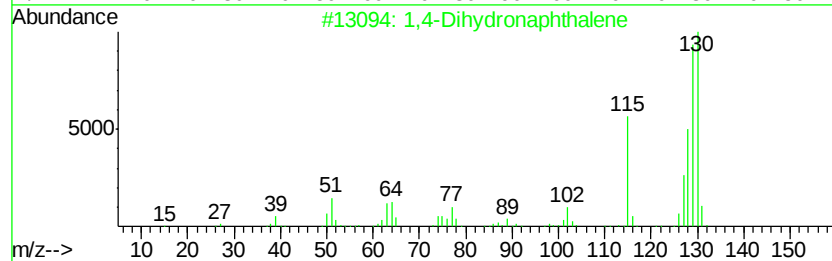
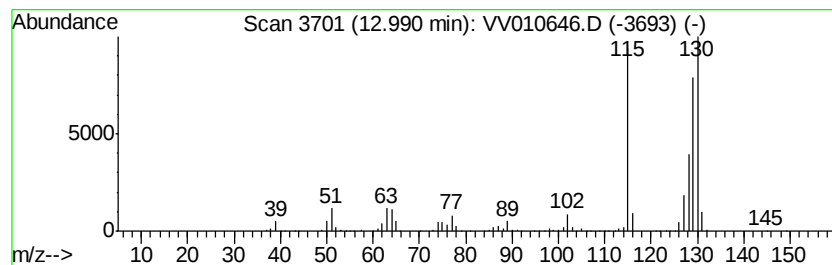
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Quant Title : VOC Analysis

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

Peak Number 20 1,4-Dihydronaphthalene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.99	140.42 ug/L	2237210	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,4-Dihydronaphthalene	130	C10H10	000612-17-9	96
2		2-Methylindene	130	C10H10	002177-47-1	94
3		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	94
4		Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	94
5		Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAJ72

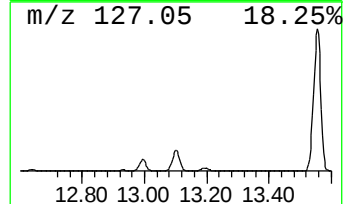
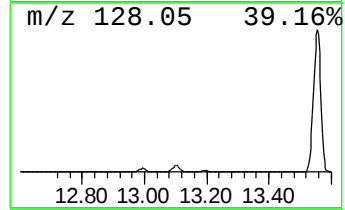
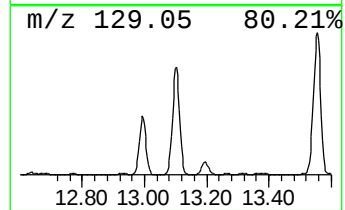
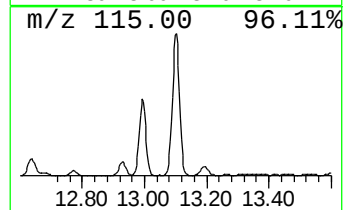
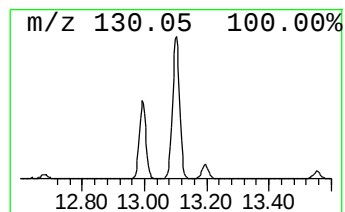
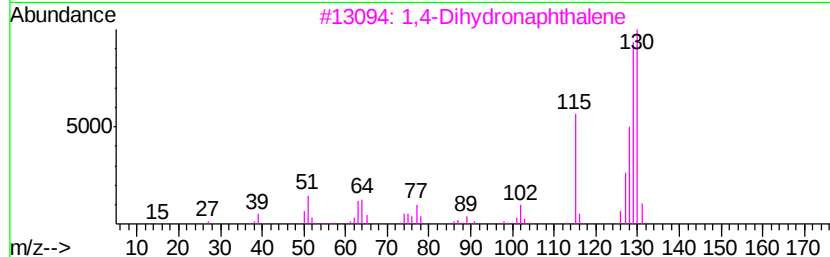
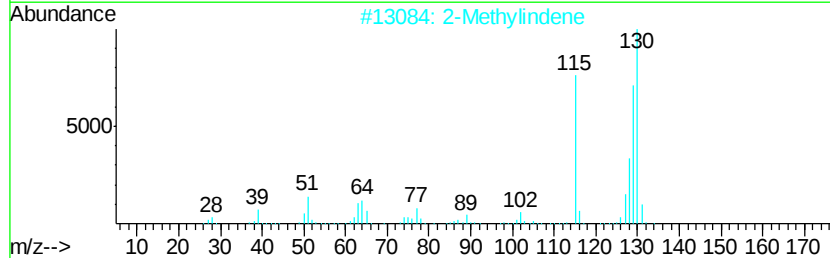
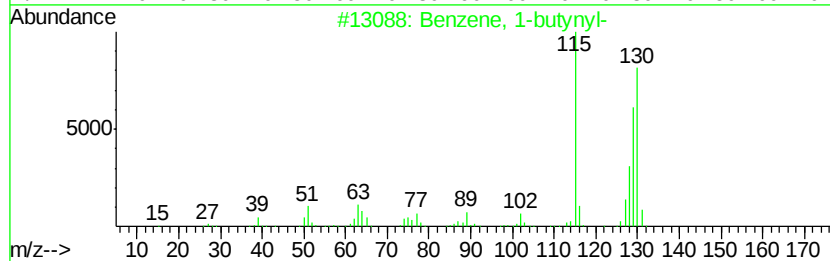
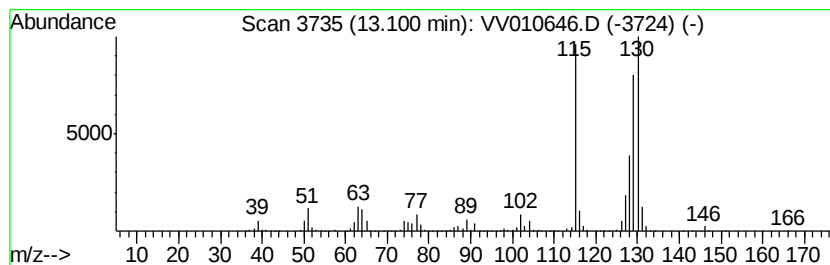
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Quant Title : VOC Analysis

