

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW121219\
 Data File : VW014326.D
 Acq On : 12 Dec 2019 15:12
 Operator : SY/VA
 Sample : K6236-14
 Misc : 5.78G/10ML/MSVOA W/SOIL
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 BFEW1

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_W\METHOD\SOM2WLM120519S.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	1.959	4	9	20	rVV3	15536	36267	0.75%	0.095%
2	2.032	20	21	26	rVB4	2325	2691	0.06%	0.007%
3	2.178	36	45	49	rBV3	26294	72175	1.50%	0.190%
4	2.349	64	73	91	rBV	390314	751091	15.57%	1.975%
5	2.562	106	108	109	rBV2	3333	2327	0.05%	0.006%
6	2.623	113	118	126	rVV2	11574	29706	0.62%	0.078%
7	2.885	151	161	175	rBV	262867	637954	13.23%	1.678%
8	3.007	179	181	183	rBV3	4480	4582	0.10%	0.012%
9	3.385	234	243	249	rBV9	3779	11455	0.24%	0.030%
10	3.471	255	257	260	rBV4	984	1173	0.02%	0.003%
11	3.532	264	267	271	rVV5	1506	2708	0.06%	0.007%
12	3.574	272	274	276	rVV3	1276	1281	0.03%	0.003%
13	3.672	287	290	296	rVB8	1596	3242	0.07%	0.009%
14	3.830	312	316	317	rVV4	920	1037	0.02%	0.003%
15	3.855	317	320	323	rVB4	1072	1386	0.03%	0.004%
16	4.013	332	346	358	rBV	519754	1656471	34.35%	4.356%
17	4.117	358	363	376	rVV	55319	153176	3.18%	0.403%
18	4.202	376	377	380	rVB3	1740	1038	0.02%	0.003%
19	4.294	391	392	396	rVB3	1182	1155	0.02%	0.003%
20	4.385	397	407	416	rVV4	9694	29294	0.61%	0.077%
21	4.501	424	426	432	rVB4	1116	2198	0.05%	0.006%
22	4.556	432	435	438	rVB4	1372	1695	0.04%	0.004%
23	4.726	459	463	465	rBV3	980	1526	0.03%	0.004%
24	4.909	481	493	511	rBV2	85917	326738	6.78%	0.859%
25	5.135	525	530	531	rBV4	1718	2389	0.05%	0.006%
26	5.159	531	534	546	rVB6	2846	7358	0.15%	0.019%
27	5.409	570	575	576	rBV4	1412	1696	0.04%	0.004%
28	5.580	598	603	606	rBV4	1028	1470	0.03%	0.004%
29	5.714	620	625	627	rBV4	1092	1861	0.04%	0.005%
30	5.757	627	632	633	rBV4	2417	3226	0.07%	0.008%
31	5.775	633	635	645	rVB9	2797	6465	0.13%	0.017%
32	5.934	650	661	669	rBV3	19659	57501	1.19%	0.151%
33	6.055	678	681	683	rBV4	892	1102	0.02%	0.003%
34	6.275	715	717	719	rBV3	1175	1068	0.02%	0.003%

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Integration Parameters: LSCINT.P

Integrator: RTE
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Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM120519S.M
 Title : VOC Analysis

35	6.500	751	754	755	rBV2	1269	1372	0.03%	0.004%
36	6.580	764	767	769	rBV3	840	1073	0.02%	0.003%
37	6.610	769	772	775	rVB5	1796	1628	0.03%	0.004%
38	6.641	775	777	779	rBV3	1266	1082	0.02%	0.003%
39	6.714	788	789	791	rBV2	1334	1248	0.03%	0.003%
40	6.793	799	802	803	rBV2	1493	1436	0.03%	0.004%
41	6.860	812	813	815	rBV2	1722	1849	0.04%	0.005%
42	6.958	826	829	831	rVB4	1699	1486	0.03%	0.004%
43	7.000	831	836	838	rVB6	1602	2806	0.06%	0.007%
44	7.074	838	848	853	rBV	142316	397310	8.24%	1.045%
45	7.128	853	857	877	rVB2	108915	352026	7.30%	0.926%
46	7.348	891	893	895	rVB3	1272	1070	0.02%	0.003%
47	7.415	901	904	905	rBV3	1997	1731	0.04%	0.005%
48	7.494	915	917	919	rBV3	2161	1958	0.04%	0.005%
49	7.519	919	921	922	rBV2	3455	2759	0.06%	0.007%
50	7.641	930	941	954	rBV	765800	1907360	39.55%	5.016%
51	7.817	967	970	971	rBV3	1944	2163	0.04%	0.006%
52	7.939	988	990	995	rVB6	3756	5590	0.12%	0.015%
53	8.104	1015	1017	1018	rBV2	2455	1133	0.02%	0.003%
54	8.147	1021	1024	1026	rBV3	2339	2942	0.06%	0.008%
55	8.171	1026	1028	1032	rVB5	2805	2585	0.05%	0.007%
56	8.268	1034	1044	1060	rBV2	1618518	4822538	100.00%	12.682%
57	8.695	1109	1114	1118	rBV8	6939	16290	0.34%	0.043%
58	8.842	1129	1138	1150	rBV	1456805	2960081	61.38%	7.784%
59	9.079	1171	1177	1182	rVB4	26189	54043	1.12%	0.142%
60	9.268	1200	1208	1220	rBV	1069179	2209781	45.82%	5.811%
61	9.537	1249	1252	1257	rVB6	12777	23687	0.49%	0.062%
62	10.030	1327	1333	1342	rBV	578781	1046198	21.69%	2.751%
63	10.323	1373	1381	1388	rBV	2098794	3707610	76.88%	9.750%
64	10.573	1417	1422	1431	rVB	339620	602096	12.49%	1.583%
65	10.701	1439	1443	1451	rVB	71477	122824	2.55%	0.323%
66	10.920	1473	1479	1486	rBV	592915	1061968	22.02%	2.793%
67	11.061	1499	1502	1507	rBV3	64205	121790	2.53%	0.320%
68	11.439	1560	1564	1570	rVB	65417	106892	2.22%	0.281%
69	11.628	1589	1595	1617	rVB	1718047	3108600	64.46%	8.175%
70	12.256	1694	1698	1704	rBV5	26615	61712	1.28%	0.162%
71	12.609	1750	1756	1761	rBV	180690	315935	6.55%	0.831%

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Integration Parameters: LSCINT.P

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72	12.688	1764	1769	1776	rVB	674412	1086892	22.54%	2.858%
73	12.865	1789	1798	1800	rBV3	137698	341996	7.09%	0.899%
74	13.304	1860	1870	1884	rBV7	196762	934035	19.37%	2.456%
75	13.457	1885	1895	1905	rVV4	207852	822604	17.06%	2.163%
76	13.554	1906	1911	1923	rVB	1160659	1887813	39.15%	4.965%
77	13.853	1954	1960	1968	rBV	1100876	1954555	40.53%	5.140%
78	14.066	1991	1995	2003	rVB2	88849	218703	4.54%	0.575%
79	14.200	2012	2017	2024	rVB5	47896	92224	1.91%	0.243%
80	14.261	2024	2027	2033	rBV6	36719	70081	1.45%	0.184%
81	14.322	2034	2037	2040	rVV5	31327	47370	0.98%	0.125%
82	14.383	2043	2047	2051	rVB3	77164	123532	2.56%	0.325%
83	14.499	2062	2066	2068	rBV4	32032	49049	1.02%	0.129%
84	14.548	2070	2074	2080	rVB7	43796	99165	2.06%	0.261%
85	14.792	2110	2114	2120	rBV3	119337	246869	5.12%	0.649%
86	14.853	2120	2124	2131	rVB4	83785	179482	3.72%	0.472%
87	14.956	2138	2141	2147	rVB5	66264	125064	2.59%	0.329%
88	15.060	2154	2158	2162	rVB8	66749	103347	2.14%	0.272%
89	15.103	2163	2165	2170	rVB2	34211	49029	1.02%	0.129%
90	15.176	2172	2177	2183	rBV2	74587	146587	3.04%	0.385%
91	15.328	2197	2202	2208	rBV	263525	418923	8.69%	1.102%
92	15.432	2214	2219	2223	rBV3	63399	136835	2.84%	0.360%
93	15.657	2248	2256	2263	rBV7	54260	176838	3.67%	0.465%
94	15.731	2264	2268	2276	rBV5	84717	185005	3.84%	0.487%
95	15.865	2285	2290	2295	rVB	134889	248111	5.14%	0.652%
96	16.169	2336	2340	2345	rBV3	82643	151711	3.15%	0.399%
97	16.285	2354	2359	2366	rVB	162383	282536	5.86%	0.743%
98	16.371	2367	2373	2379	rVV3	202402	474123	9.83%	1.247%
99	16.438	2380	2384	2396	rVB6	132012	361152	7.49%	0.950%
100	16.633	2410	2416	2417	rBV6	94277	161406	3.35%	0.424%

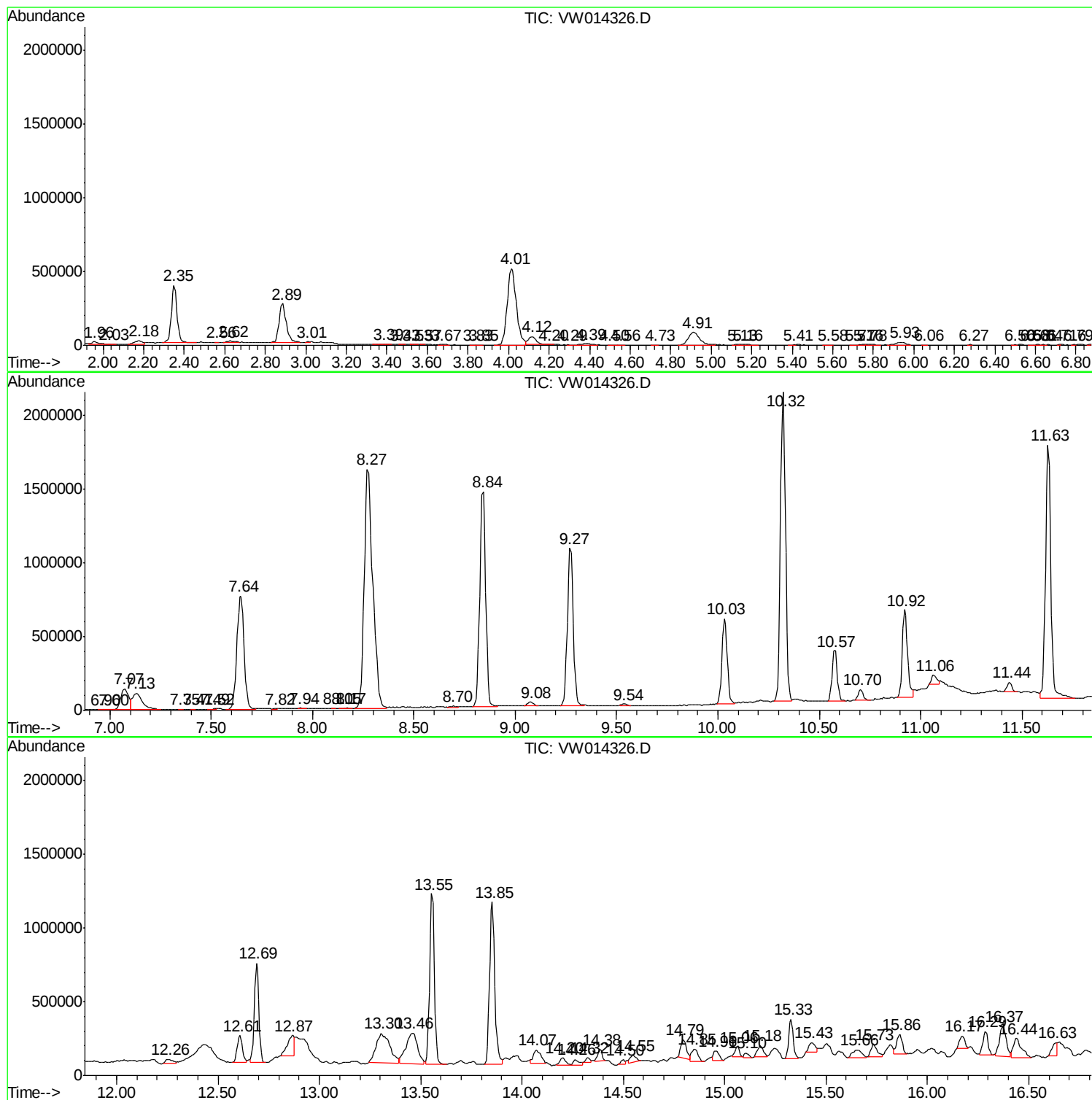
Sum of corrected areas: 38026191

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Instrument :
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ClientSampleId :
BFEW1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM120519S.M
Quant Title : VOC Analysis

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



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ALS Vial : 2 Sample Multiplier: 1

