

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW102818\
 Data File : VW006408.D
 Acq On : 26 Oct 2018 15:40
 Operator : SY/AP
 Sample : J5635-25
 Misc : 4.88G/10ML/MSVOA W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 A40G1

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\SOM2WLM100918S.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	1.952	5	8	26	rVB2	14637	32824	0.26%	0.092%
2	2.093	26	31	37	rVB3	1912	3339	0.03%	0.009%
3	2.361	65	75	86	rBV	54757	126960	1.02%	0.354%
4	2.458	88	91	94	rVV2	2303	3340	0.03%	0.009%
5	2.495	94	97	105	rVB3	2195	3919	0.03%	0.011%
6	2.574	105	110	119	rBV8	1794	4538	0.04%	0.013%
7	2.812	144	149	154	rBV	3535	6534	0.05%	0.018%
8	2.897	155	163	177	rVB	47337	131231	1.05%	0.366%
9	3.403	238	246	247	rBV5	2165	4180	0.03%	0.012%
10	3.726	293	299	308	rVB6	1861	5685	0.05%	0.016%
11	4.013	335	346	358	rBV	108848	319711	2.56%	0.892%
12	4.135	358	366	379	rVB2	20410	61084	0.49%	0.171%
13	4.373	398	405	417	rBV2	3441	10635	0.09%	0.030%
14	4.909	484	493	494	rBV2	4741	8970	0.07%	0.025%
15	5.787	627	637	644	rBV5	2540	7823	0.06%	0.022%
16	5.933	653	661	662	rBV3	1760	2637	0.02%	0.007%
17	6.177	697	701	704	rVV4	1509	3431	0.03%	0.010%
18	7.086	842	850	857	rVV	28345	82005	0.66%	0.229%
19	7.153	857	861	873	rVB2	11981	34795	0.28%	0.097%
20	7.366	883	896	897	rBV6	1801	5433	0.04%	0.015%
21	7.653	932	943	952	rBV	171303	412567	3.31%	1.152%
22	8.152	1014	1025	1026	rBV5	3369	7640	0.06%	0.021%
23	8.281	1026	1046	1066	rVV2	334461	1085662	8.70%	3.031%
24	8.433	1066	1071	1077	rVV5	2905	7701	0.06%	0.021%
25	8.634	1088	1104	1109	rBV3	11351	43067	0.35%	0.120%
26	8.695	1109	1114	1122	rVV5	10279	29703	0.24%	0.083%
27	8.756	1122	1124	1130	rVB4	2805	5195	0.04%	0.015%
28	8.848	1130	1139	1150	rBV	369608	749956	6.01%	2.093%
29	9.280	1199	1210	1228	rBV	223144	509192	4.08%	1.421%
30	9.494	1241	1245	1246	rBV3	2210	2332	0.02%	0.007%
31	9.549	1252	1254	1261	rVB3	3759	7272	0.06%	0.020%
32	9.792	1274	1294	1297	rBV3	11720	50997	0.41%	0.142%
33	10.036	1322	1334	1354	rVB	121921	271536	2.18%	0.758%
34	10.201	1354	1361	1362	rBV6	2067	3818	0.03%	0.011%

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Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM100918S.M
 Title : VOC Analysis

35	10.329	1374	1382	1388	rBV	460315	818889	6.56%	2.286%
36	10.402	1388	1394	1403	rVV3	43403	101612	0.81%	0.284%
37	10.512	1403	1412	1417	rVV3	29420	71286	0.57%	0.199%
38	10.579	1417	1423	1429	rVV	80534	159069	1.27%	0.444%
39	10.628	1429	1431	1437	rVB2	7734	10392	0.08%	0.029%
40	10.695	1437	1442	1448	rBV2	20321	34904	0.28%	0.097%
41	10.786	1451	1457	1465	rBV5	7153	17033	0.14%	0.048%
42	10.926	1474	1480	1498	rVB	120944	210901	1.69%	0.589%
43	11.182	1516	1522	1526	rBV7	2788	5240	0.04%	0.015%
44	11.457	1534	1567	1569	rBV3	108795	505400	4.05%	1.411%
45	11.634	1590	1596	1608	rVB	518256	986646	7.91%	2.754%
46	11.731	1608	1612	1619	rVB2	6903	14892	0.12%	0.042%
47	11.847	1622	1631	1641	rBV	30203	57678	0.46%	0.161%
48	12.170	1676	1684	1692	rVB	47748	82896	0.66%	0.231%
49	12.255	1693	1698	1704	rVB4	2011	3334	0.03%	0.009%
50	12.469	1723	1733	1739	rBV	248932	421065	3.37%	1.175%
51	12.694	1739	1770	1796	rVB3	288802	1623875	13.01%	4.533%
52	12.902	1800	1804	1806	rBV2	14288	17429	0.14%	0.049%
53	12.950	1806	1812	1821	rVB	115421	198807	1.59%	0.555%
54	13.042	1823	1827	1830	rBV6	3131	5504	0.04%	0.015%
55	13.127	1830	1841	1842	rBV3	28067	70837	0.57%	0.198%
56	13.200	1844	1853	1858	rVV2	66719	232368	1.86%	0.649%
57	13.255	1860	1862	1868	rVB	57437	75657	0.61%	0.211%
58	13.304	1868	1870	1875	rVB	16460	23374	0.19%	0.065%
59	13.414	1875	1888	1892	rBV2	62565	137307	1.10%	0.383%
60	13.450	1892	1894	1896	rVB	8787	6899	0.06%	0.019%
61	13.505	1896	1903	1908	rBV	1173152	1872813	15.01%	5.228%
62	13.584	1908	1916	1926	rVV	7661943	12477168	100.00%	34.829%
63	13.725	1934	1939	1941	rBV3	22931	33304	0.27%	0.093%
64	13.773	1941	1947	1950	rBV5	42194	81426	0.65%	0.227%
65	13.865	1955	1962	1972	rVV2	415804	949222	7.61%	2.650%
66	13.956	1972	1977	1982	rVB5	25082	46881	0.38%	0.131%
67	14.048	1989	1992	1994	rBV2	7256	8878	0.07%	0.025%
68	14.109	1994	2002	2013	rVB4	68531	218635	1.75%	0.610%
69	14.206	2013	2018	2023	rBV4	17875	37181	0.30%	0.104%
70	14.261	2023	2027	2029	rBV2	19338	29879	0.24%	0.083%
71	14.401	2036	2050	2058	rBV3	174541	723325	5.80%	2.019%