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

Peak Number 21 Benzene, 1-butynyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.10	298.76 ug/L	4759990	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-butynyl-	130	C10H10	000622-76-4	95
2		2-Methylindene	130	C10H10	002177-47-1	95
3		1,4-Dihydronaphthalene	130	C10H10	000612-17-9	95
4		Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	94
5		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampled :
GAJ72

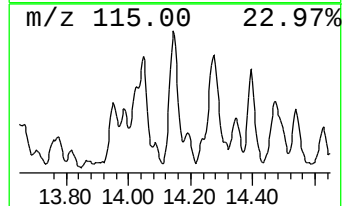
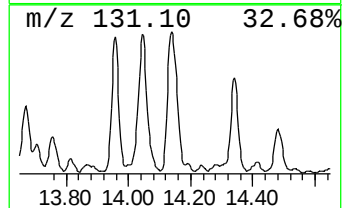
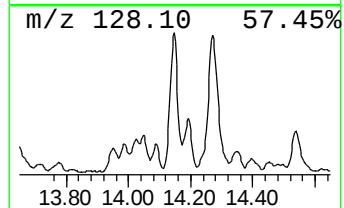
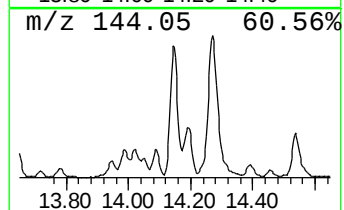
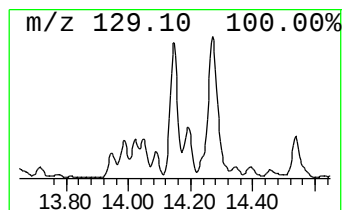
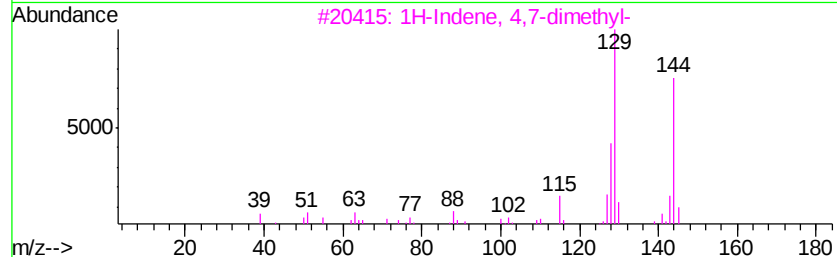
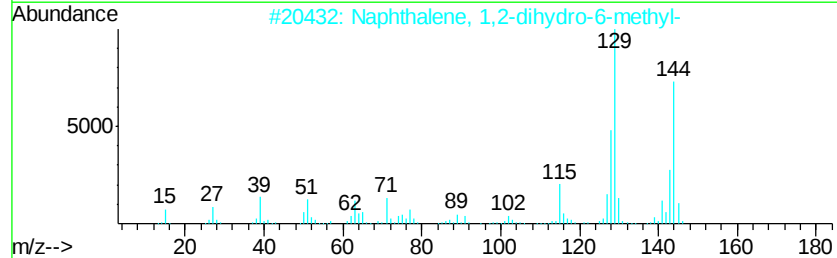
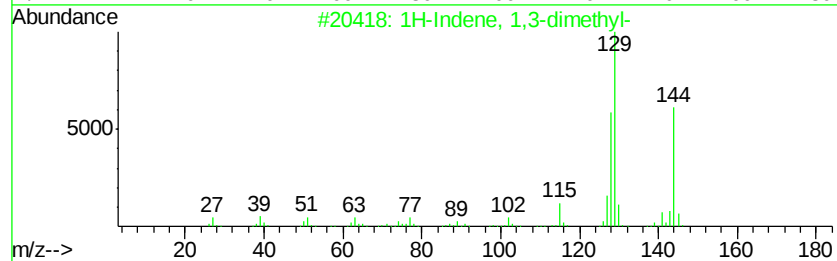
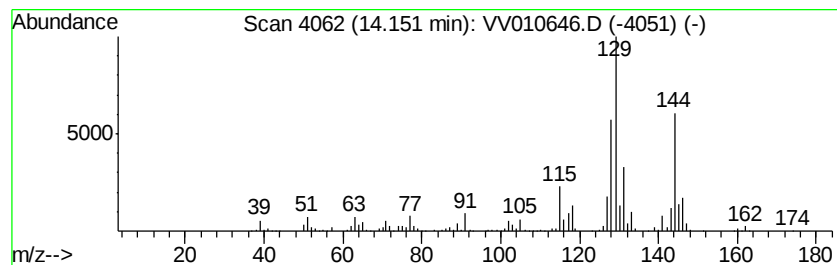
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Quant Title : VOC Analysis

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

Peak Number 24 1H-Indene, 1,3-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.15	41.54 ug/L	661900	1,4-Dichlorobenzene-d4	11.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 1,3-dimethyl-	144	C11H12	002177-48-2	94
2			Naphthalene, 1,2-dihydro-6-methyl-	144	C11H12	002717-47-7	90
3			1H-Indene, 4,7-dimethyl-	144	C11H12	006974-97-6	81
4			1H-Indene, 2,3-dimethyl-	144	C11H12	004773-82-4	81
5			1H-Indene, 1,1-dimethyl-	144	C11H12	018636-55-0	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAJ72