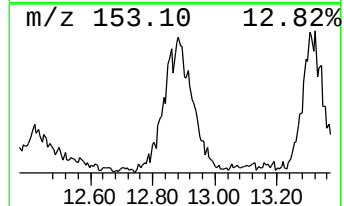
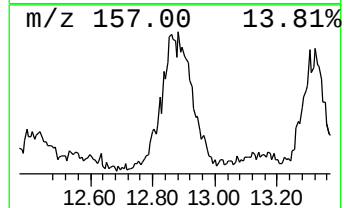
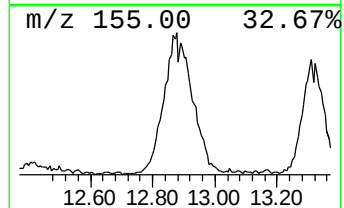
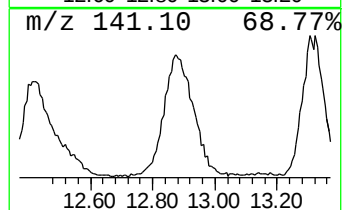
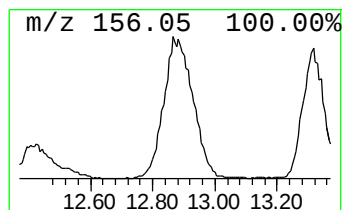
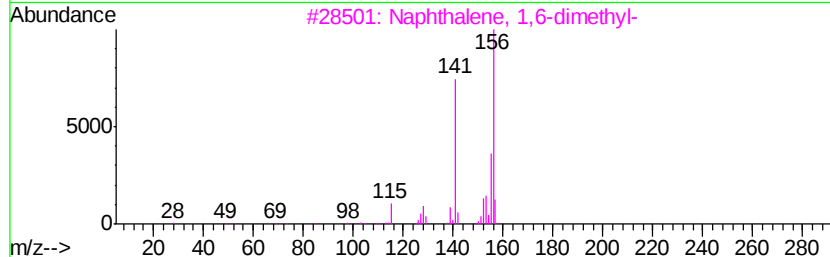
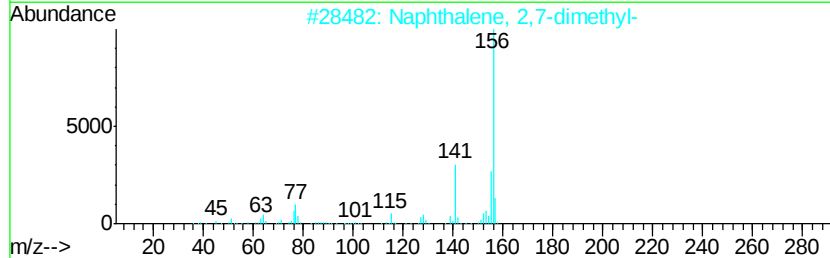
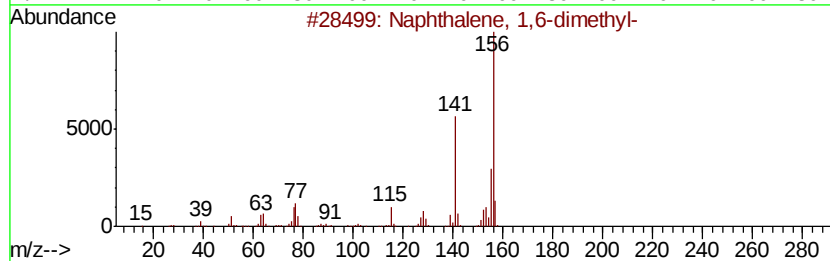
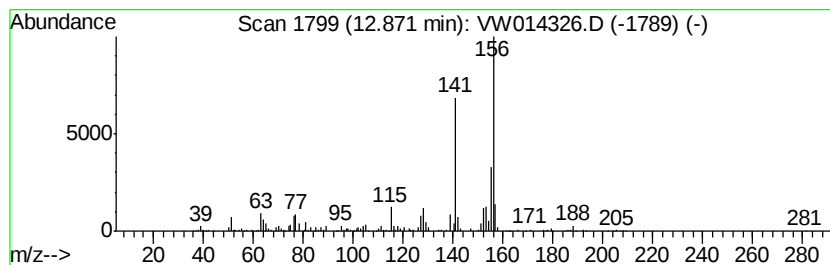
Instrument :
MSVOA_W
ClientSampleId :
BFEW1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM120519S.M
Quant Title : VOC Analysis

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

Peak Number 3 Naphthalene, 1,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.87	4.53 ug/L	341996	1,4-Dichlorobenzene-d4	13.55	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95



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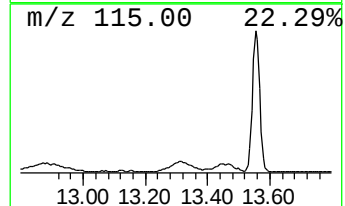
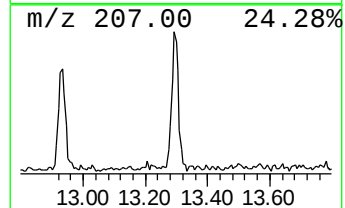
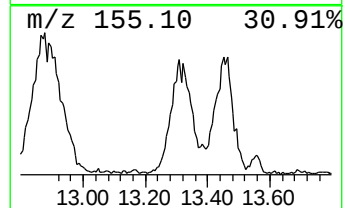
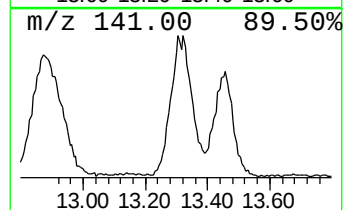
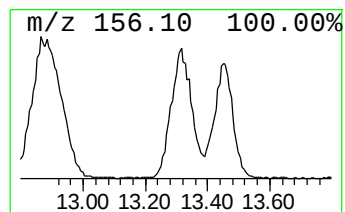
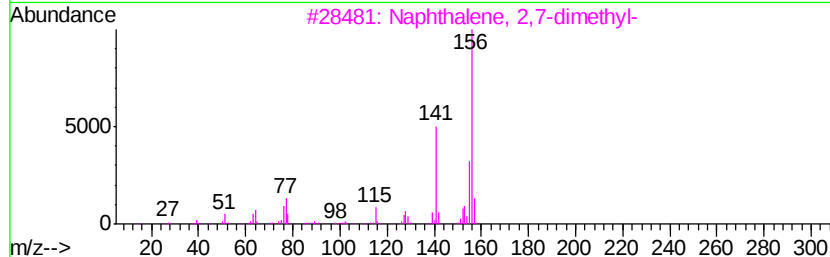
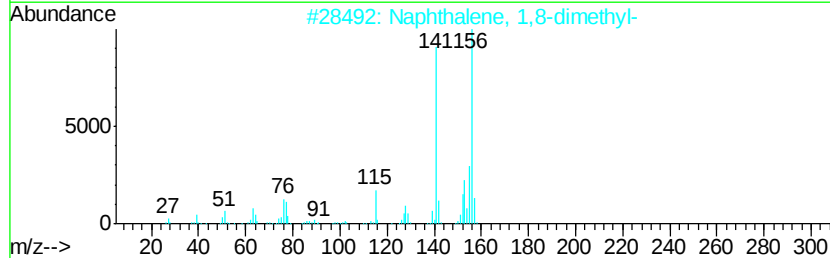
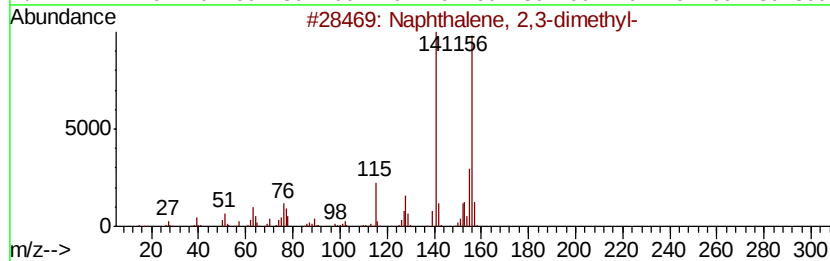
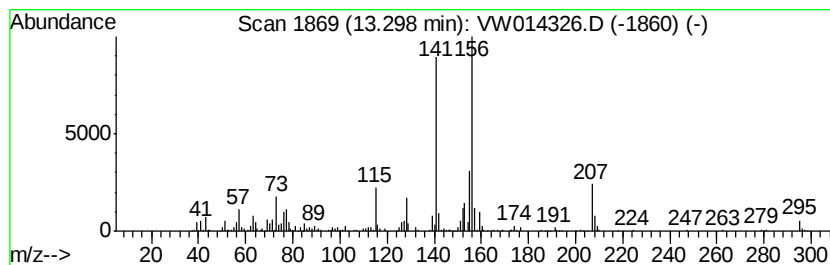
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Quant Title : VOC Analysis

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

Peak Number 4 Naphthalene, 2,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.30	12.37 ug/L	934035	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2		Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	98
3		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97



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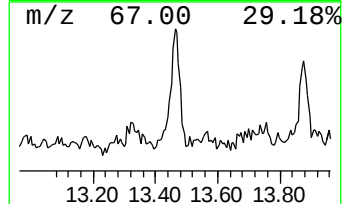
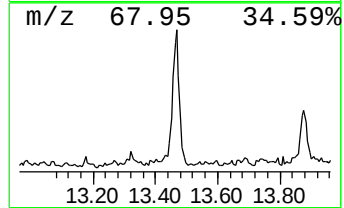
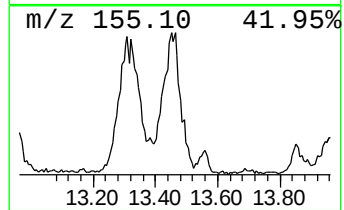
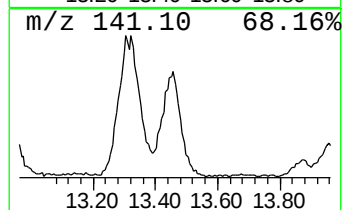
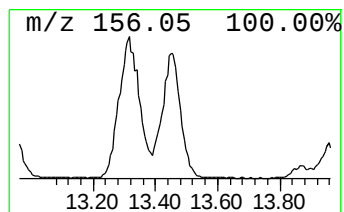
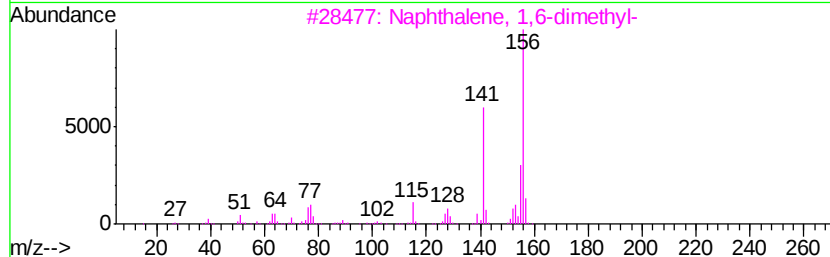
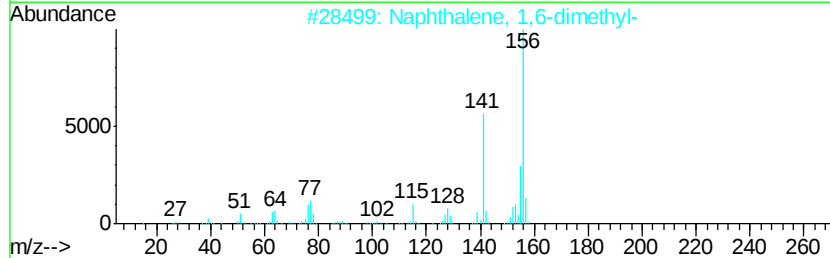
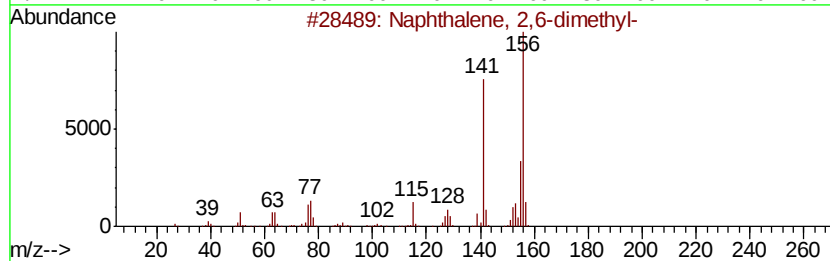
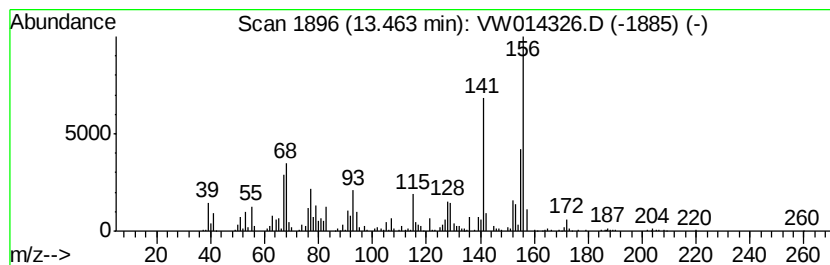
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Quant Title : VOC Analysis

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

Peak Number 5 Naphthalene, 2,6-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.46	10.89 ug/L	822604	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
2		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
4		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