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 Title : VOC Analysis

72	14.468	2058	2061	2062	rVV	68865	88377	0.71%	0.247%
73	14.529	2062	2071	2087	rVB2	883740	3142596	25.19%	8.772%
74	14.645	2087	2090	2094	rBV3	13143	17927	0.14%	0.050%
75	14.694	2094	2098	2101	rVB5	10276	12693	0.10%	0.035%
76	14.731	2101	2104	2108	rBV	52199	70378	0.56%	0.196%
77	14.798	2108	2115	2125	rVV4	73790	237769	1.91%	0.664%
78	14.913	2125	2134	2144	rVB	168955	482364	3.87%	1.346%
79	15.072	2156	2160	2163	rBV5	8103	12232	0.10%	0.034%
80	15.139	2166	2171	2175	rVV2	21593	40803	0.33%	0.114%
81	15.182	2176	2178	2187	rVB7	10410	22593	0.18%	0.063%
82	15.267	2188	2192	2199	rVB9	6475	11870	0.10%	0.033%
83	15.346	2199	2205	2206	rBV2	13086	21208	0.17%	0.059%
84	15.377	2206	2210	2214	rVB	23129	39309	0.32%	0.110%
85	15.413	2214	2216	2222	rVB5	6332	9358	0.08%	0.026%
86	15.511	2226	2232	2236	rBV	7819	13272	0.11%	0.037%
87	15.578	2236	2243	2248	rBV3	37636	87747	0.70%	0.245%
88	15.651	2248	2255	2263	rVV3	19342	53623	0.43%	0.150%
89	15.724	2264	2267	2279	rVB6	13847	41790	0.33%	0.117%
90	15.822	2279	2283	2286	rBV2	16034	25458	0.20%	0.071%
91	15.865	2286	2290	2293	rBV2	13628	17867	0.14%	0.050%
92	16.011	2302	2314	2317	rBV5	49916	160670	1.29%	0.449%
93	16.182	2335	2342	2346	rBV	378288	684908	5.49%	1.912%
94	16.230	2346	2350	2356	rVV	609131	1234894	9.90%	3.447%
95	16.285	2356	2359	2368	rVB3	203190	371300	2.98%	1.036%
96	16.364	2368	2372	2375	rBV2	4910	5828	0.05%	0.016%
97	16.432	2375	2383	2389	rBV	480281	917411	7.35%	2.561%
98	16.511	2389	2396	2404	rVB	468935	867121	6.95%	2.421%
99	16.614	2405	2413	2423	rBV	204197	395416	3.17%	1.104%
100	16.761	2424	2437	2441	rBV3	79069	319334	2.56%	0.891%

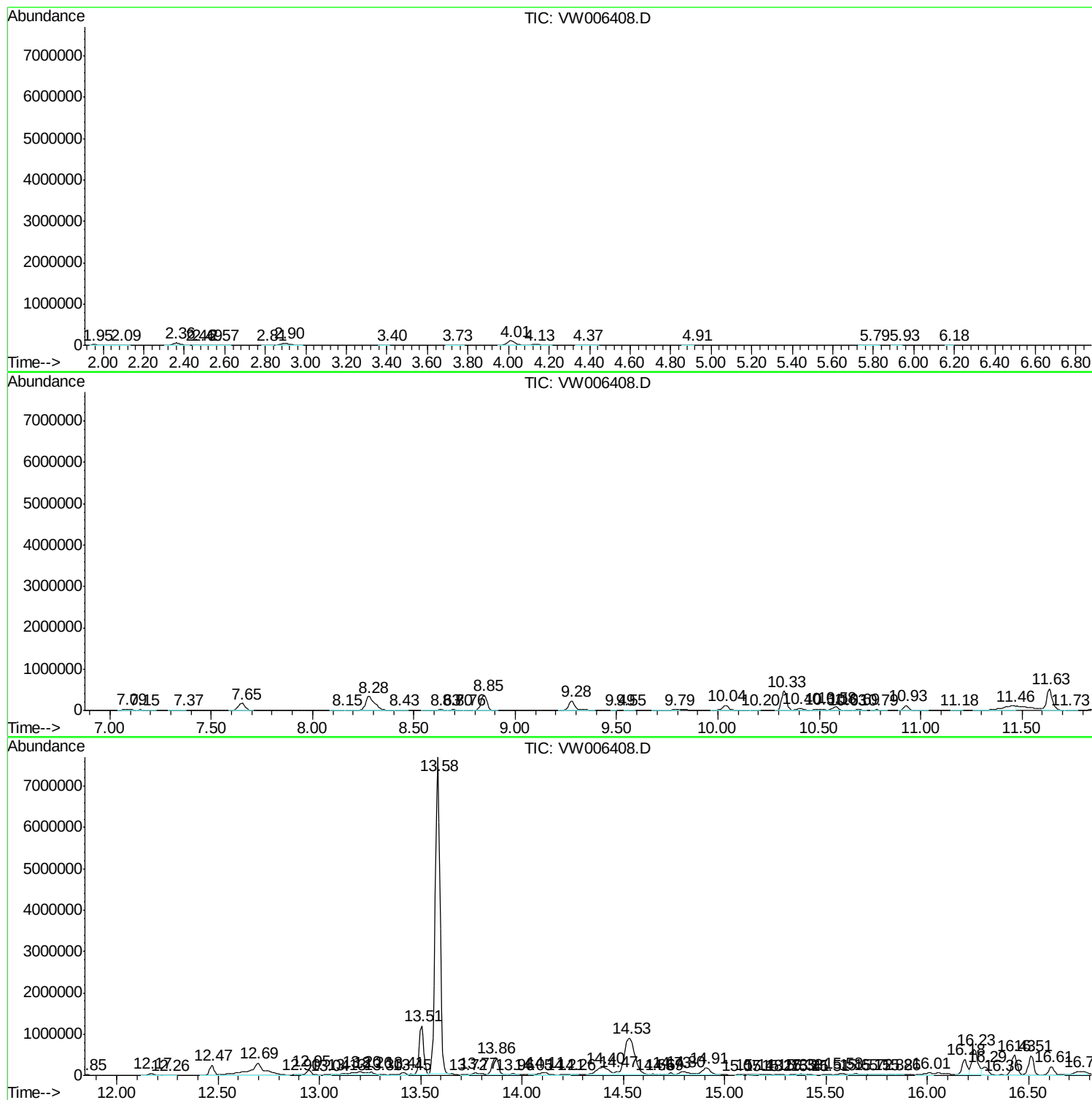
Sum of corrected areas: 35823834

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM100918S.M
Quant Title : VOC Analysis