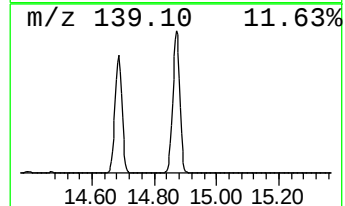
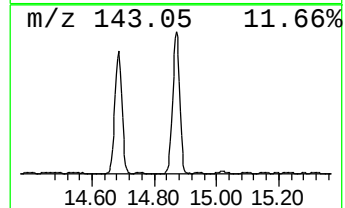
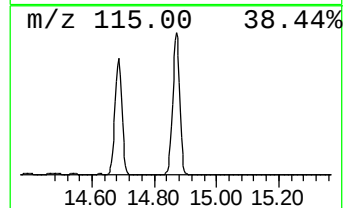
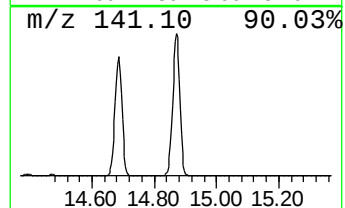
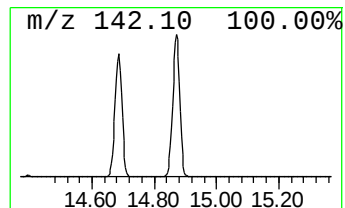
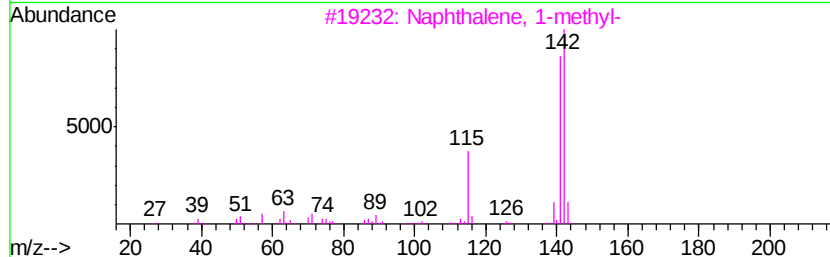
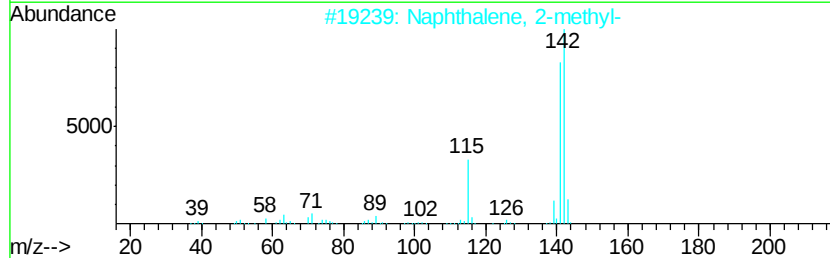
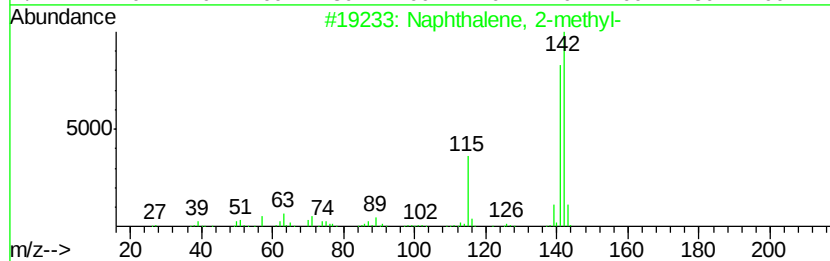
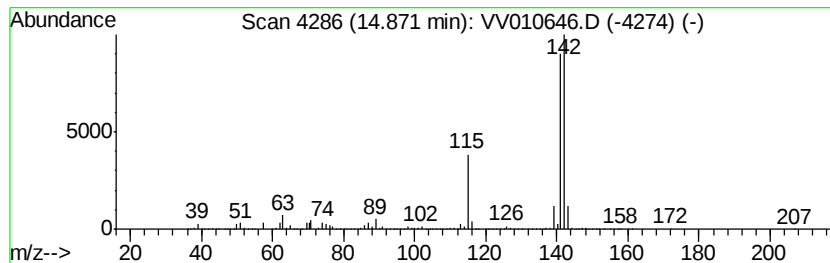
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Quant Title : VOC Analysis

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

Peak Number 26 Naphthalene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	941.79 ug/L	15005000	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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, 1-methyl-	142	C11H10	000090-12-0	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAJ72