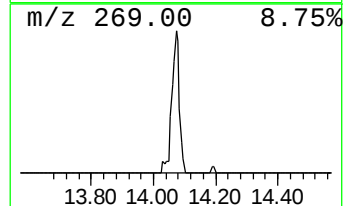
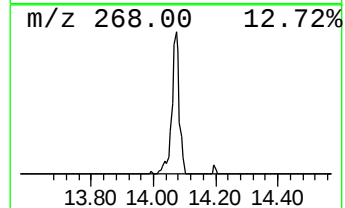
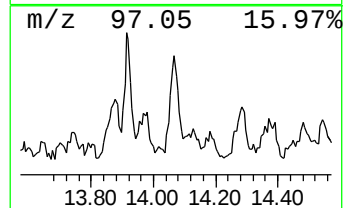
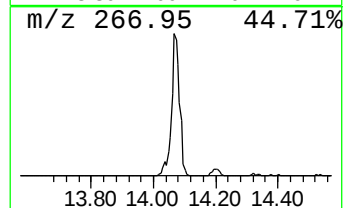
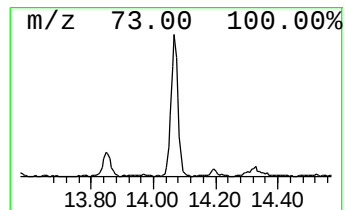
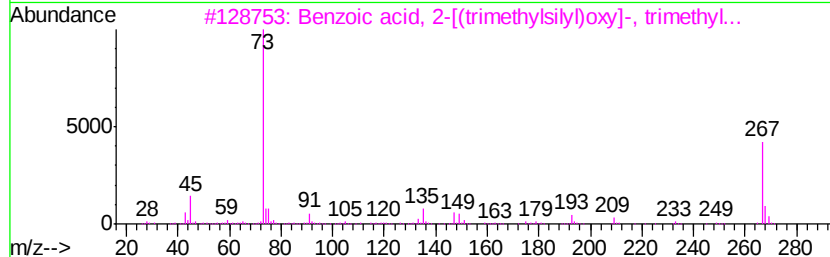
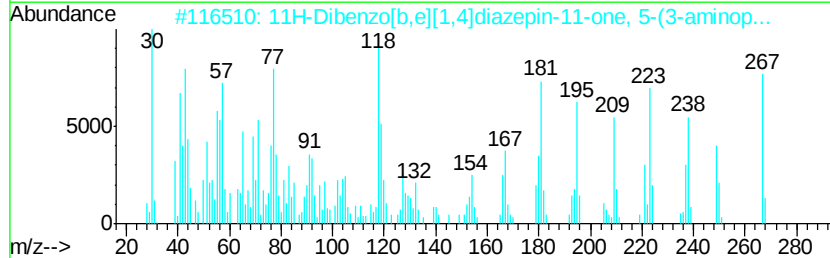
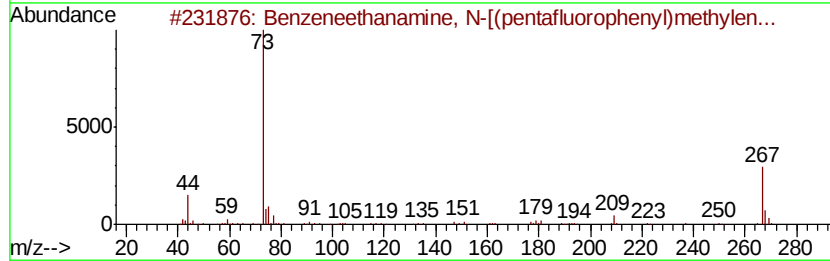
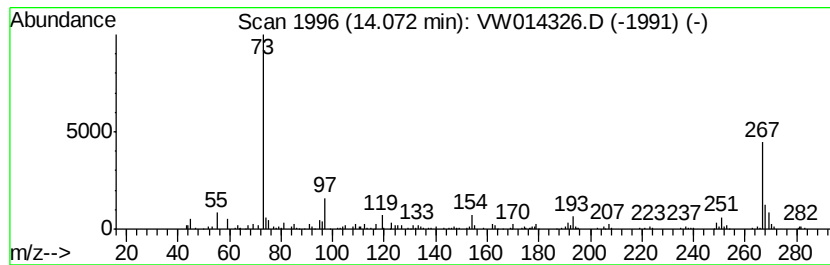
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ClientSampleId :
BFEW1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM120519S.M
Quant Title : VOC Analysis

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

Peak Number 6 unknown-01 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.07	2.90 ug/L	218703	1,4-Dichlorobenzene-d4	13.55
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzeneethanamine, N-[(pentafluoro...]	475 C21H26F5N02Si2	055429-85-1 45
2		11H-Dibenzo[b,e][1,4]diazepin-11...	267 C16H17N3O	013450-73-2 38
3		Benzoic acid, 2-[(trimethylsilyl)...]	282 C13H22O3Si2	003789-85-3 38
4		1-(3-Chlorophenyl)piperazine, 4-...	268 C13H21ClN2Si	1000373-99-3 11
5		1,3,7-Octatriene, 2-methyl-3-tri...	194 C12H22Si	1000152-32-4 10



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEW1

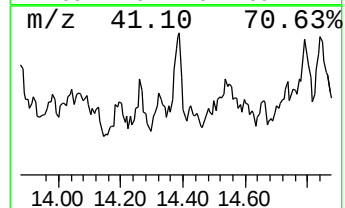
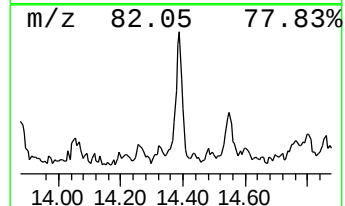
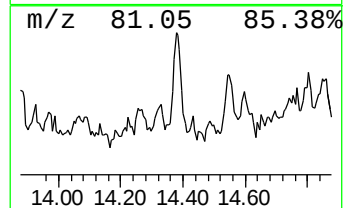
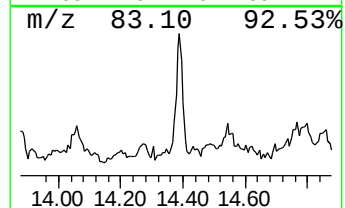
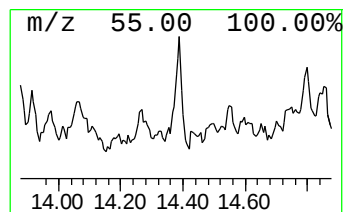
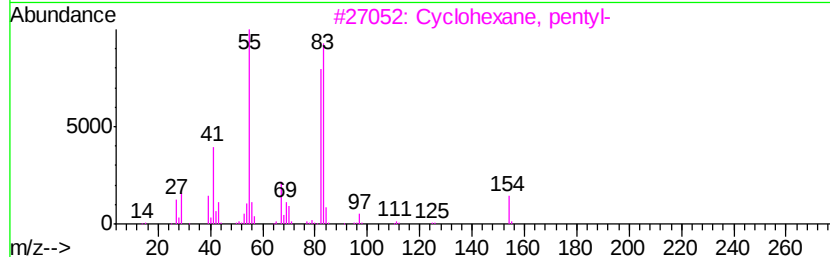
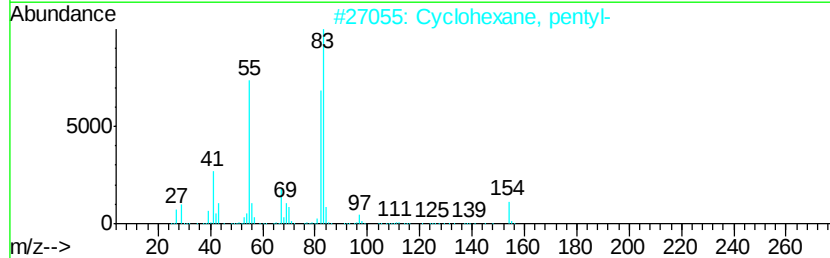
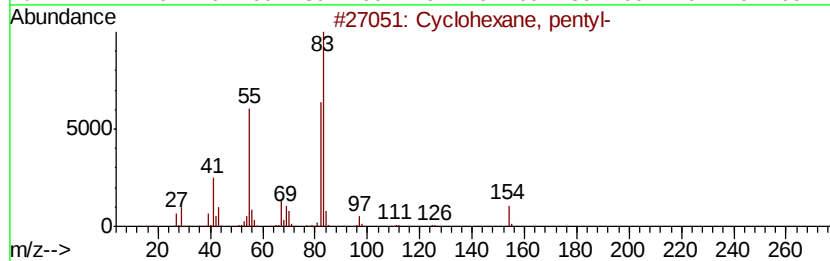
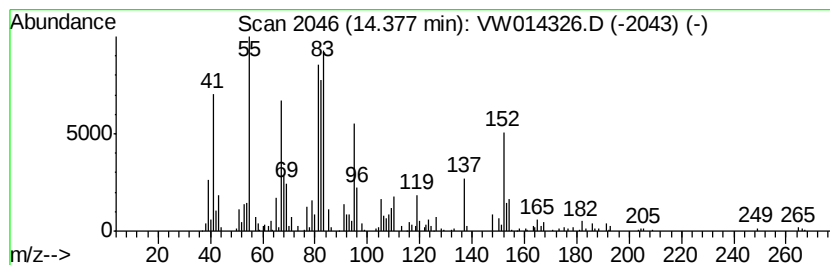
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Quant Title : VOC Analysis

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

Peak Number 7 (DEL) Alkane: Cyclic14.38 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.38	1.64 ug/L	123532	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, pentyl-	154	C11H22	004292-92-6	49
2		Cyclohexane, pentyl-	154	C11H22	004292-92-6	49
3		Cyclohexane, pentyl-	154	C11H22	004292-92-6	49
4		Cyclohexane, 1,1'-(1,4-butanediol...	222	C16H30	006165-44-2	49
5		1,1'-Bicyclohexyl	166	C12H22	000092-51-3	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEW1

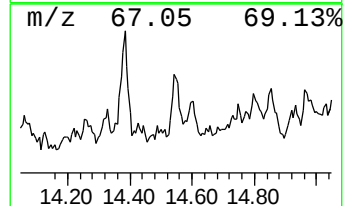
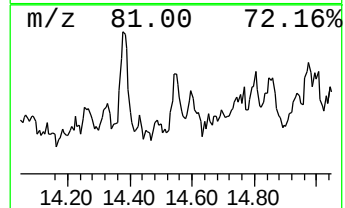
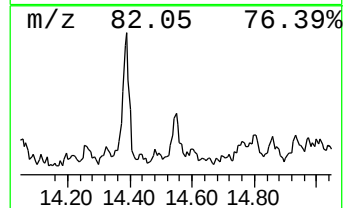
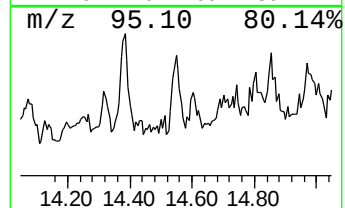
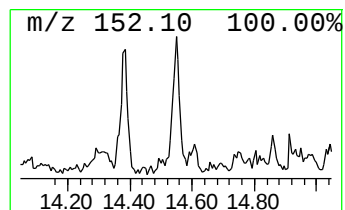
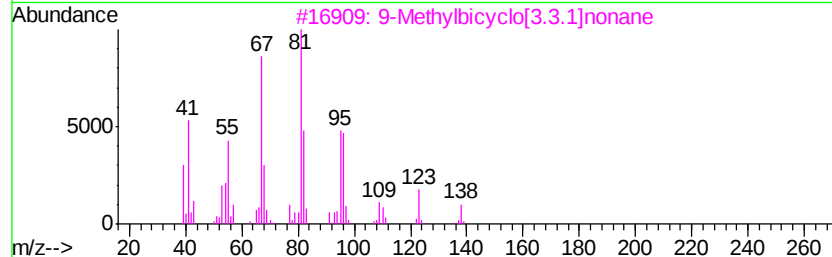
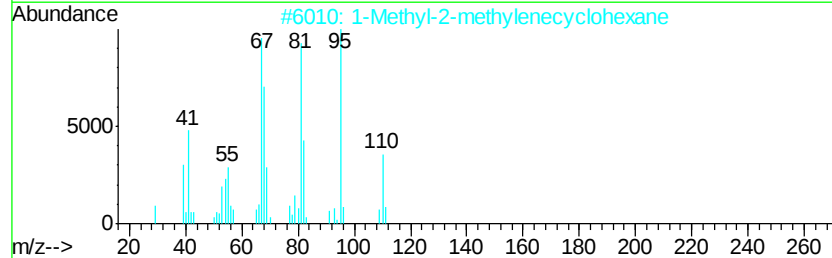
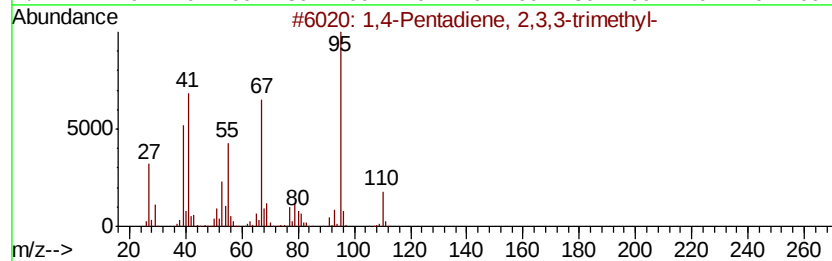
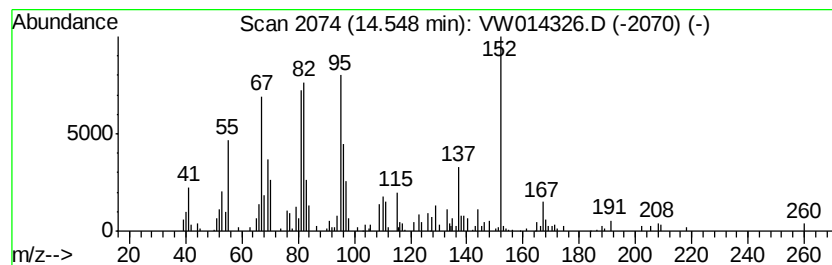
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Quant Title : VOC Analysis