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



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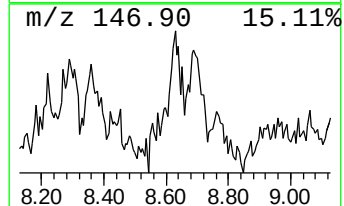
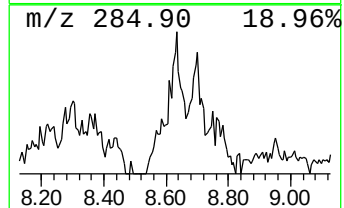
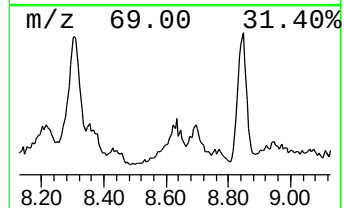
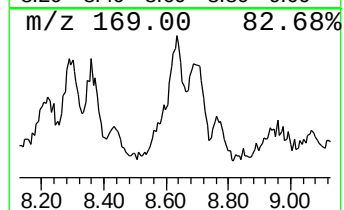
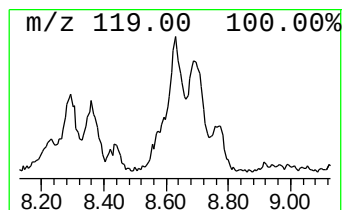
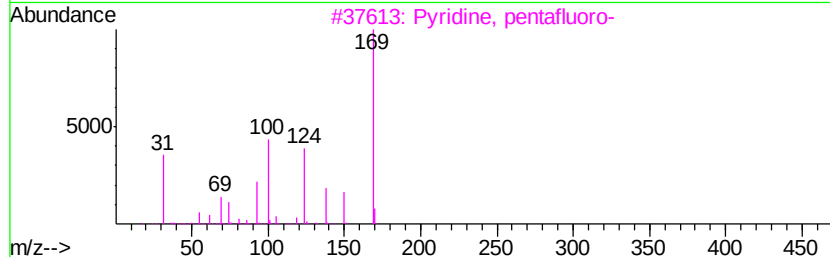
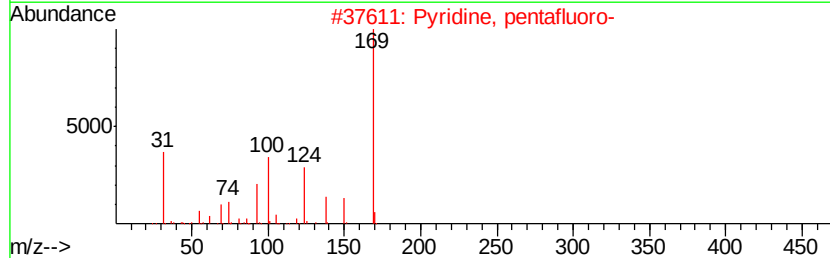
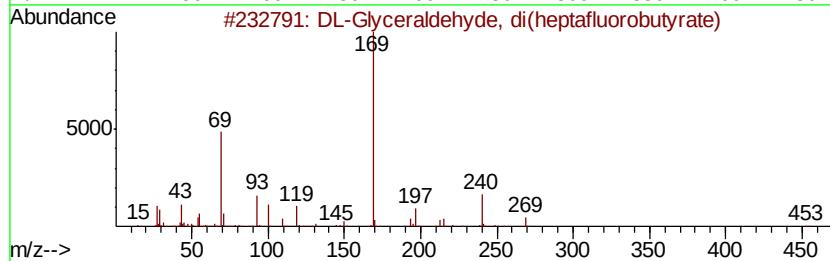
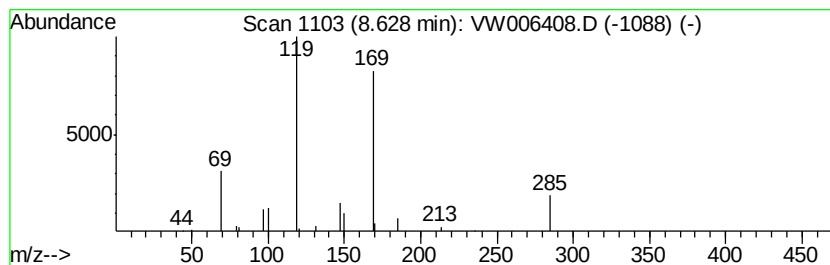
Instrument :
MSVOA_W
ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM100918S.M
Quant Title : VOC Analysis

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

Peak Number 1 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
8.63	1.44 ug/L	43067	1,4-Difluorobenzene	8.85		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		DL-Glyceraldehyde, di(heptafluor...	482	C11H4F14O5	1000365-33-0	17
2		Pyridine, pentafluoro-	169	C5F5N	000700-16-3	9
3		Pyridine, pentafluoro-	169	C5F5N	000700-16-3	9
4		Bicyclo[2.2.1]heptan-2-ol, 3-ami...	169	C10H19NO	032344-86-8	9
5		2-(2,2,3,3,4,4,4-Heptafluorobuta...	454	C10H4F14O4	1000352-03-2	9



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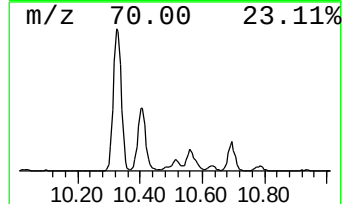
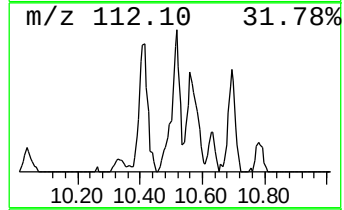
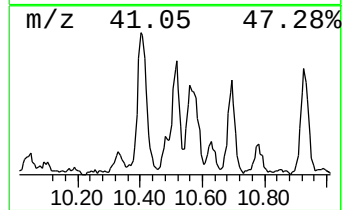
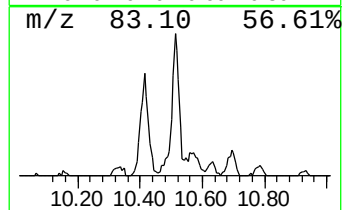
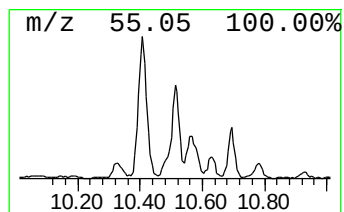
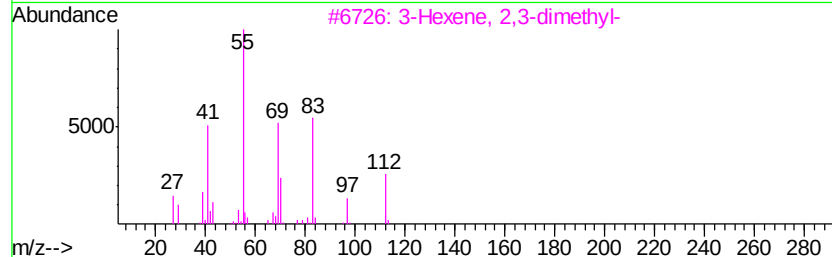
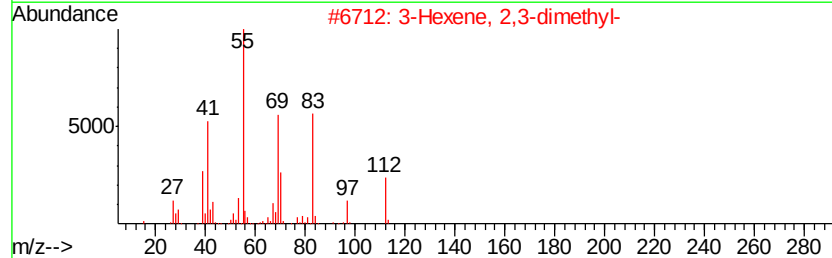
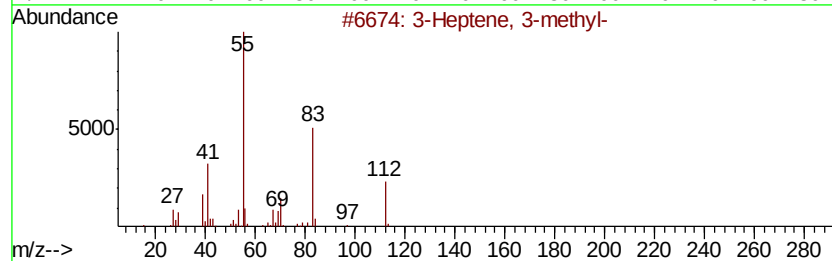
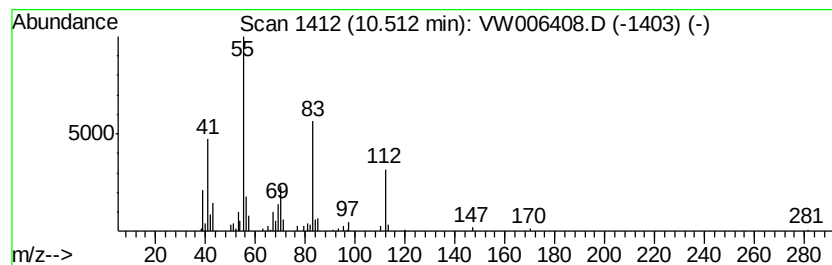
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM100918S.M
Quant Title : VOC Analysis