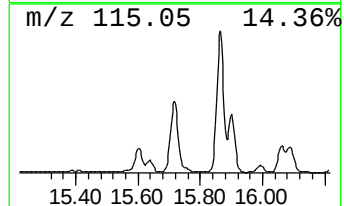
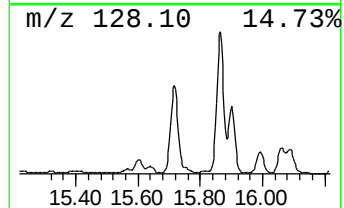
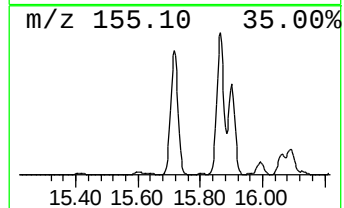
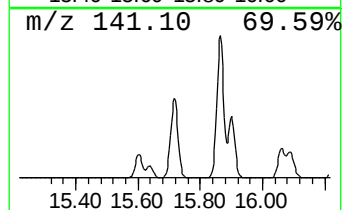
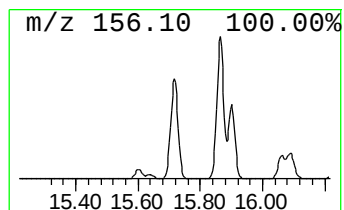
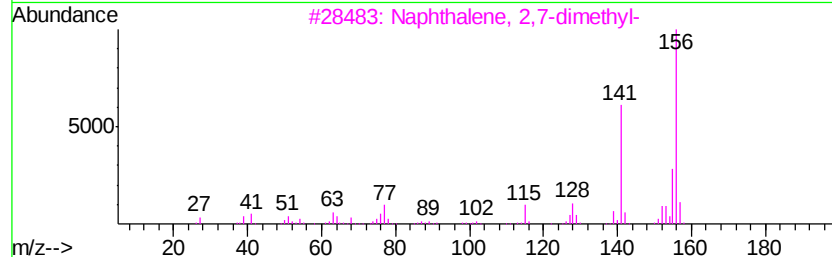
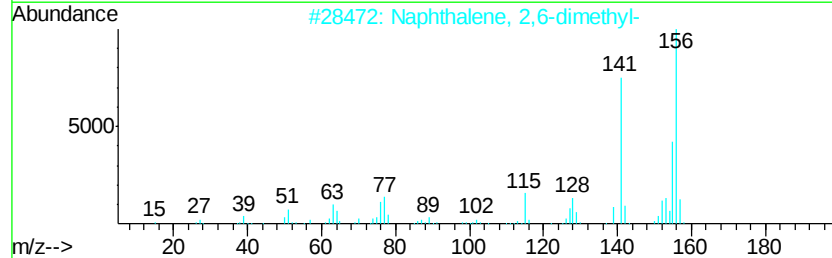
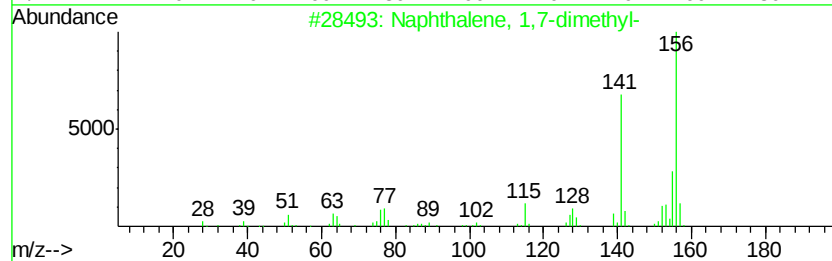
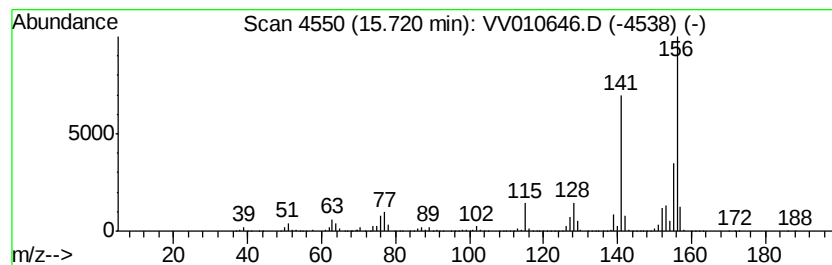
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Quant Title : VOC Analysis

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

Peak Number 27 Naphthalene, 1,7-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.72	238.64 ug/L	3802090	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
2		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
5		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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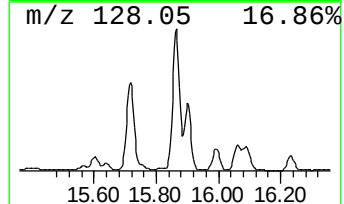
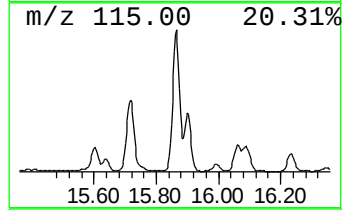
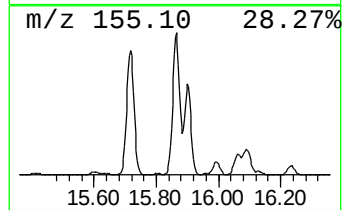
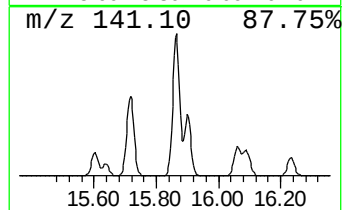
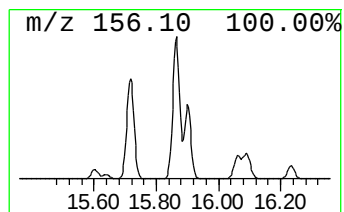
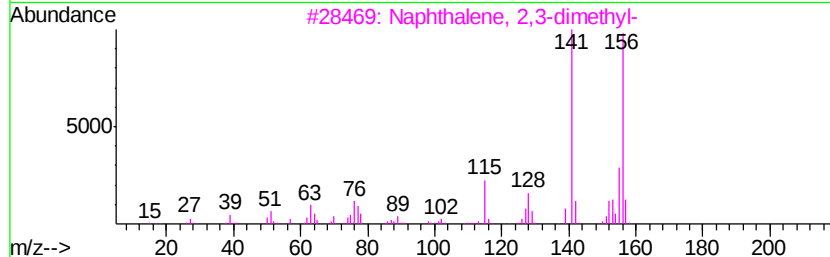
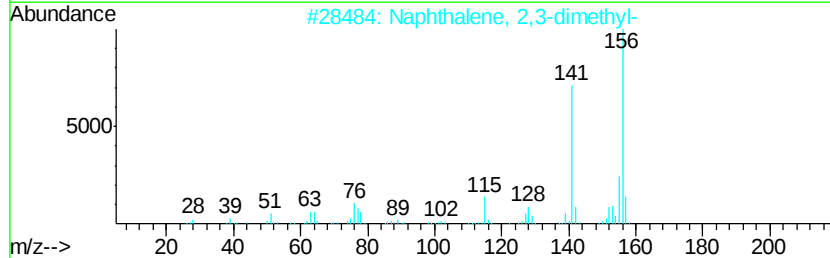
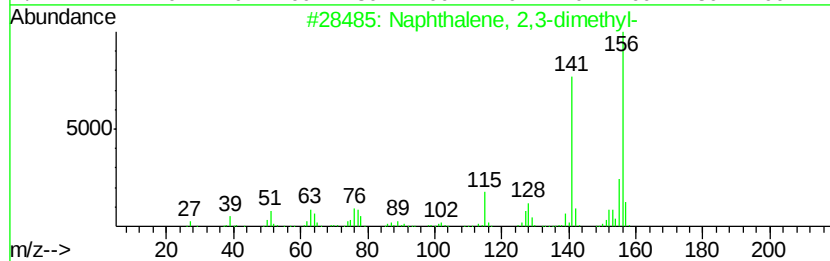
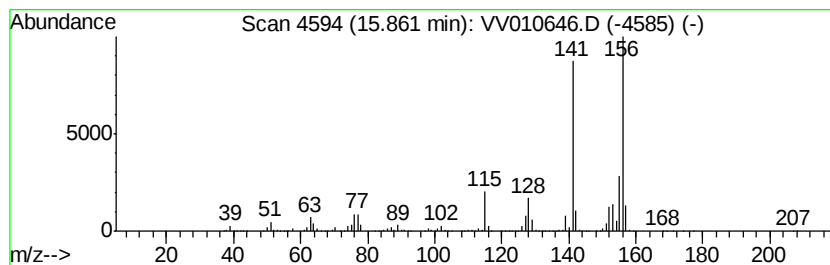
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Quant Title : VOC Analysis