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

Peak Number 8 unknown-02 Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.55	1.31 ug/L	99165	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,4-Pentadiene, 2,3,3-trimethyl-	110	C8H14	000756-02-5	35
2		1-Methyl-2-methylenecyclohexane	110	C8H14	002808-75-5	25
3		9-Methylbicyclo[3.3.1]nonane	138	C10H18	025107-01-1	25
4		cis,cis-2,8-Dimethylspiro[5.5]un...	180	C13H24	095530-75-9	22
5		Cyclobutene, 1,2,3,4-tetramethyl-	110	C8H14	090346-45-5	18



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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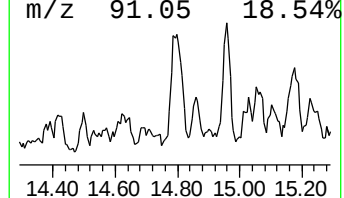
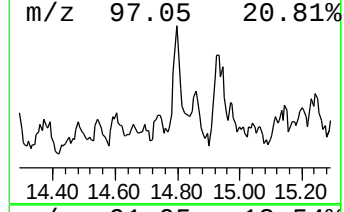
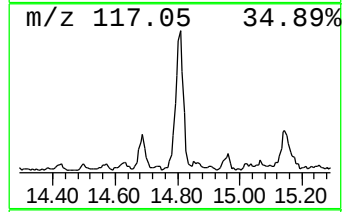
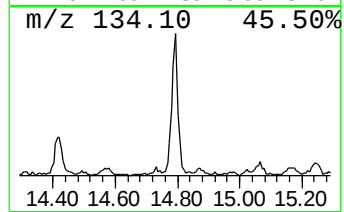
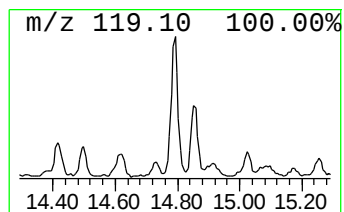
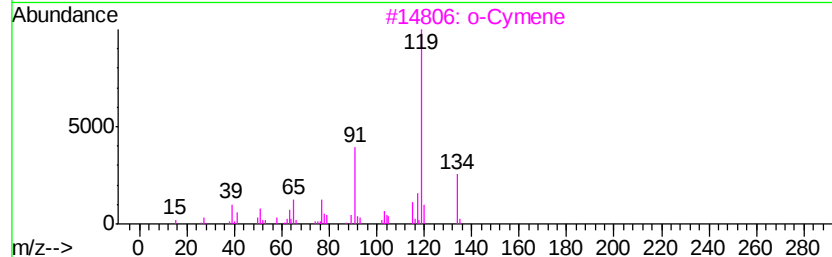
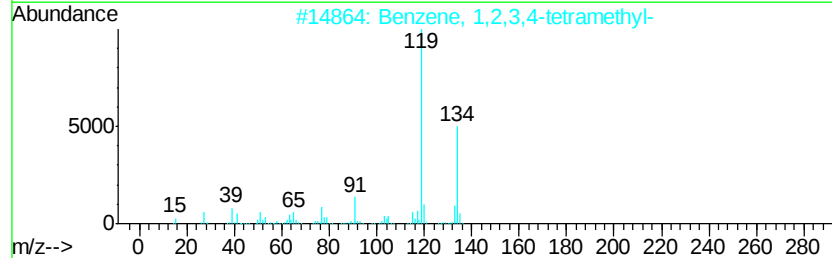
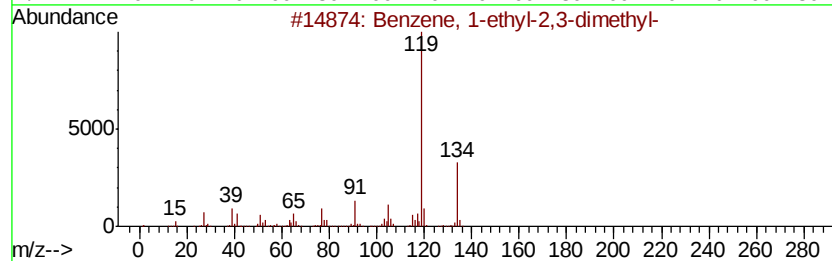
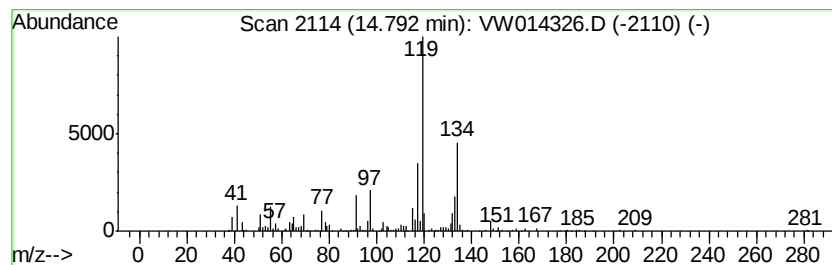
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Quant Title : VOC Analysis

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

Peak Number 9 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.79	3.27 ug/L	246869	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	70
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	70
3		o-Cymene	134	C10H14	000527-84-4	70
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	64
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEW1

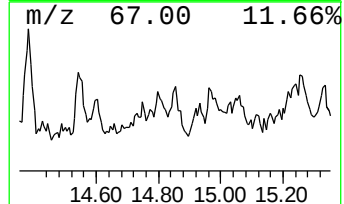
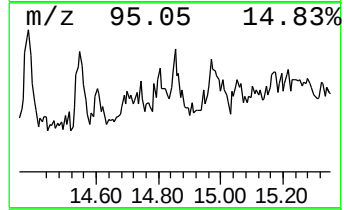
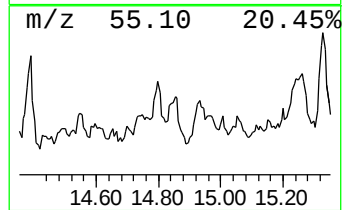
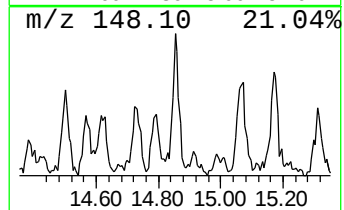
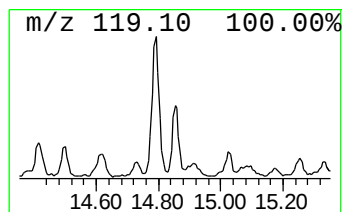
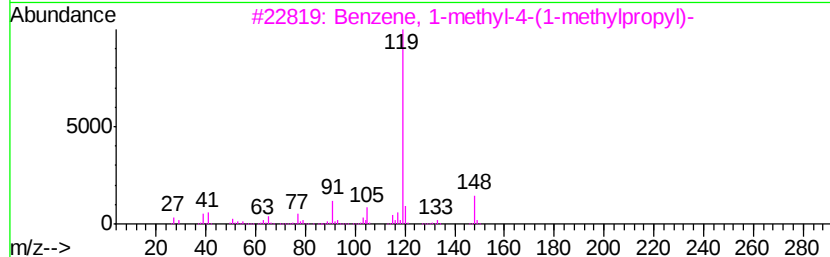
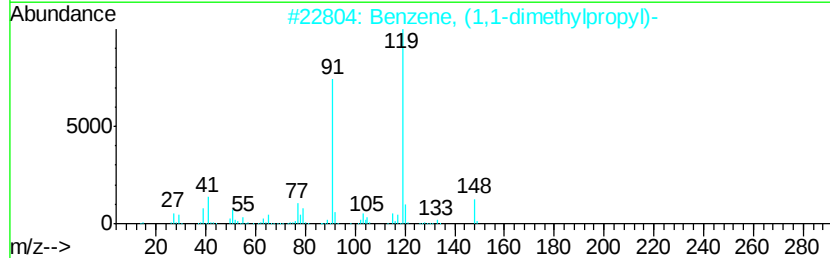
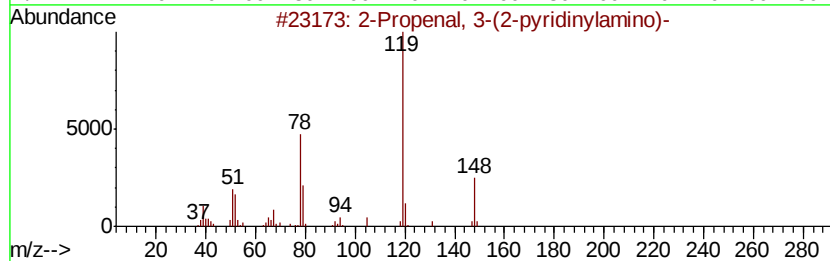
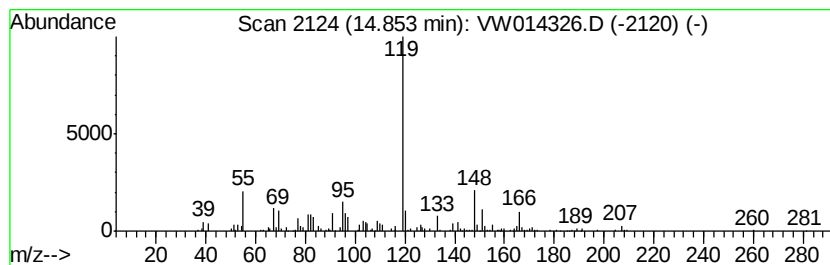
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Quant Title : VOC Analysis

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

Peak Number 10 2-Propenal, 3-(2-pyridinyla... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	2.38 ug/L	179482	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propenal, 3-(2-pyridinylamino)-	148	C8H8N2O	068970-82-1	50
2		Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	50
3		Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	50
4		Glycine, N-(m-toluoy)-, methyl e...	207	C11H13NO3	1000299-63-8	47
5		Benzene, 1,1'-(1,1,2,2-tetrameth...	238	C18H22	001889-67-4	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BFEW1

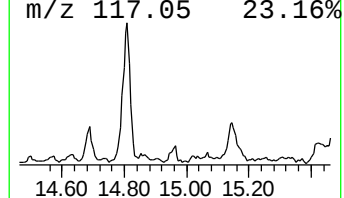
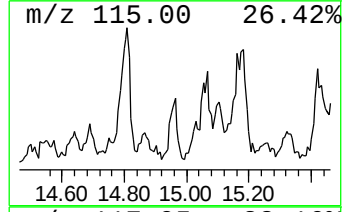
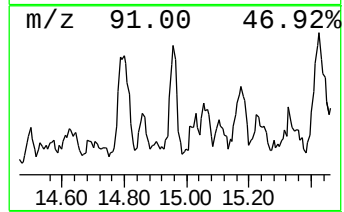
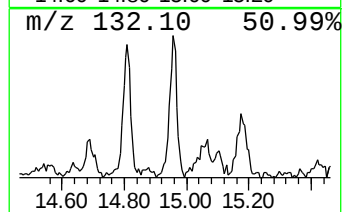
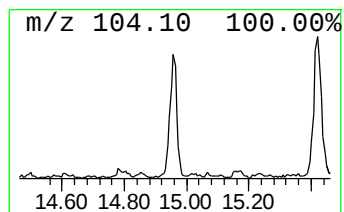
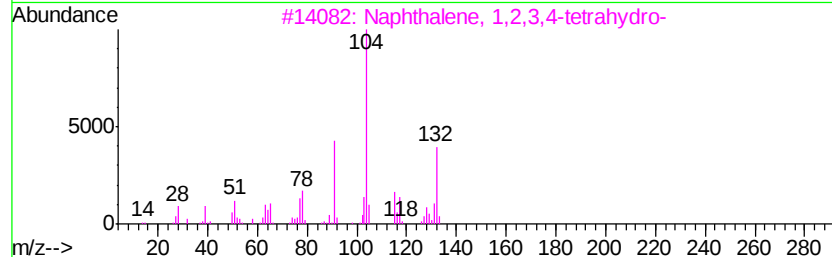
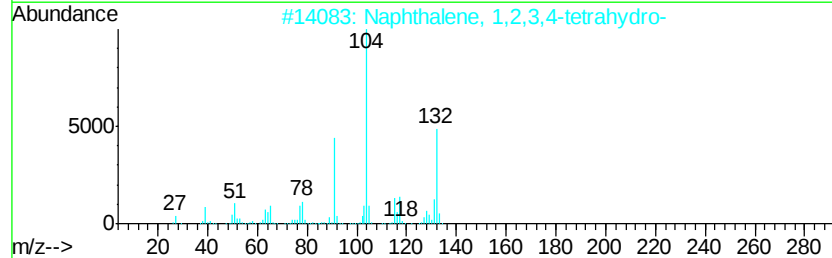
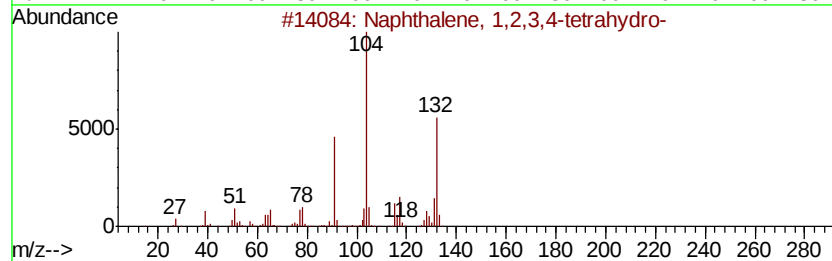
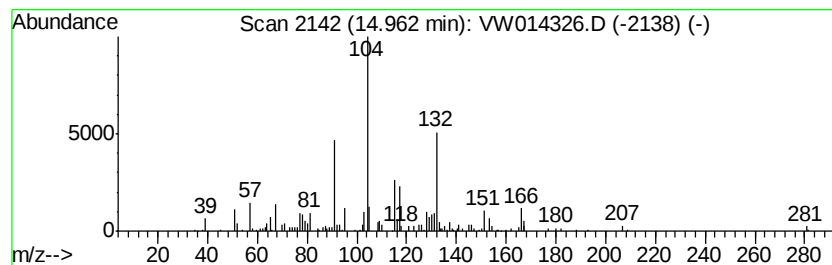
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Quant Title : VOC Analysis