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

Peak Number 4 3-Heptene, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.51	1.81 ug/L	71286	Chlorobenzene-d5	11.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Heptene, 3-methyl-	112	C8H16	007300-03-0	83
2		3-Hexene, 2,3-dimethyl-	112	C8H16	007145-23-5	81
3		3-Hexene, 2,3-dimethyl-	112	C8H16	007145-23-5	74
4		3-Heptene, 5-methyl-	112	C8H16	013172-91-3	68
5		2-Pentene, 3-ethyl-2-methyl-	112	C8H16	019780-67-7	58



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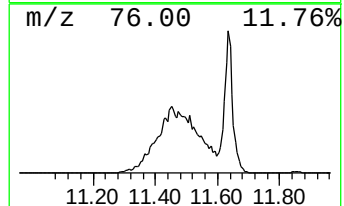
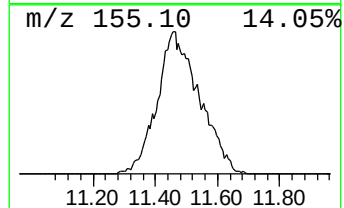
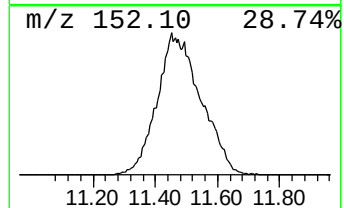
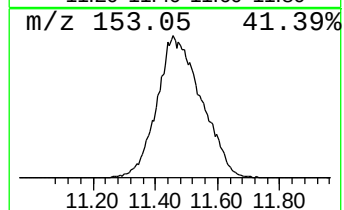
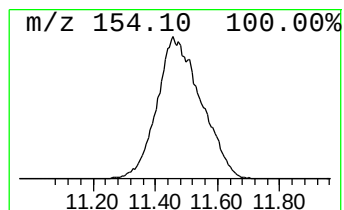
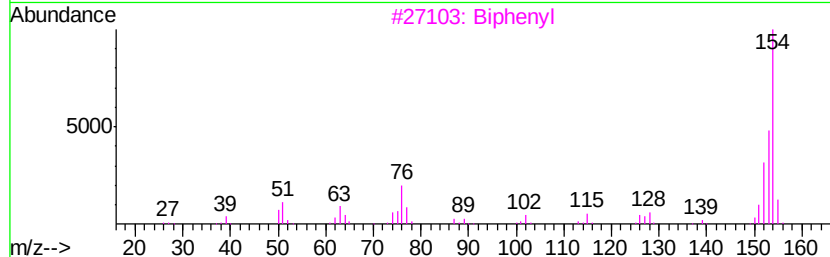
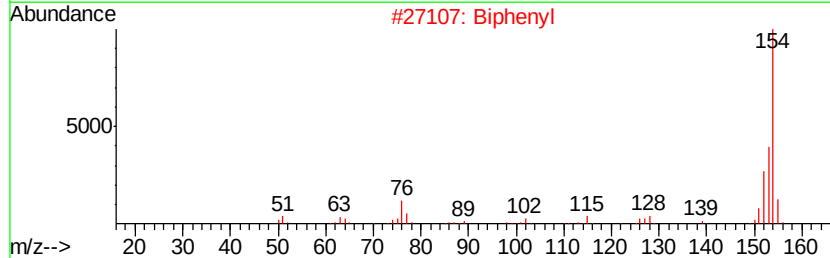
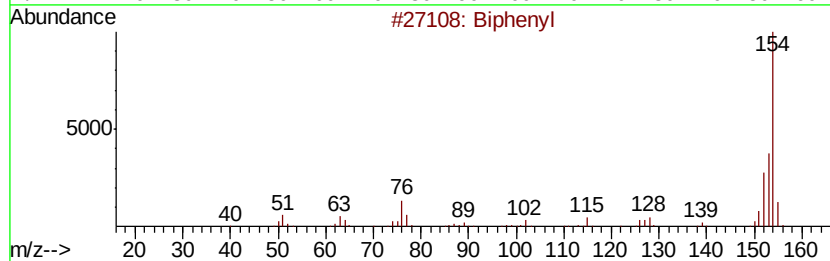
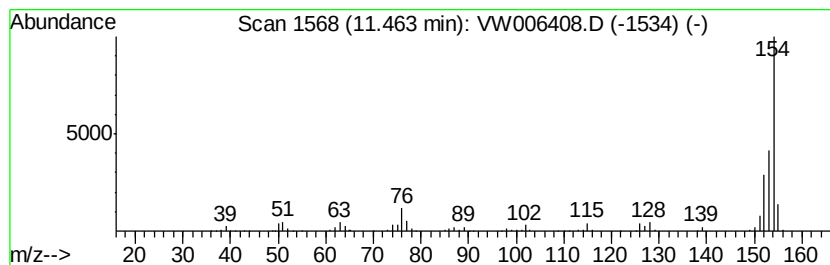
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Quant Title : VOC Analysis

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

Peak Number 5 Biphenyl Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.46	12.81 ug/L	505400	Chlorobenzene-d5	11.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Biphenyl	154	C12H10	000092-52-4	95
2		Biphenyl	154	C12H10	000092-52-4	95
3		Biphenyl	154	C12H10	000092-52-4	93
4		Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	87
5		Biphenyl	154	C12H10	000092-52-4	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW102818\
Data File : VW006408.D
Acq On : 26 Oct 2018 15:40
Operator : SY/AP
Sample : J5635-25
Misc : 4.88G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