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

Peak Number 28 Naphthalene, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.86	310.94 ug/L	4953990	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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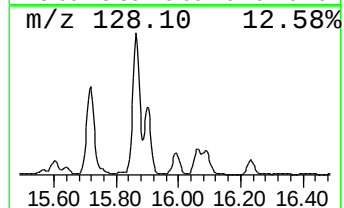
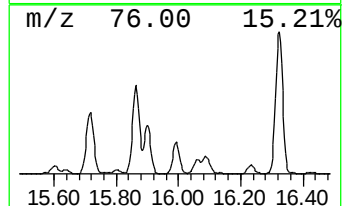
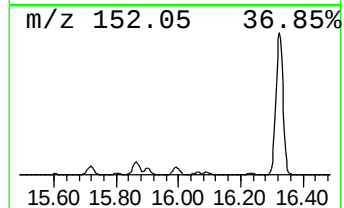
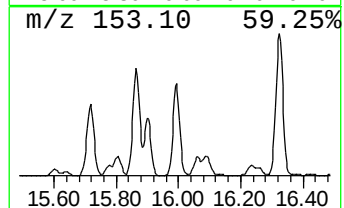
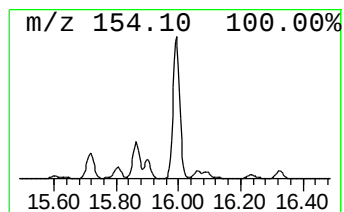
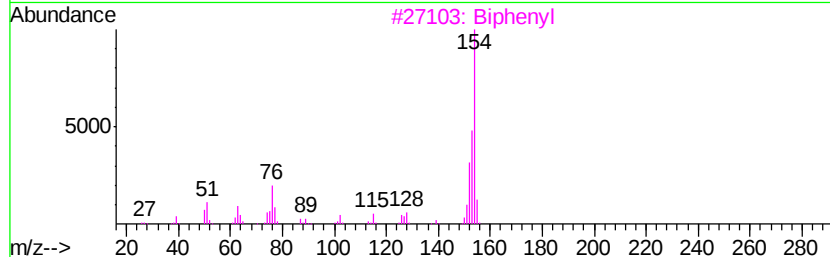
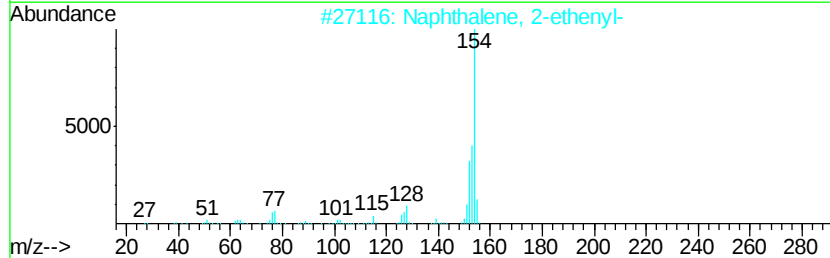
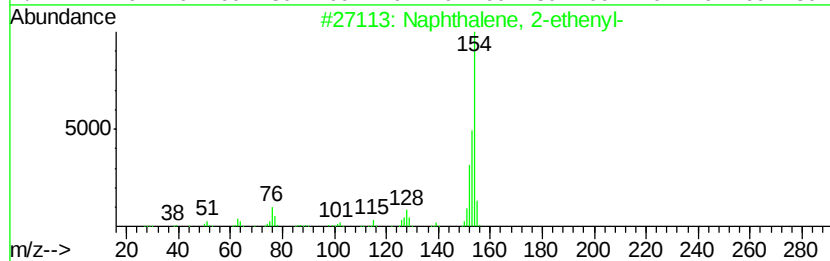
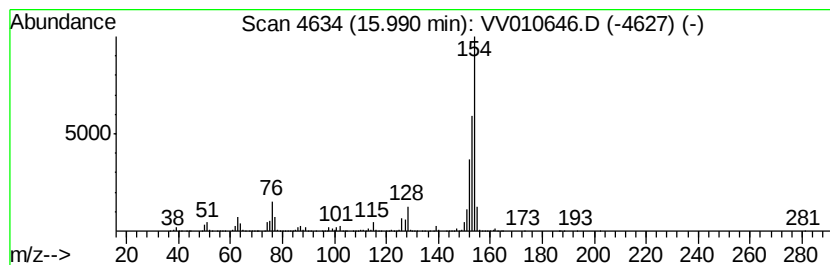
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Quant Title : VOC Analysis