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

Peak Number 11 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	1.66 ug/L	125064	1,4-Dichlorobenzene-d4	13.55

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	91
2			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	91
3			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	81
4			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	74
5			Indan, epoxide	132	C9H8O	000768-22-9	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEW1

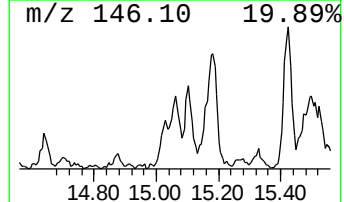
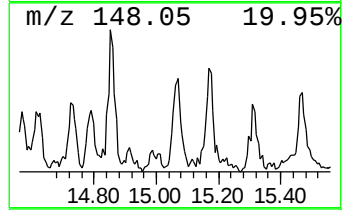
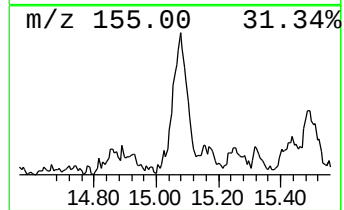
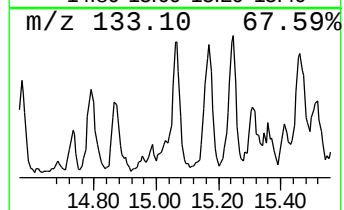
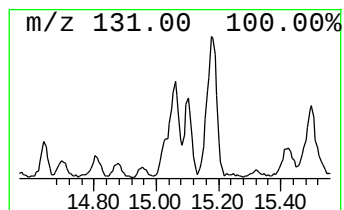
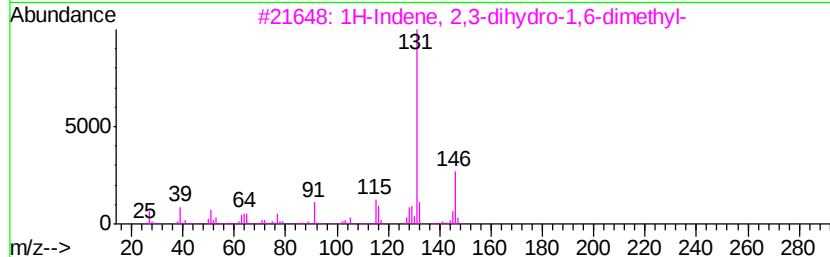
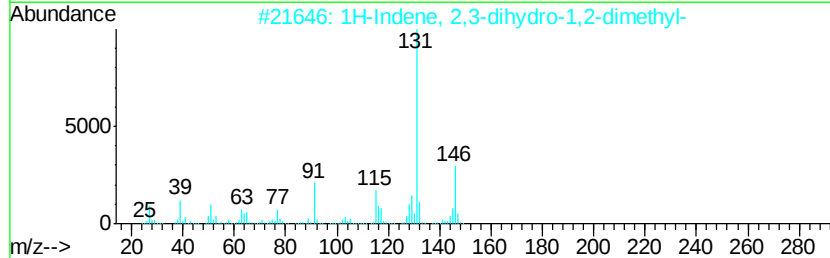
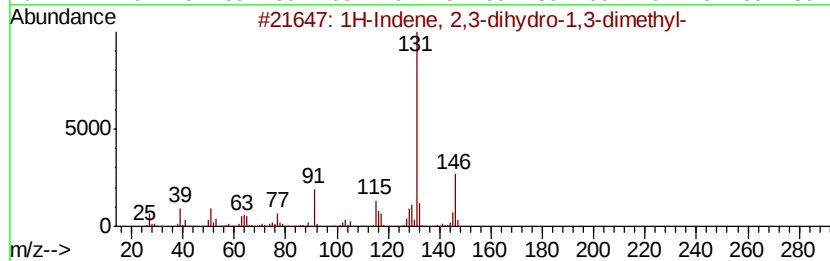
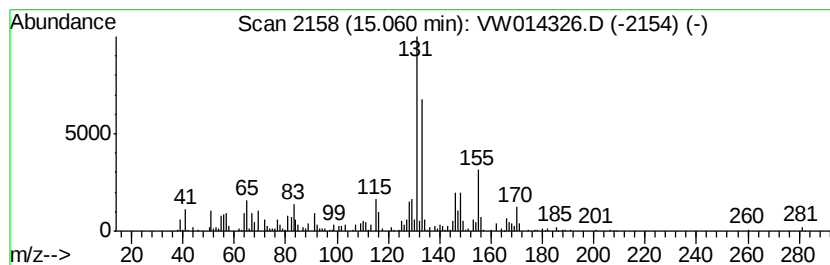
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Quant Title : VOC Analysis

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

Peak Number 12 1H-Indene, 2,3-dihydro-1,3-... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	1.37 ug/L	103347	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	55
2		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	55
3		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	49
4		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	49
5		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEW1

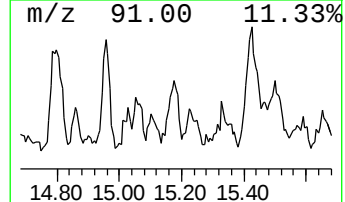
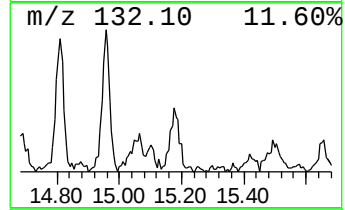
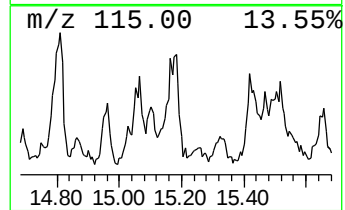
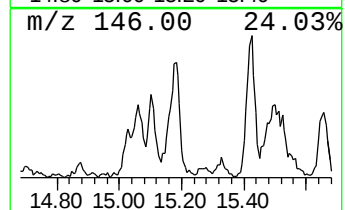
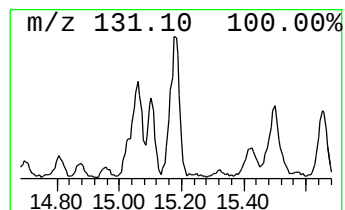
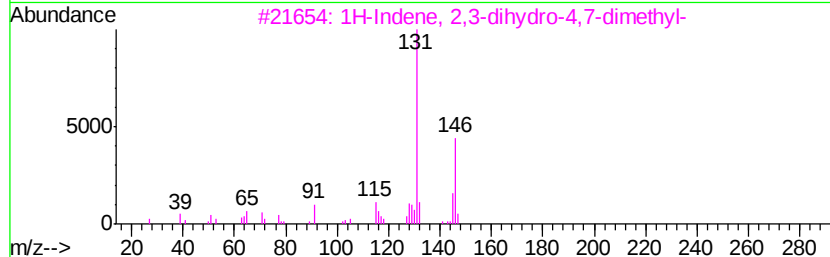
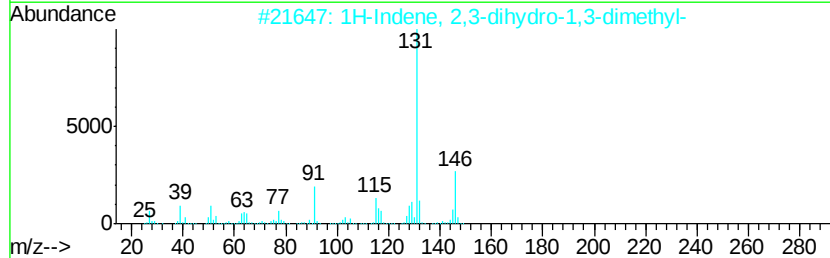
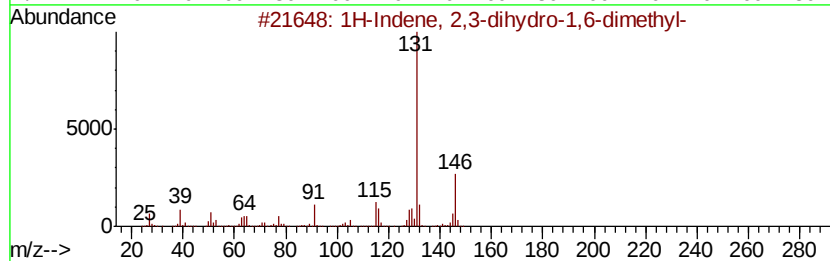
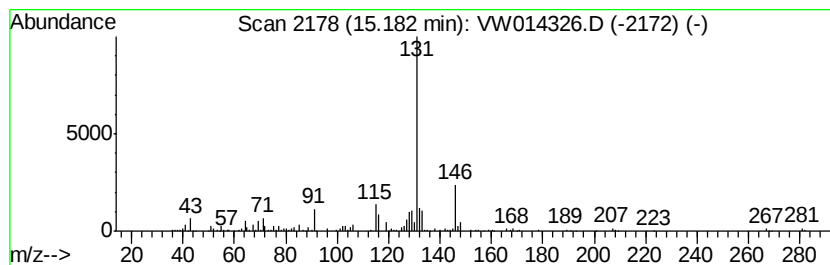
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Quant Title : VOC Analysis

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

Peak Number 13 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.18	1.94 ug/L	146587	1,4-Dichlorobenzene-d4	13.55

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90
2			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	81
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	81
5			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEW1

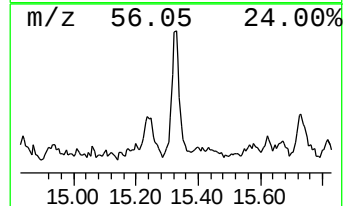
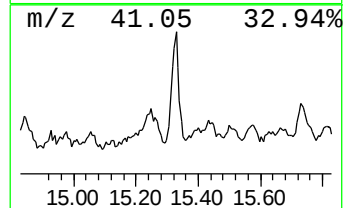
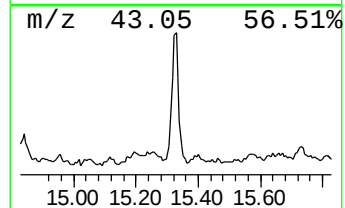
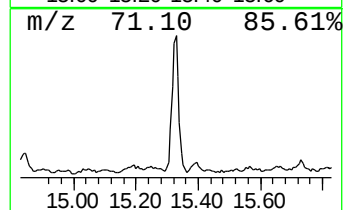
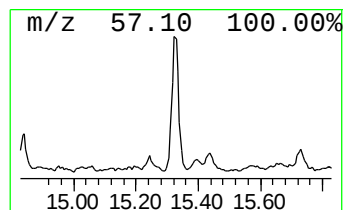
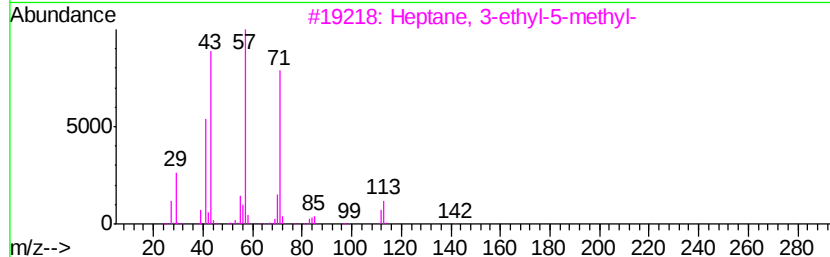
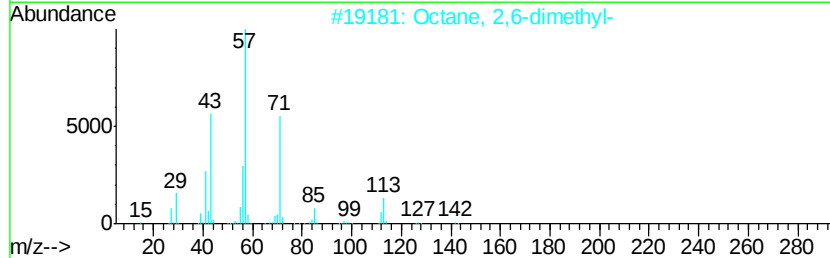
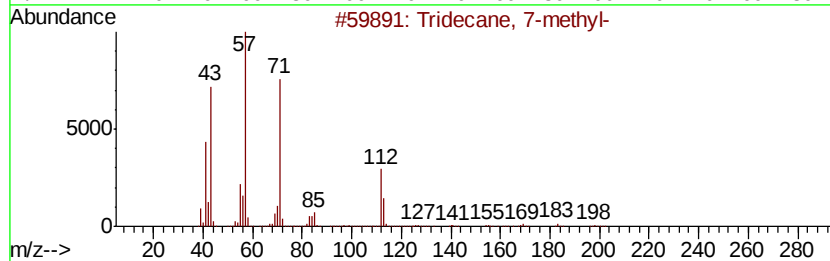
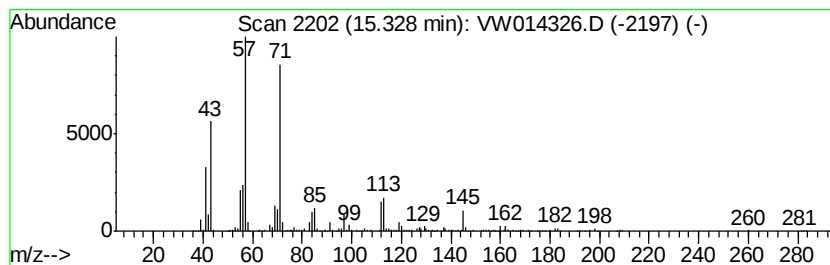
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Quant Title : VOC Analysis