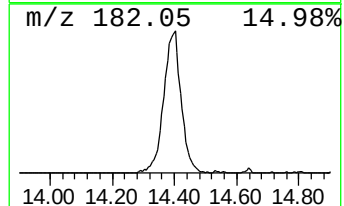
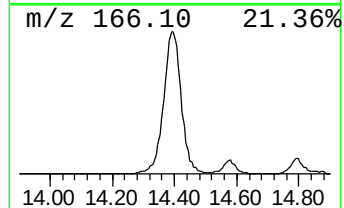
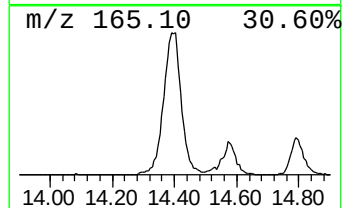
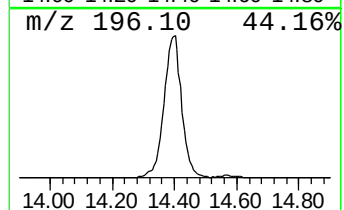
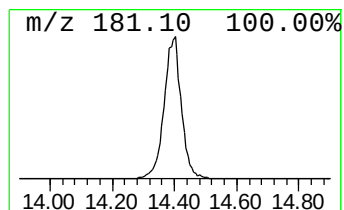
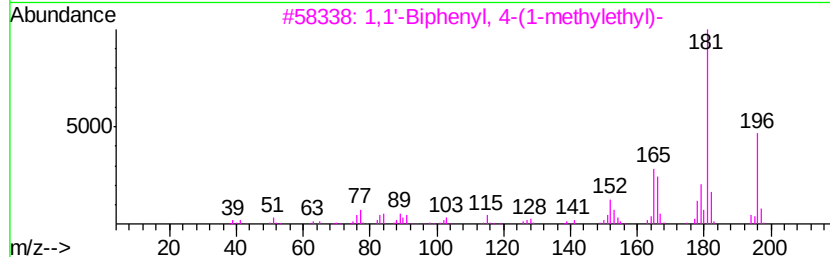
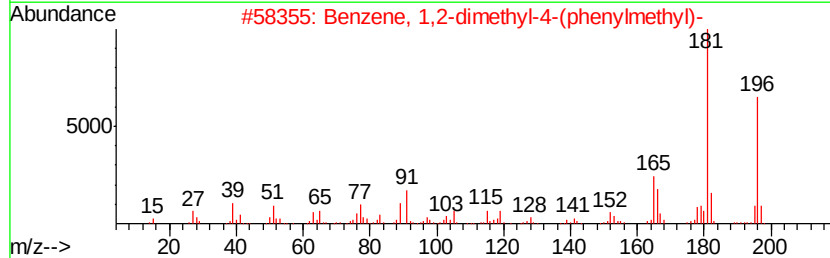
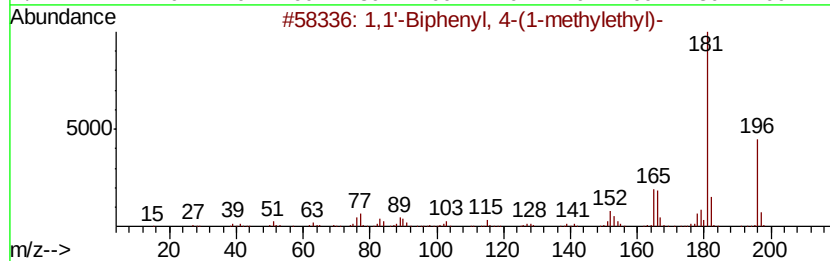
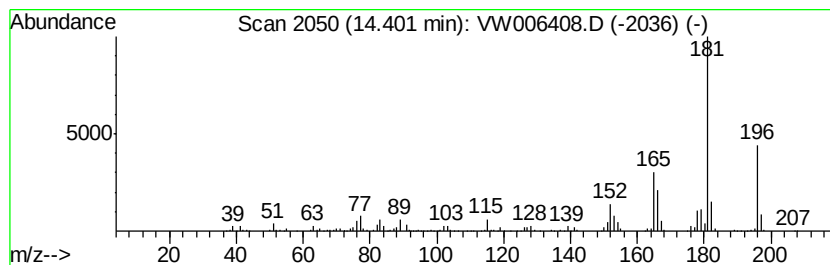
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM100918S.M
Quant Title : VOC Analysis

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

Peak Number 6 1,1'-Biphenyl, 4-(1-methyle... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.40	1.45 ug/L	723325	1,4-Dichlorobenzene-d4	13.57	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4-(1-methylethyl)-	196	C15H16	007116-95-2	97
2	Benzene, 1,2-dimethyl-4-(phenylm...	196	C15H16	013540-56-2	94
3	1,1'-Biphenyl, 4-(1-methylethyl)-	196	C15H16	007116-95-2	93
4	Methane, di-p-tolyl-	196	C15H16	004957-14-6	91
5	Benzene, 1-methyl-3-[(4-methylph...	196	C15H16	021895-16-9	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW102818\
Data File : VW006408.D
Acq On : 26 Oct 2018 15:40
Operator : SY/AP
Sample : J5635-25
Misc : 4.88G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

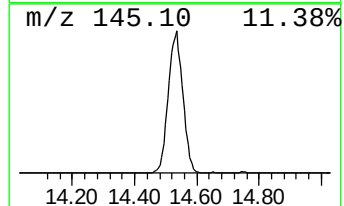
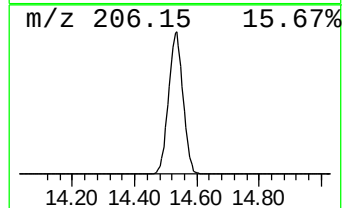
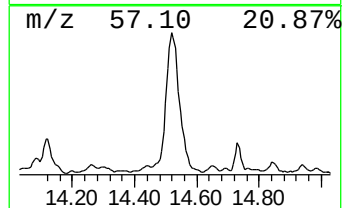
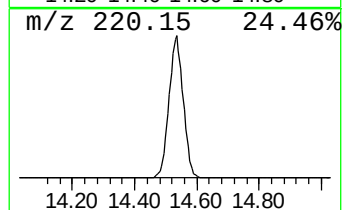
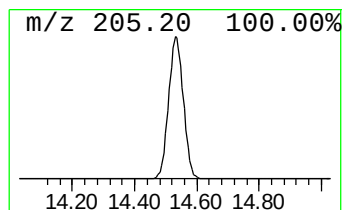
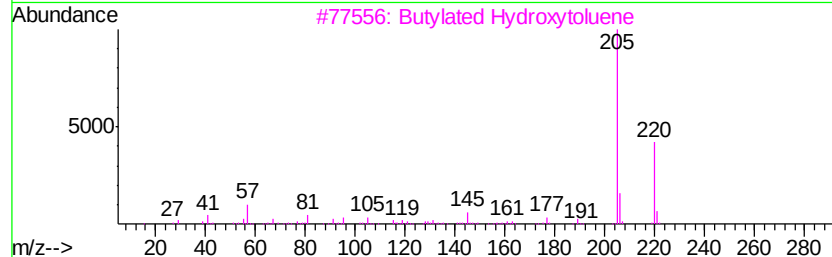
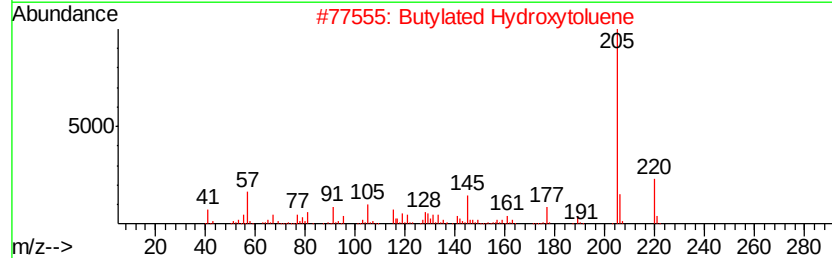
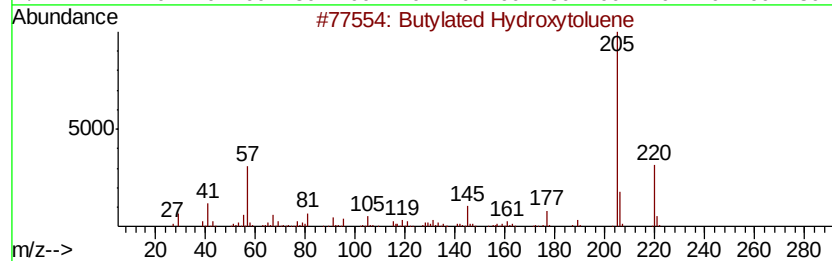
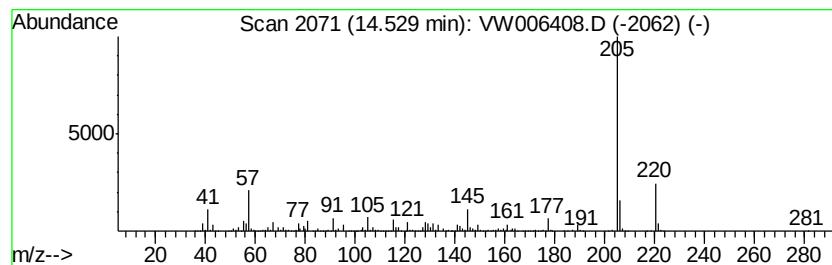
Instrument :
MSVOA_W
ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM100918S.M
Quant Title : VOC Analysis