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

Peak Number 29 Naphthalene, 2-ethenyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.99	50.93 ug/L	811395	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	94
2		Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	93
3		Biphenyl	154	C12H10	000092-52-4	90
4		Acenaphthene	154	C12H10	000083-32-9	81
5		Biphenyl	154	C12H10	000092-52-4	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAJ72

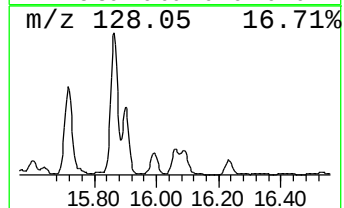
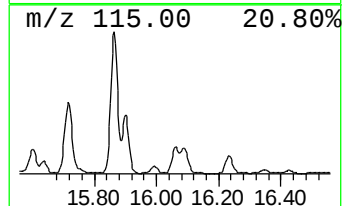
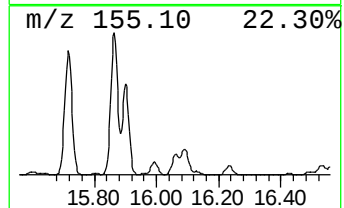
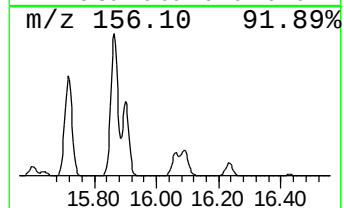
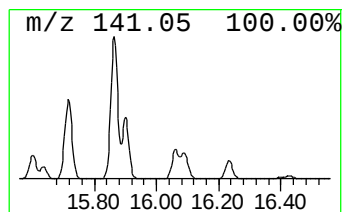
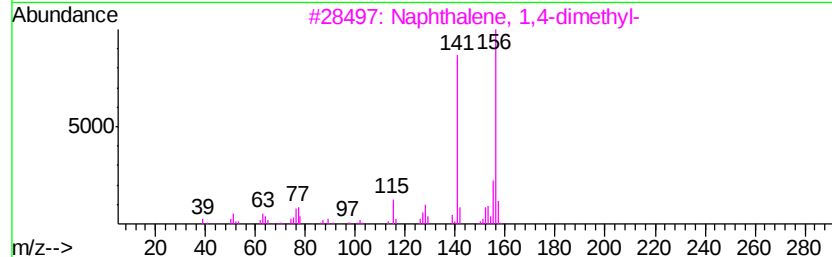
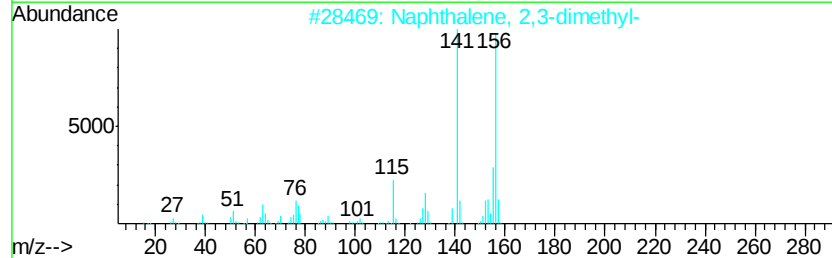
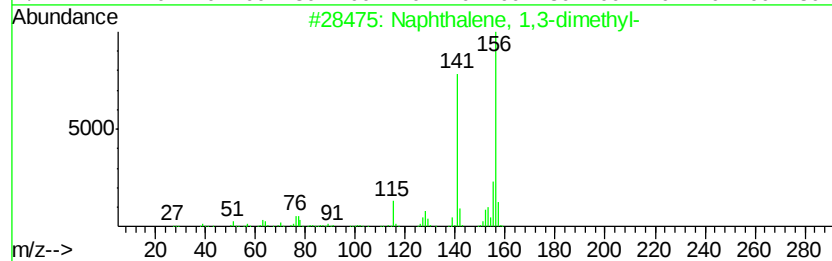
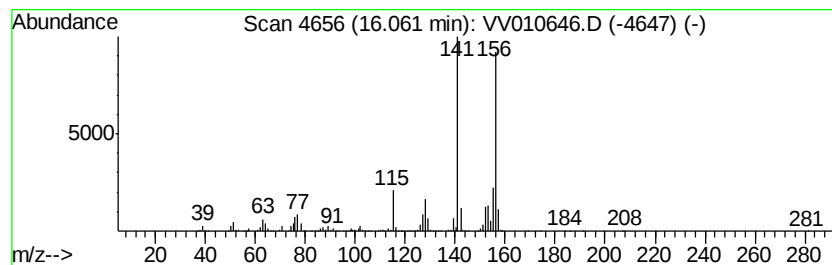
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Quant Title : VOC Analysis

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

Peak Number 30 Naphthalene, 1,3-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.06	55.20 ug/L	879479	1,4-Dichlorobenzene-d4	11.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
5			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAJ72