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.33	5.55 ug/L	418923	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 7-methyl-	198	C14H30	026730-14-3	81
2		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
3		Heptane, 3-ethyl-5-methyl-	142	C10H22	052896-90-9	64
4		Nonane, 4-methyl-5-propyl-	184	C13H28	062185-55-1	64
5		Sulfurous acid, 2-ethylhexyl tri...	376	C21H44O3S	1000309-19-6	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEW1

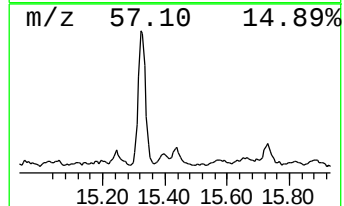
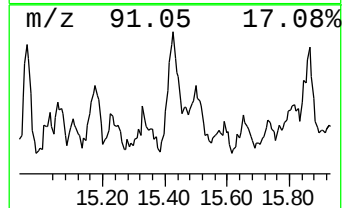
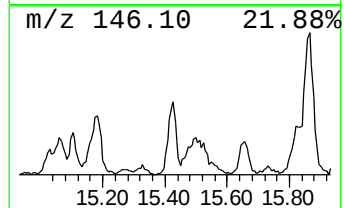
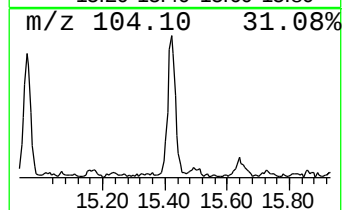
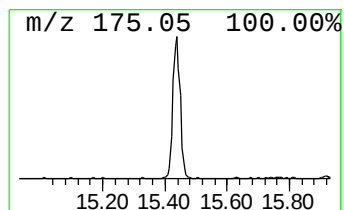
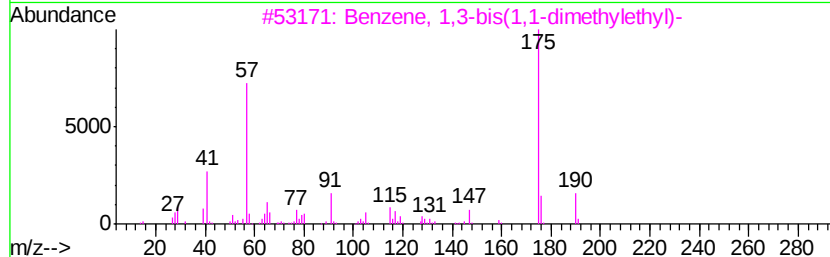
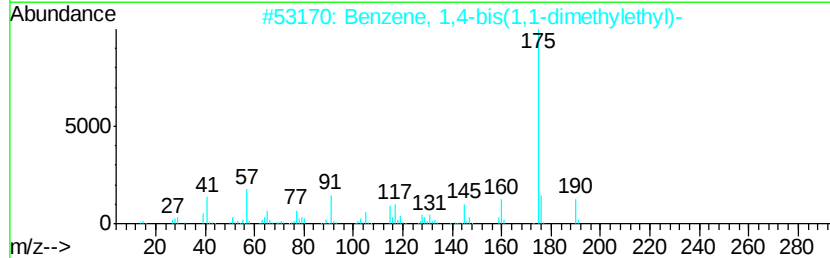
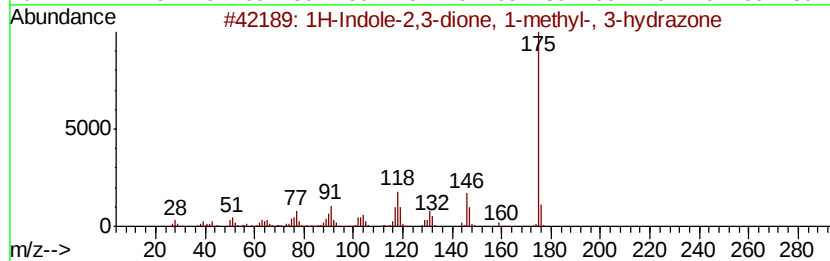
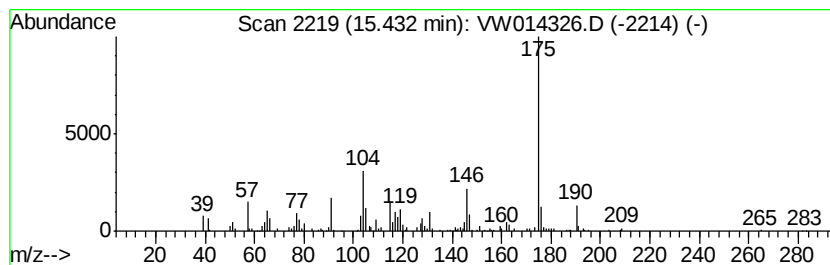
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Quant Title : VOC Analysis

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

Peak Number 15 1H-Indole-2,3-dione, 1-meth... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.43	1.81 ug/L	136835	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indole-2,3-dione, 1-methyl-, ...	175	C9H9N3O	003265-23-4	72
2		Benzene, 1,4-bis(1,1-dimethyleth...	190	C14H22	001012-72-2	64
3		Benzene, 1,3-bis(1,1-dimethyleth...	190	C14H22	001014-60-4	60
4		Benzene, 1,2,4-tripropyl-	204	C15H24	041898-97-9	50
5		2,2'-Ethylidenebis(5-methylfuran)	190	C12H14O2	003209-79-8	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEW1

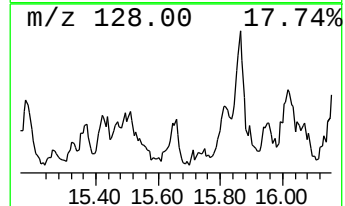
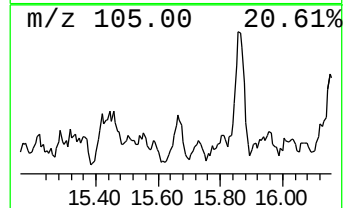
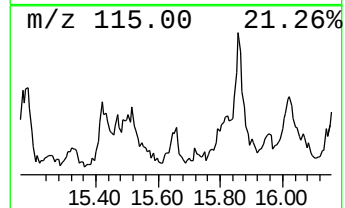
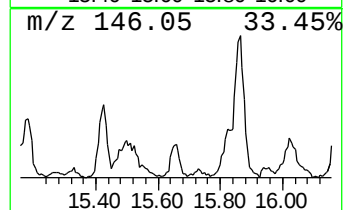
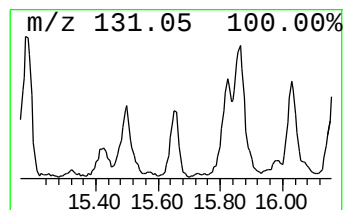
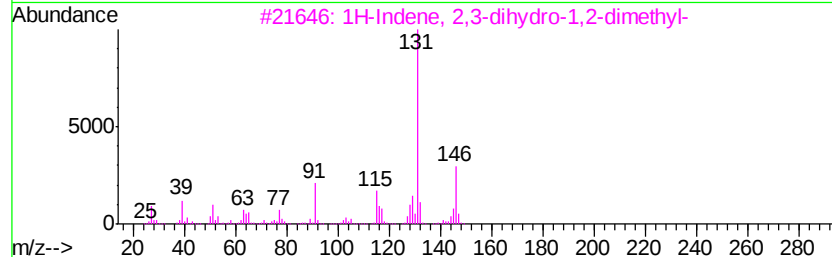
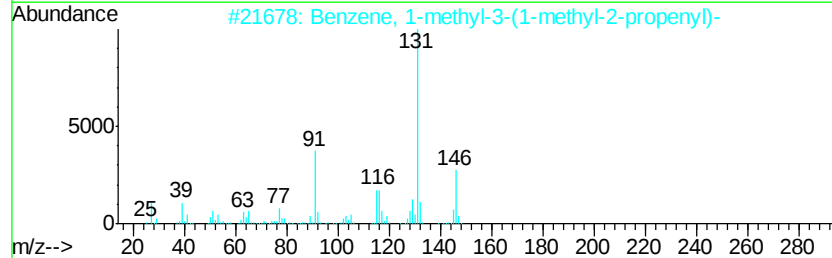
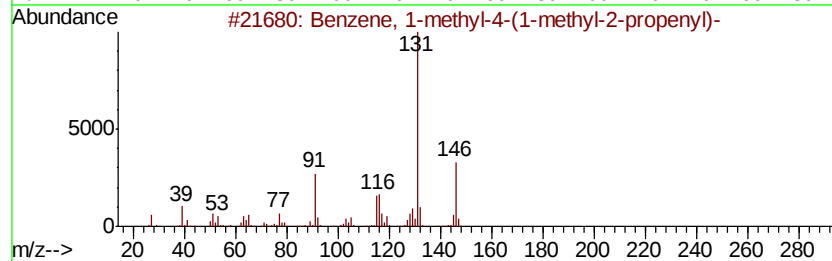
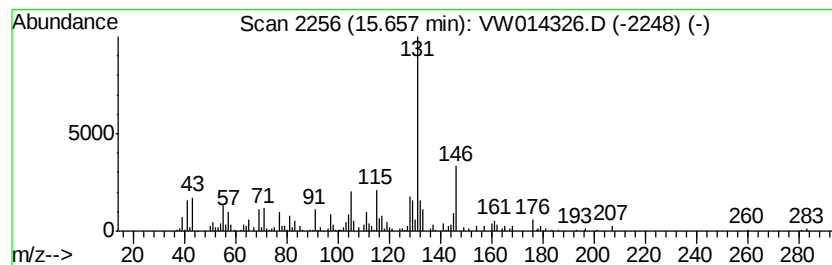
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Quant Title : VOC Analysis

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

Peak Number 16 Benzene, 1-methyl-4-(1-meth... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	2.34 ug/L	176838	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	76
2		Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	76
3		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	70
4		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	68
5		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	68



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

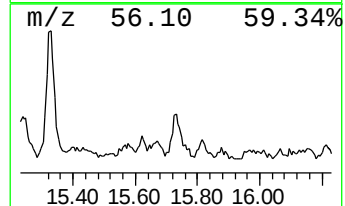
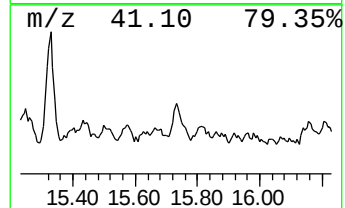
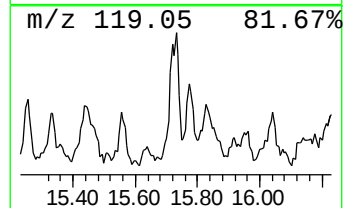
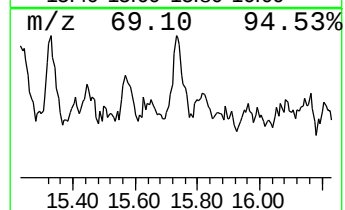
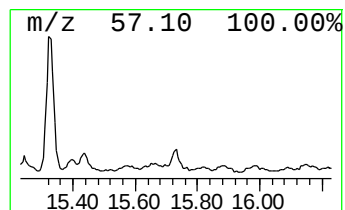
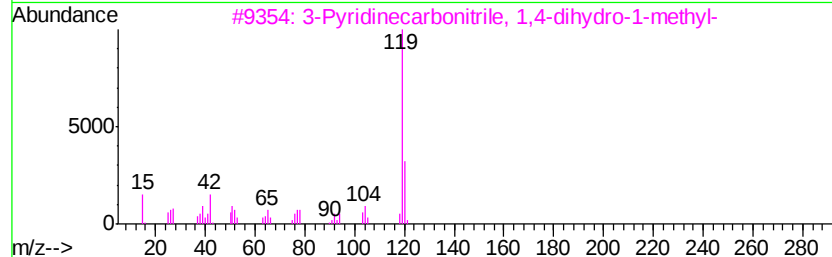
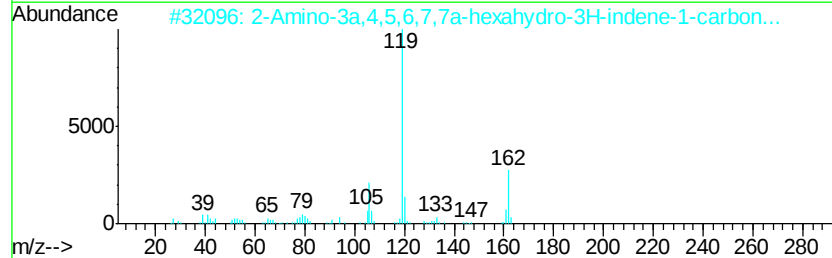
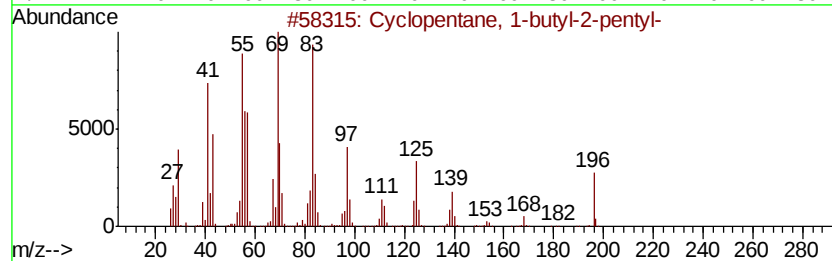
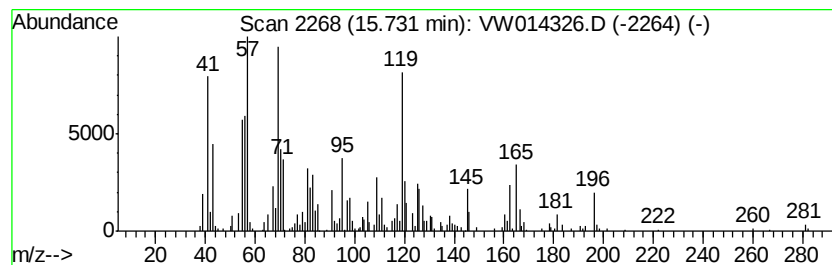
Instrument :
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ClientSampleId :
BFEW1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM120519S.M
Quant Title : VOC Analysis