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

Peak Number 7 Butylated Hydroxytoluene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.53	6.30 ug/L	3142600	1,4-Dichlorobenzene-d4	13.57		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	97	
2	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	96	
3	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	94	
4	Phenol, 4,6-di(1,1-dimethylethyl...	220	C15H24O	000616-55-7	81	
5	Propanoic acid, 2-methyl-3-[4-t-...	220	C14H20O2	1000131-87-0	72	



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW102818\
Data File : VW006408.D
Acq On : 26 Oct 2018 15:40
Operator : SY/AP
Sample : J5635-25
Misc : 4.88G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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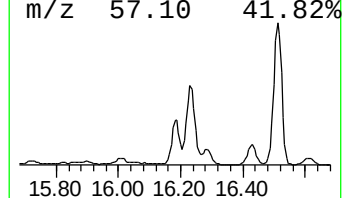
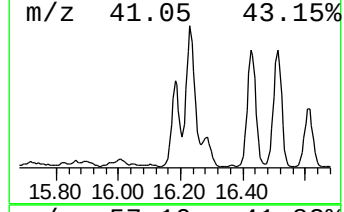
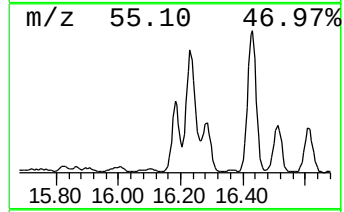
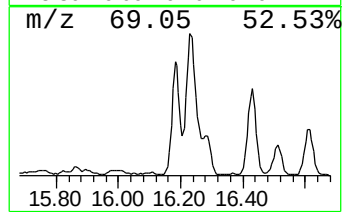
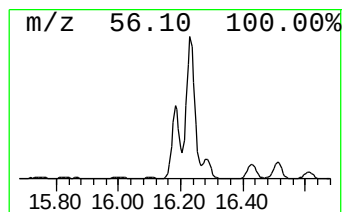
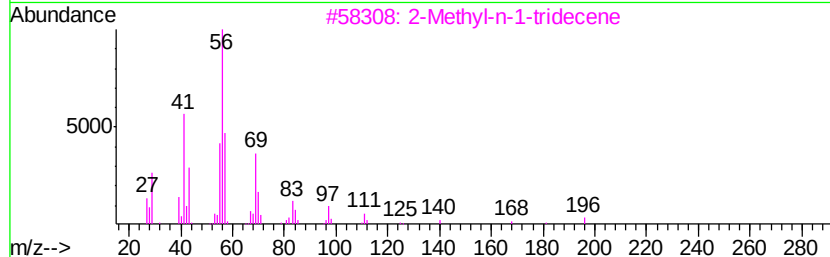
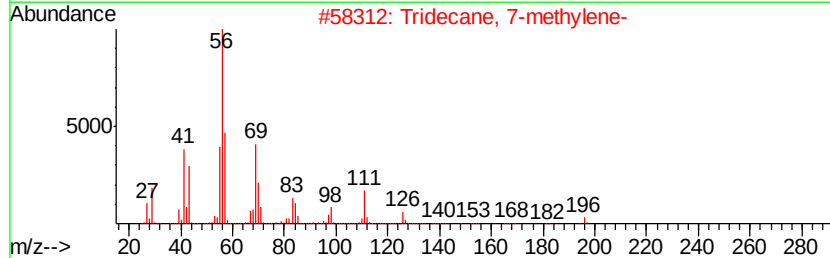
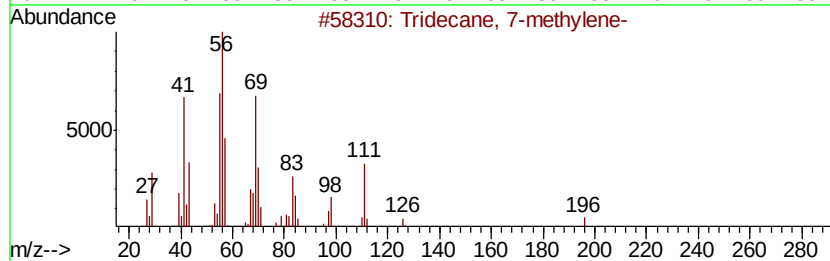
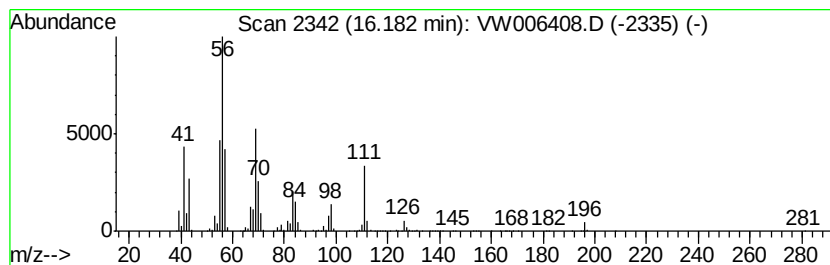
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Quant Title : VOC Analysis