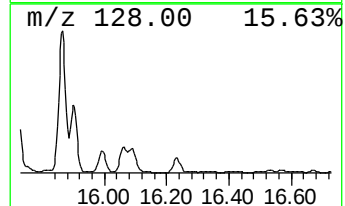
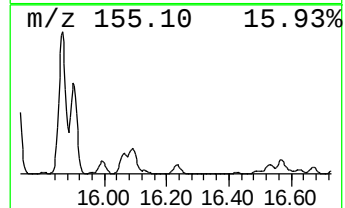
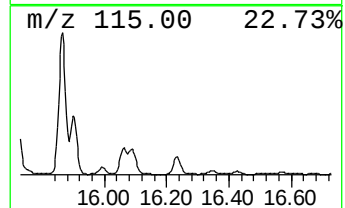
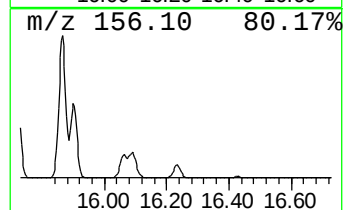
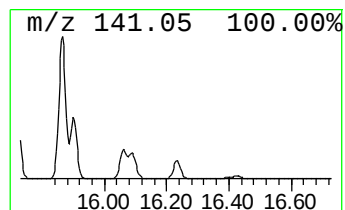
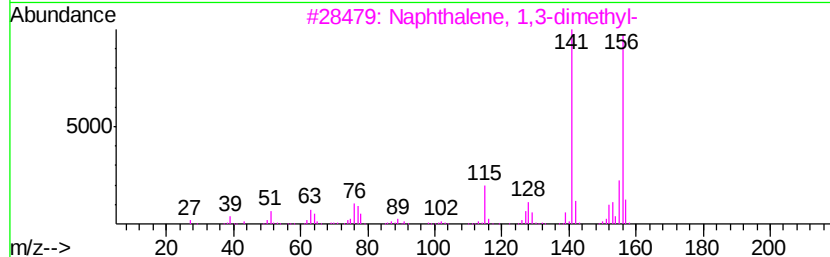
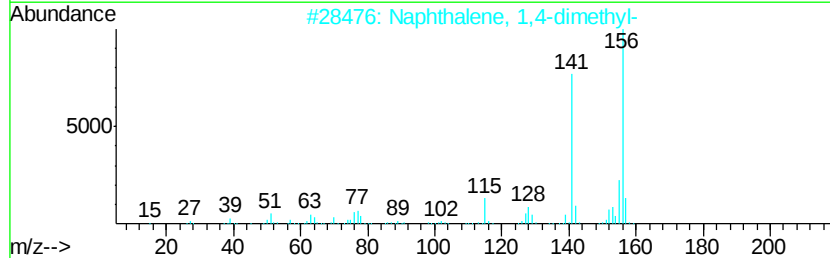
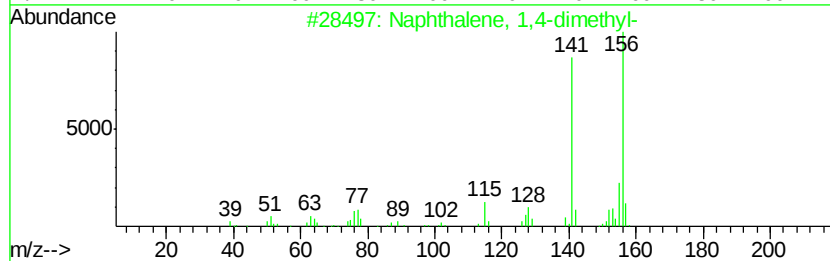
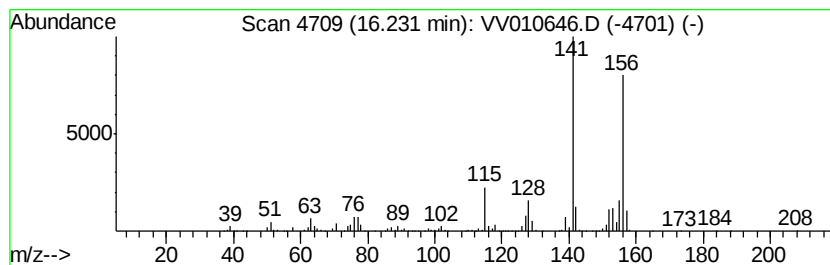
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Quant Title : VOC Analysis

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

Peak Number 31 Naphthalene, 1,4-dimethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.23	35.35 ug/L	563211	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
2		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
3		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
4		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
5		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV050219\
Data File : VV010646.D
Acq On : 03 May 2019 03:51
Operator : SY/MD
Sample : K2603-20
Misc : 5.00G/10ML/MSVOA V/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAJ72