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

Peak Number 17 (DEL) Alkane: Cyclic15.73 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.73	2.45 ug/L	185005	1,4-Dichlorobenzene-d4	13.55
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclopentane, 1-butyl-2-pentyl-	196 C14H28	061142-52-7 25
2		2-Amino-3a,4,5,6,7,7a-hexahydro-...	162 C10H14N2	1000185-29-8 22
3		3-Pyridinecarbonitrile, 1,4-dihv...	120 C7H8N2	019424-15-8 20
4		1-Heptanol, 2,4-diethyl-	172 C11H24O	080192-55-8 18
5		Cyclopentane, iodo-	196 C5H9I	001556-18-9 18



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEW1

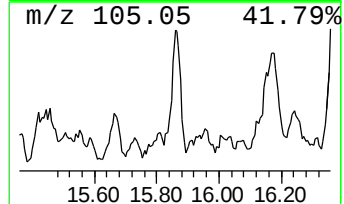
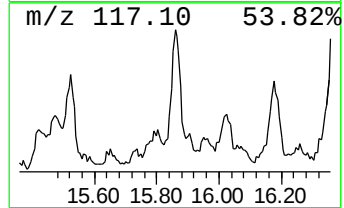
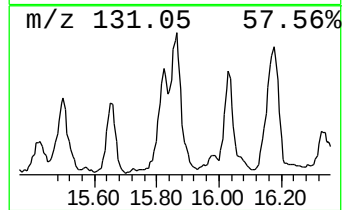
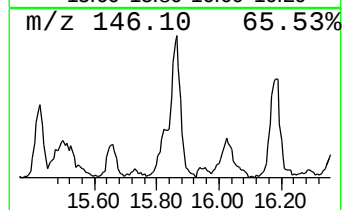
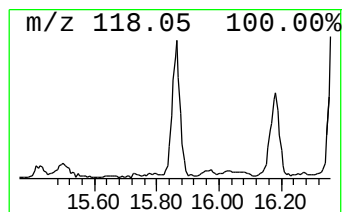
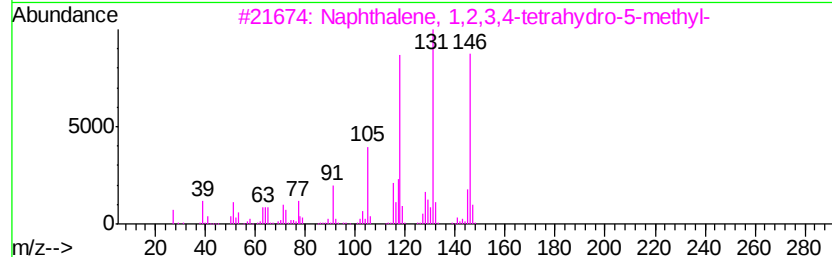
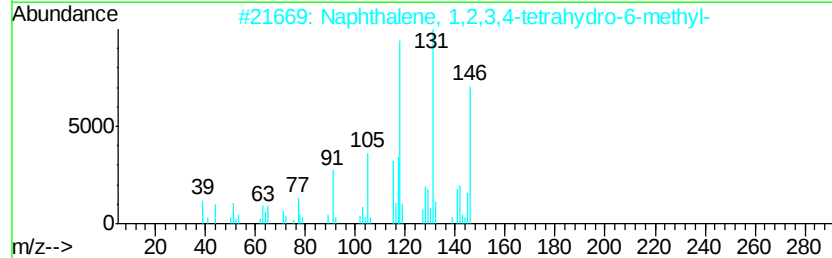
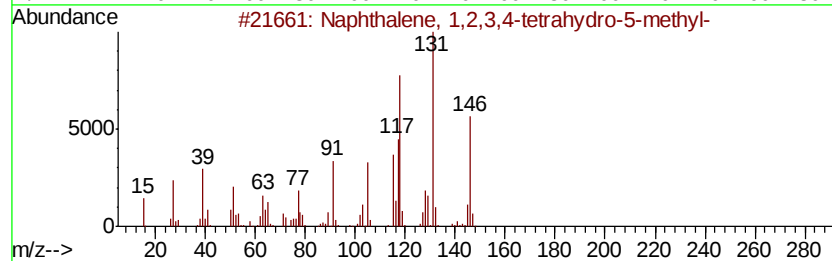
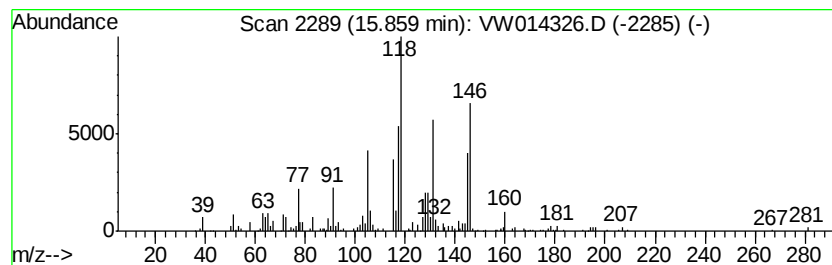
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Quant Title : VOC Analysis

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

Peak Number 18 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.86	3.29 ug/L	248111	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	68
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	62
3		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	58
4		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	52
5		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	52



Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEW1

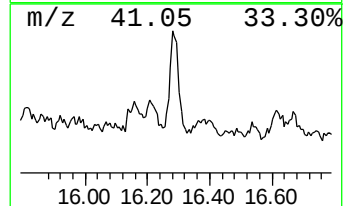
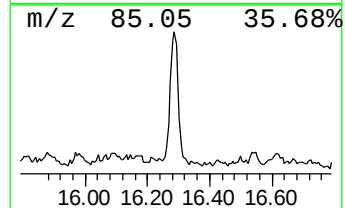
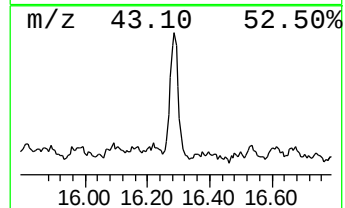
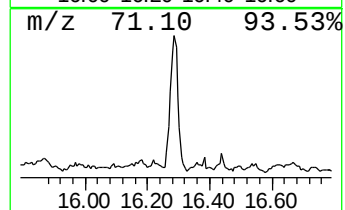
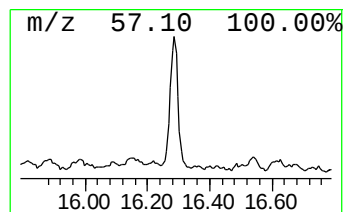
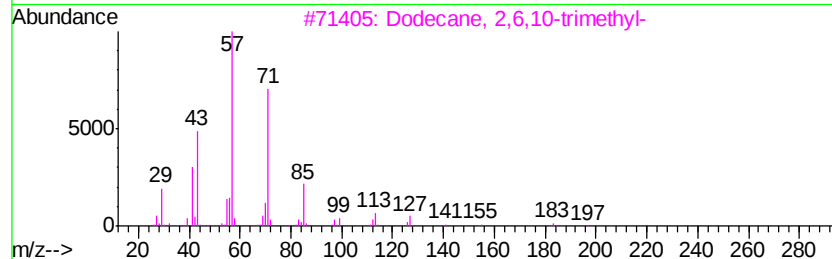
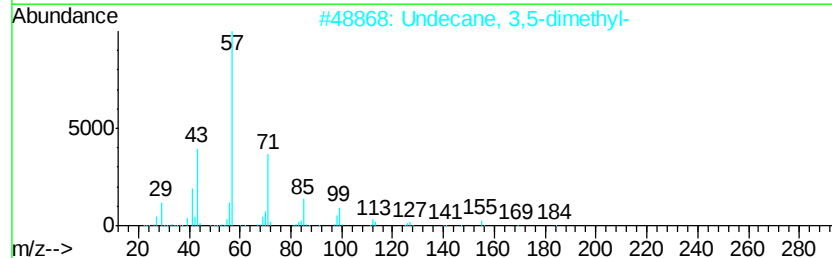
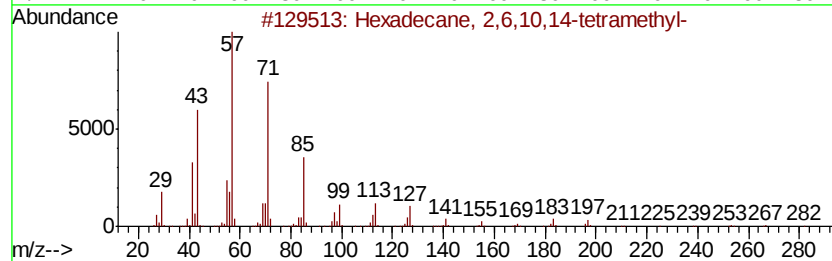
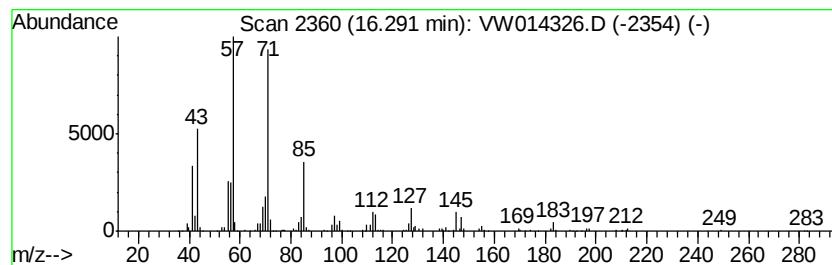
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Quant Title : VOC Analysis

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.29	3.74 ug/L	282536	1,4-Dichlorobenzene-d4	13.55

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	86
2			Undecane, 3,5-dimethyl-	184	C13H28	017312-81-1	68
3			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	64
4			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
5			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEW1

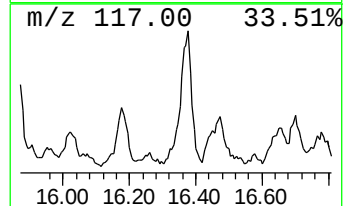
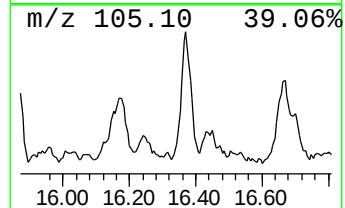
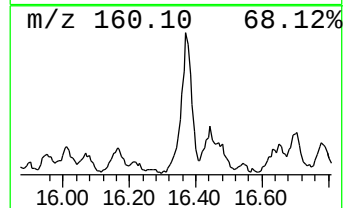
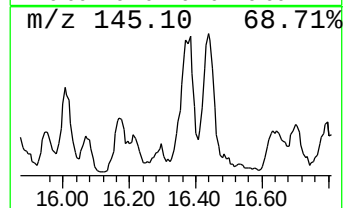
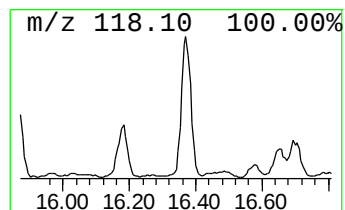
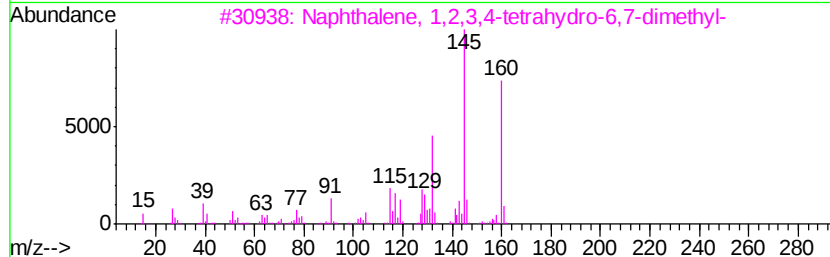
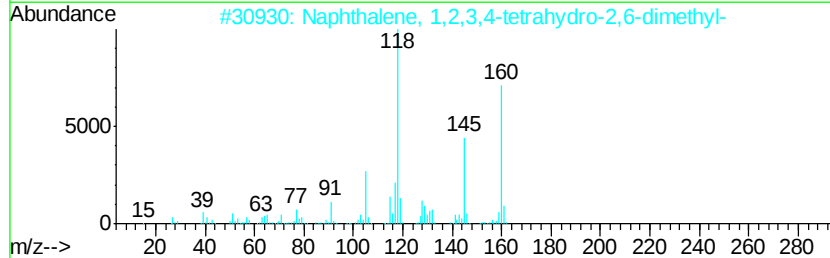
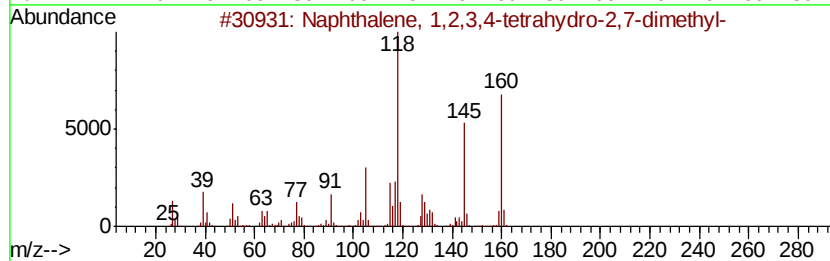
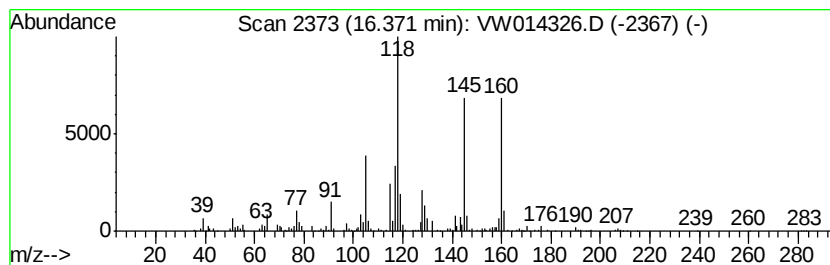
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Quant Title : VOC Analysis