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.18	1.37 ug/L	684908	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 7-methylene-	196	C14H28	019780-80-4	96
2		Tridecane, 7-methylene-	196	C14H28	019780-80-4	95
3		2-Methyl-n-1-tridecene	196	C14H28	018094-01-4	83
4		5-Undecene, 7-methyl-, (Z)-	168	C12H24	074630-62-9	72
5		5-Undecene, 7-methyl-, (E)-	168	C12H24	074630-66-3	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW102818\
Data File : VW006408.D
Acq On : 26 Oct 2018 15:40
Operator : SY/AP
Sample : J5635-25
Misc : 4.88G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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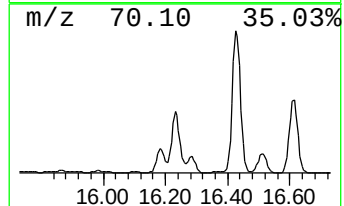
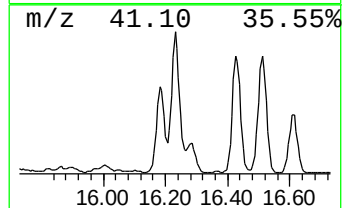
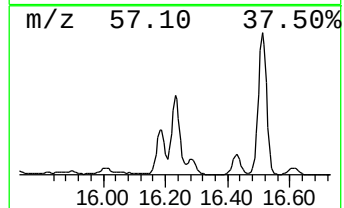
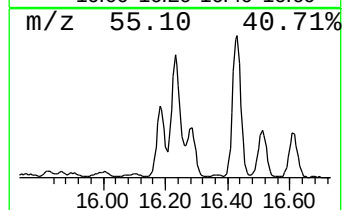
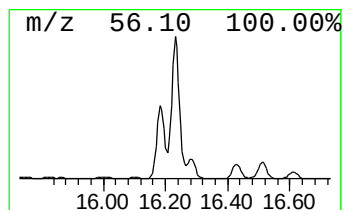
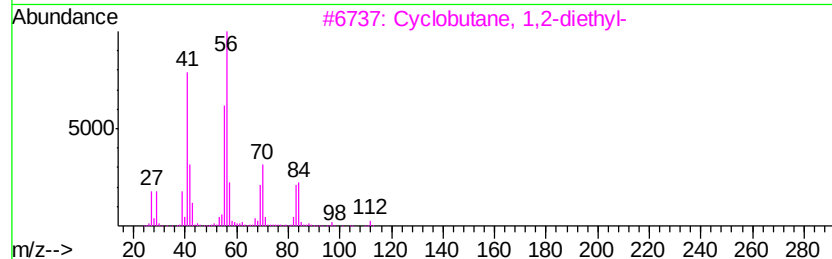
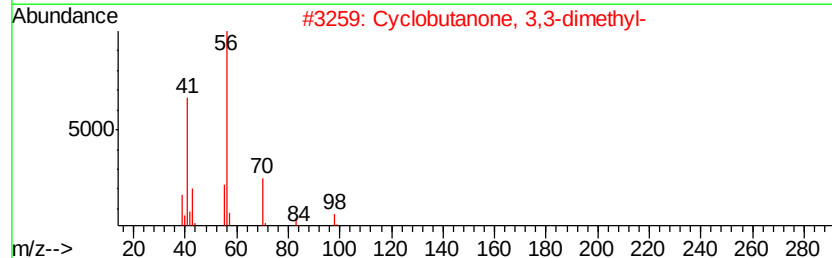
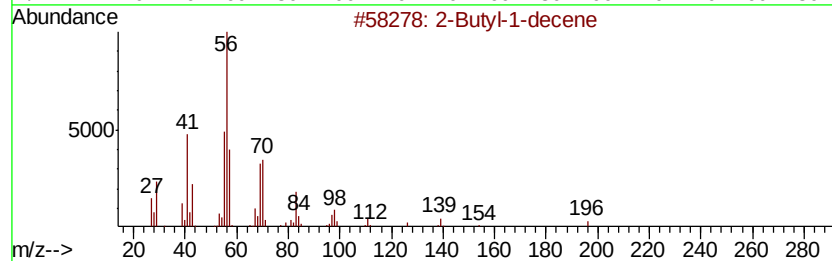
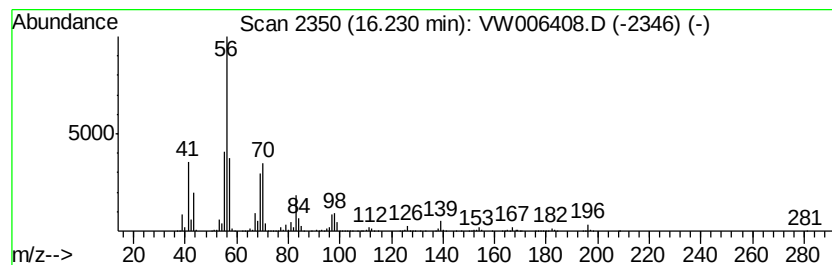
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Quant Title : VOC Analysis

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

Peak Number 9 2-Butyl-1-decene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.23	2.47 ug/L	1234890	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Butyl-1-decene	196	C14H28	051655-65-3	95
2		Cyclobutanone, 3,3-dimethyl-	98	C6H10O	001192-33-2	64
3		Cyclobutane, 1,2-diethyl-	112	C8H16	061141-83-1	59
4		5-Methyl-2-hexanol, trifluoroace...	212	C9H15F3O2	1000364-20-2	53
5		Undecane, 5-methylene-	168	C12H24	005698-48-6	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW102818\
Data File : VW006408.D
Acq On : 26 Oct 2018 15:40
Operator : SY/AP
Sample : J5635-25
Misc : 4.88G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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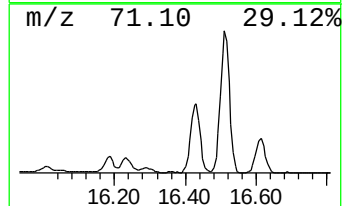
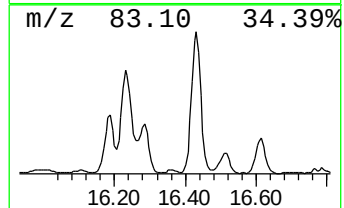
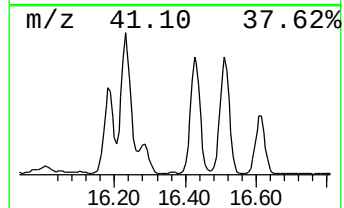
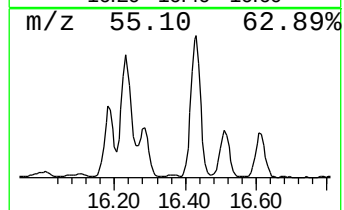
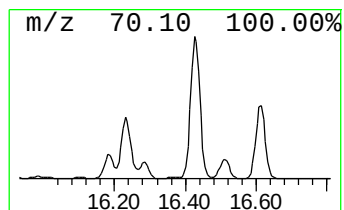
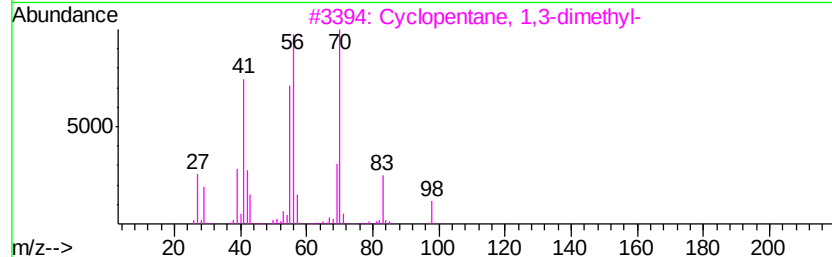
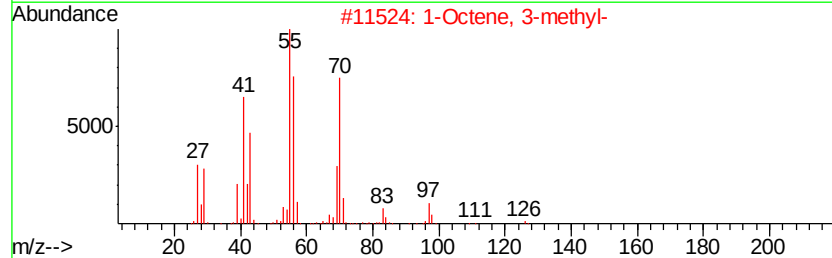
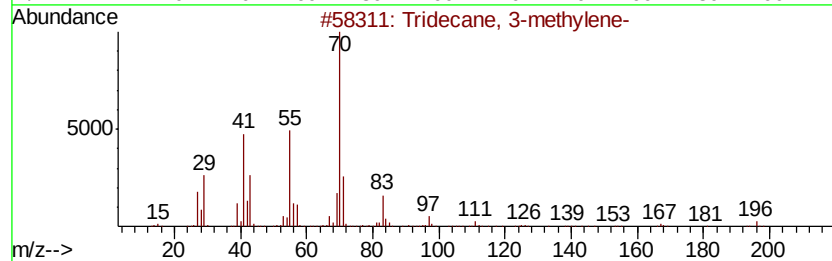
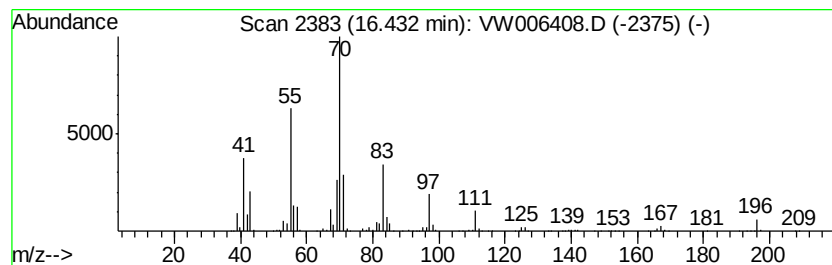
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Quant Title : VOC Analysis