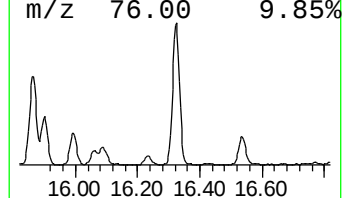
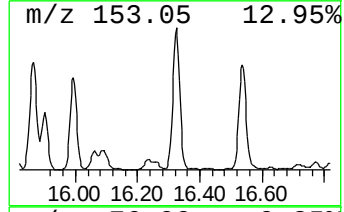
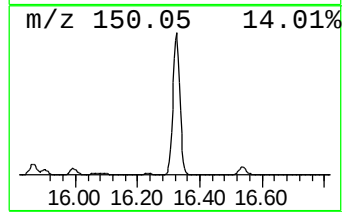
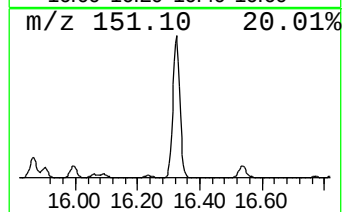
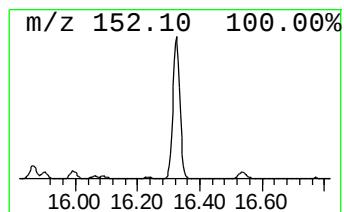
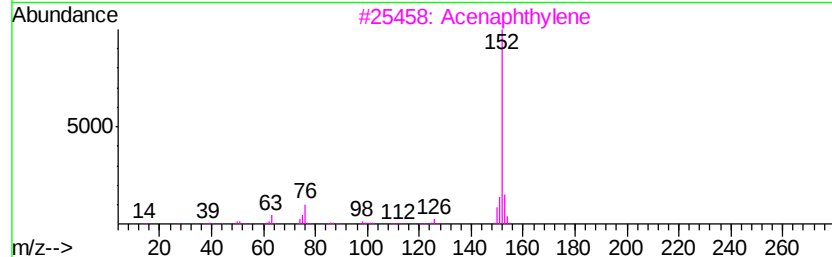
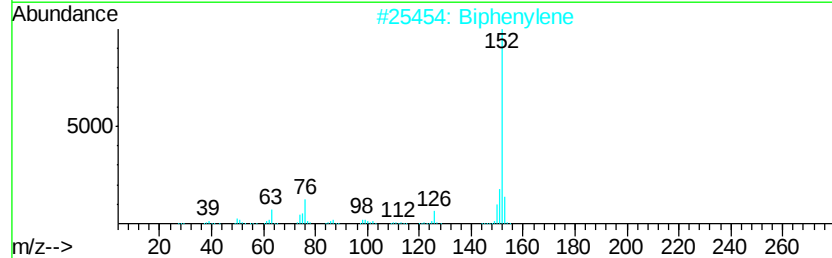
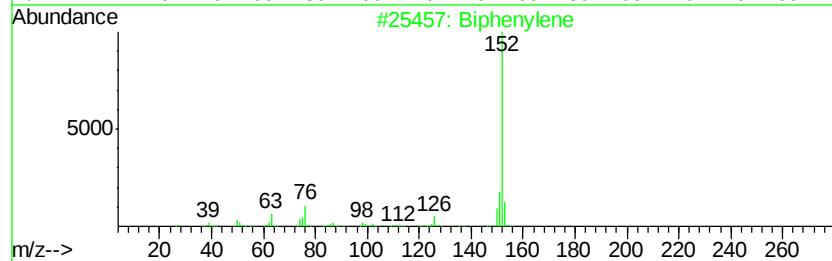
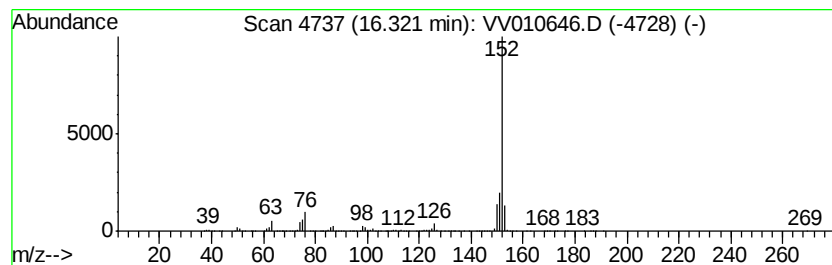
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM050219S.M
Quant Title : VOC Analysis

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

Peak Number 32 Biphenylene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.32	225.50 ug/L	3592810	1,4-Dichlorobenzene-d4	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Biphenylene	152	C12H8	000259-79-0	93
2		Biphenylene	152	C12H8	000259-79-0	91
3		Acenaphthylene	152	C12H8	000208-96-8	90
4		Acenaphthylene	152	C12H8	000208-96-8	83
5		Benzaldehyde, 2-hydroxy-3-methoxy-	152	C8H8O3	000148-53-8	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV050219\
 Data File : VV010646.D
 Acq On : 03 May 2019 03:51
 Operator : SY/MD
 Sample : K2603-20
 Misc : 5.00G/10ML/MSVOA_V/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 GAJ72

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOM2VLM050219S.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-...	10.49	3.3	ug/L	53218	3	11.29	398311	25.0
Benzene, 1,2,4-tr...	10.58	8.3	ug/L	131890	3	11.29	398311	25.0
Benzene, 1,2,3-tr...	10.96	12.0	ug/L	191203	3	11.29	398311	25.0
Benzene, 1-etheny...	11.03	54.9	ug/L	874765	3	11.29	398311	25.0
Benzene, 1-propenyl-	11.44	5.3	ug/L	84172	3	11.29	398311	25.0
Indane	11.58	25.6	ug/L	408185	3	11.29	398311	25.0
Indene	11.79	295.7	ug/L	4711080	3	11.29	398311	25.0
Benzene, 1-methyl...	11.98	15.8	ug/L	251002	3	11.29	398311	25.0
Bicyclo[4.2.0]oct...	12.04	12.8	ug/L	203566	3	11.29	398311	25.0
Benzene, (2-methy...	12.24	80.5	ug/L	1283130	3	11.29	398311	25.0
Benzene, 1,2,3,4-...	12.46	22.1	ug/L	352535	3	11.29	398311	25.0
Benzene, 1,2,3,5-...	12.51	96.4	ug/L	1535190	3	11.29	398311	25.0
Benzene, 4-etheny...	12.64	74.6	ug/L	1188220	3	11.29	398311	25.0
Benzene, 1-etheny...	12.77	21.4	ug/L	341526	3	11.29	398311	25.0
Benzene, 1-ethyl-...	12.93	106.1	ug/L	1690220	3	11.29	398311	25.0
1,4-Dihydronaphth...	12.99	140.4	ug/L	2237210	3	11.29	398311	25.0
Benzene, 1-butyryl-	13.10	298.8	ug/L	4759990	3	11.29	398311	25.0
1H-Indene, 1,3-di...	14.15	41.5	ug/L	661900	3	11.29	398311	25.0
Naphthalene, 1-me...	14.87	941.8	ug/L	15005000	3	11.29	398311	25.0
Naphthalene, 1,7-...	15.72	238.6	ug/L	3802090	3	11.29	398311	25.0
Naphthalene, 2,3-...	15.86	310.9	ug/L	4953990	3	11.29	398311	25.0
Naphthalene, 2-et...	15.99	50.9	ug/L	811395	3	11.29	398311	25.0
Naphthalene, 1,3-...	16.06	55.2	ug/L	879479	3	11.29	398311	25.0
Naphthalene, 1,4-...	16.23	35.4	ug/L	563211	3	11.29	398311	25.0
Biphenylene	16.32	225.5	ug/L	3592810	3	11.29	398311	25.0