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

Peak Number 21 Naphthalene, 1,2,3,4-tetra... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.37	6.28 ug/L	474123	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	94
2		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	81
3		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	78
4		1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	014944-23-1	74
5		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	68



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEW1

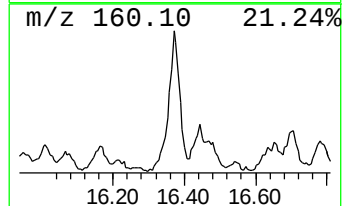
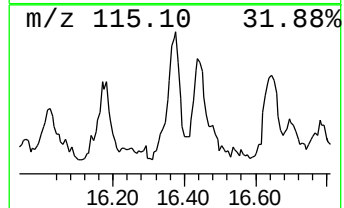
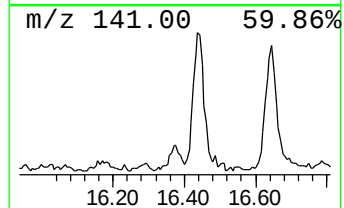
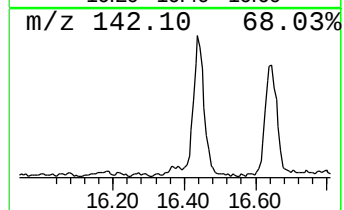
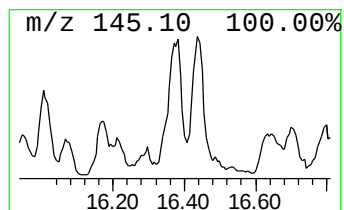
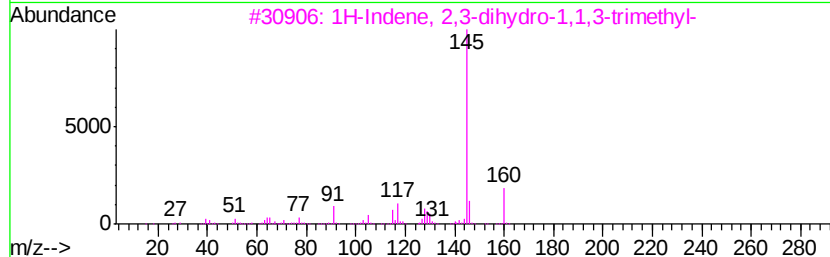
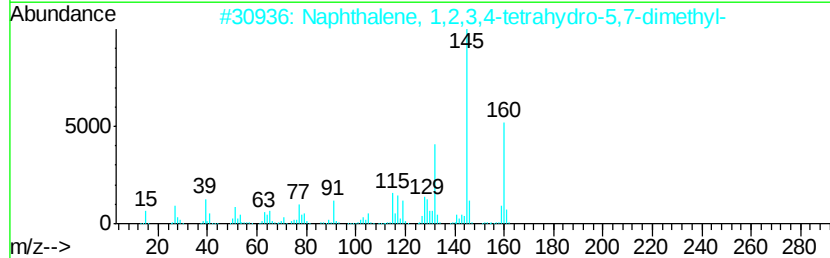
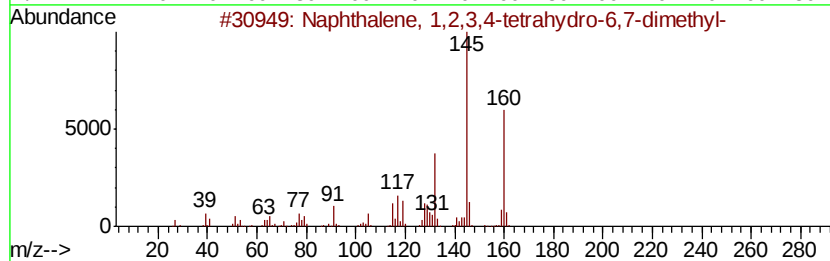
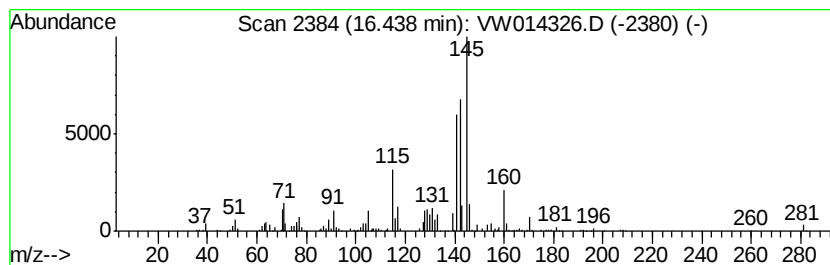
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Quant Title : VOC Analysis

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

Peak Number 22 Naphthalene, 1,2,3,4-tetra... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.44	4.78 ug/L	361152	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	53
2		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021693-54-9	53
3		1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	50
4		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001985-59-7	50
5		Benzene, (2,2-dimethyl-1-methyle...	160	C12H16	005676-29-9	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121219\
Data File : VW014326.D
Acq On : 12 Dec 2019 15:12
Operator : SY/VA
Sample : K6236-14
Misc : 5.78G/10ML/MSVOA W/SOIL
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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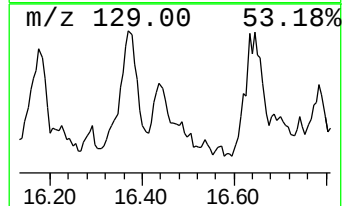
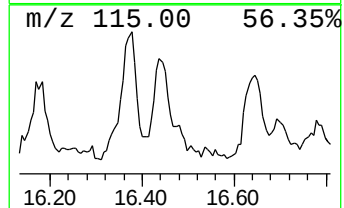
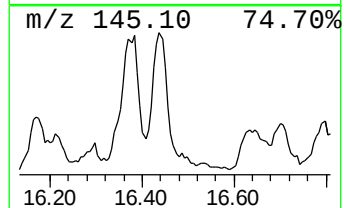
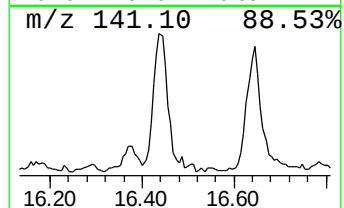
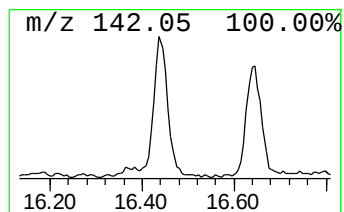
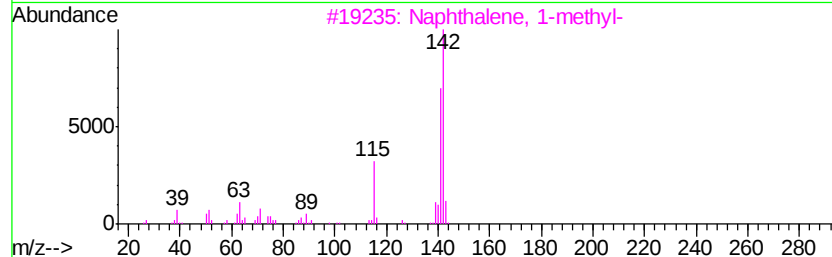
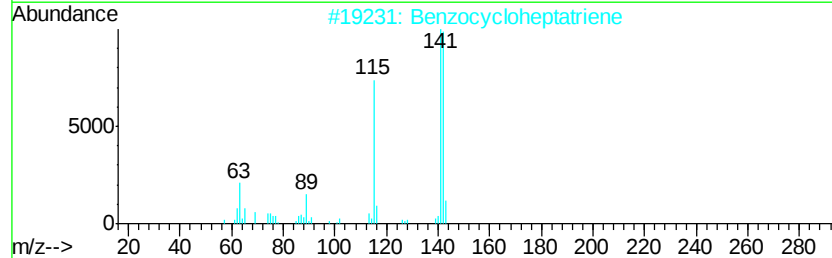
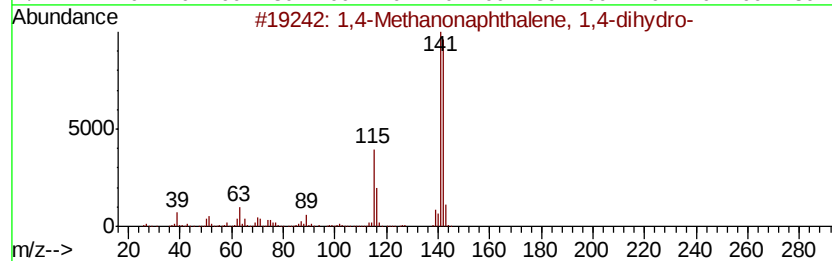
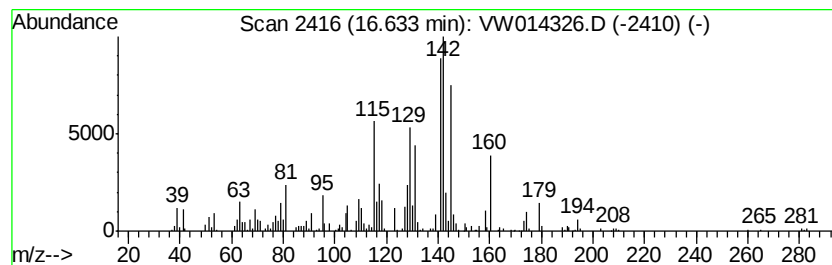
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Quant Title : VOC Analysis

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

Peak Number 23 unknown-03 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.63	2.14 ug/L	161406	1,4-Dichlorobenzene-d4	13.55

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	42
2			Benzocycloheptatriene	142	C11H10	000264-09-5	41
3			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	38
4			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	38
5			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	30



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW121219\
 Data File : VW014326.D
 Acq On : 12 Dec 2019 15:12
 Operator : SY/VA
 Sample : K6236-14
 Misc : 5.78G/10ML/MSVOA_W/SOIL
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 BFEW1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM120519S.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
Naphthalene, 1,6-...	12.87	4.5	ug/L	341996	3	13.55	1887810	25.0
Naphthalene, 2,3-...	13.30	12.4	ug/L	934035	3	13.55	1887810	25.0
Naphthalene, 2,6-...	13.46	10.9	ug/L	822604	3	13.55	1887810	25.0
unknown-01	14.07	2.9	ug/L	218703	3	13.55	1887810	25.0
(DEL) Alkane: Cyc...	14.38	1.6	ug/L	123532	3	13.55	1887810	25.0
unknown-02	14.55	1.3	ug/L	99165	3	13.55	1887810	25.0
Benzene, 1-ethyl-...	14.79	3.3	ug/L	246869	3	13.55	1887810	25.0
2-Propenal, 3-(2-...	14.85	2.4	ug/L	179482	3	13.55	1887810	25.0
Naphthalene, 1,2,...	14.96	1.7	ug/L	125064	3	13.55	1887810	25.0
1H-Indene, 2,3-di...	15.06	1.4	ug/L	103347	3	13.55	1887810	25.0
1H-Indene, 2,3-di...	15.18	1.9	ug/L	146587	3	13.55	1887810	25.0
(DEL) Alkane: Str...	15.33	5.5	ug/L	418923	3	13.55	1887810	25.0
1H-Indole-2,3-dio...	15.43	1.8	ug/L	136835	3	13.55	1887810	25.0
Benzene, 1-methyl...	15.66	2.3	ug/L	176838	3	13.55	1887810	25.0
(DEL) Alkane: Cyc...	15.73	2.5	ug/L	185005	3	13.55	1887810	25.0
Naphthalene, 1,2,...	15.86	3.3	ug/L	248111	3	13.55	1887810	25.0
(DEL) Alkane: Str...	16.29	3.7	ug/L	282536	3	13.55	1887810	25.0
Naphthalene, 1,2,...	16.37	6.3	ug/L	474123	3	13.55	1887810	25.0
Naphthalene, 1,2,...	16.44	4.8	ug/L	361152	3	13.55	1887810	25.0
unknown-03	16.63	2.1	ug/L	161406	3	13.55	1887810	25.0