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.43	1.84 ug/L	917411	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 3-methylene-	196	C14H28	019780-34-8	74
2		1-Octene, 3-methyl-	126	C9H18	013151-08-1	59
3		Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	50
4		2,3-Dimethyl-1-hexene	112	C8H16	016746-86-4	47
5		Hexanal, 3-methyl-	114	C7H14O	019269-28-4	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW102818\
Data File : VW006408.D
Acq On : 26 Oct 2018 15:40
Operator : SY/AP
Sample : J5635-25
Misc : 4.88G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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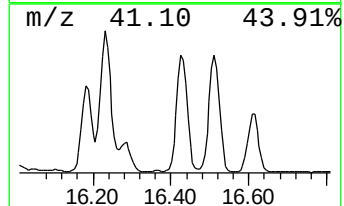
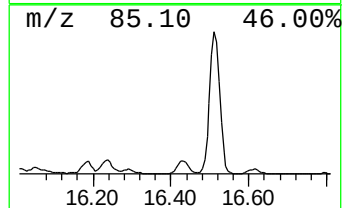
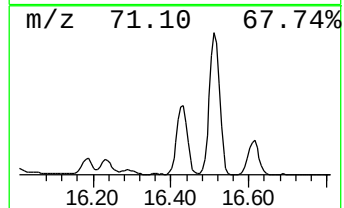
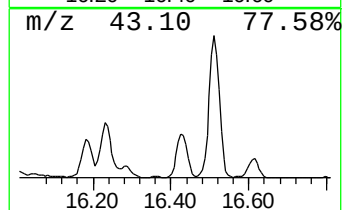
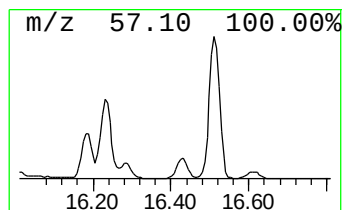
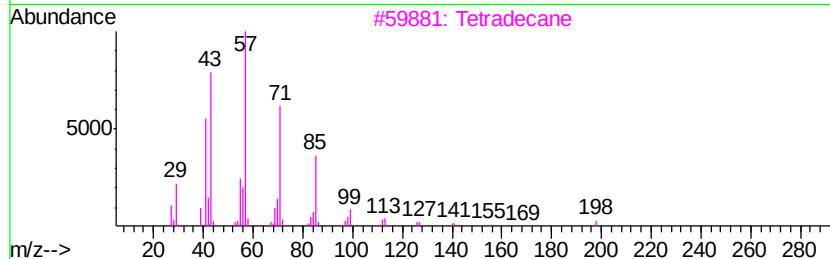
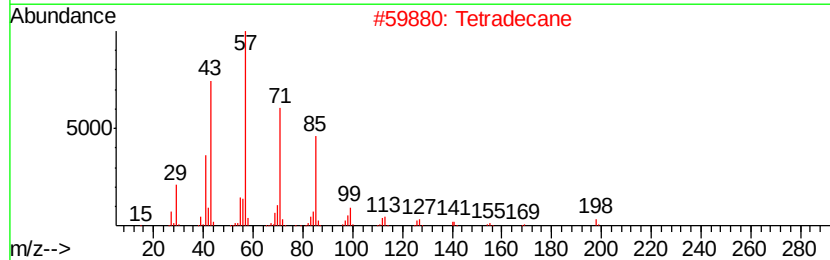
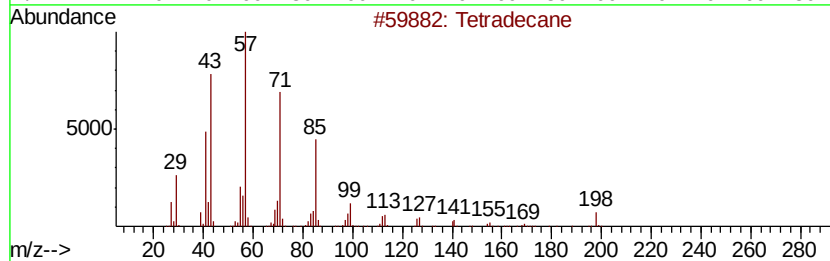
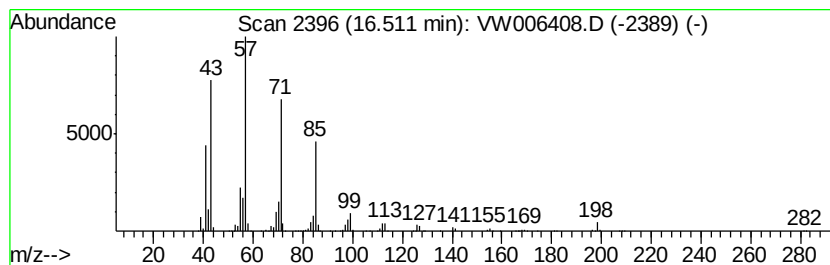
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Quant Title : VOC Analysis

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.51	1.74 ug/L	867121	1,4-Dichlorobenzene-d4	13.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	97
2			Tetradecane	198	C14H30	000629-59-4	96
3			Tetradecane	198	C14H30	000629-59-4	96
4			Hexadecane	226	C16H34	000544-76-3	90
5			Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	86



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW102818\
Data File : VW006408.D
Acq On : 26 Oct 2018 15:40
Operator : SY/AP
Sample : J5635-25
Misc : 4.88G/10ML/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
A40G1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM100918S.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
unknown-01	8.63	1.4	ug/L	43067	1	8.85	749956	25.0
3-Heptene, 3-methyl-	10.51	1.8	ug/L	71286	2	11.63	986646	25.0
Biphenyl	11.46	12.8	ug/L	505400	2	11.63	986646	25.0
1,1'-Biphenyl, 4-...	14.40	1.5	ug/L	723325	3	13.57	12477200	25.0
Butylated Hydroxy...	14.53	6.3	ug/L	3142600	3	13.57	12477200	25.0
(DEL) Alkane: Str...	16.18	1.4	ug/L	684908	3	13.57	12477200	25.0
2-Butyl-1-decene	16.23	2.5	ug/L	1234890	3	13.57	12477200	25.0
(DEL) Alkane: Str...	16.43	1.8	ug/L	917411	3	13.57	12477200	25.0
(DEL) Alkane: Str...	16.51	1.7	ug/L	867121	3	13.57	12477200	25.0