

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW092518\
 Data File : VW005619.D
 Acq On : 25 Sep 2018 18:44
 Operator : VA/AP
 Sample : J5066-09
 Misc : 4.83G/5ML/MSVOA W/SOIL
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EP1-Q1-H4

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % 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\82W092118S.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.891	970	982	986	rBV	221176	506778	3.43%	0.278%
2	7.952	986	992	1003	rVB	407009	889907	6.03%	0.489%
3	8.311	1039	1051	1060	rBV	227846	496459	3.36%	0.273%
4	8.848	1130	1139	1148	rBV	621608	1207702	8.18%	0.664%
5	10.329	1374	1382	1390	rBV	844457	1467372	9.94%	0.806%
6	11.238	1524	1531	1537	rBV	262202	504779	3.42%	0.277%
7	11.439	1559	1564	1573	rVB	165597	292160	1.98%	0.161%
8	11.634	1588	1596	1606	rBV	804305	1494390	10.12%	0.821%
9	11.768	1613	1618	1622	rBV	195517	319556	2.16%	0.176%
10	11.823	1622	1627	1632	rBV	207583	375802	2.54%	0.206%
11	11.878	1632	1636	1648	rVB3	303630	956174	6.47%	0.525%
12	11.981	1648	1653	1660	rBV2	195034	546041	3.70%	0.300%
13	12.067	1660	1667	1671	rVB4	133455	351128	2.38%	0.193%
14	12.146	1671	1680	1686	rBV3	1160631	2696206	18.25%	1.481%
15	12.256	1690	1698	1703	rVB2	985132	2179442	14.76%	1.197%
16	12.371	1703	1717	1722	rBV2	2315610	6430023	43.54%	3.533%
17	12.457	1722	1731	1739	rBV7	799921	2815028	19.06%	1.547%
18	12.567	1747	1749	1753	rVB	256752	303610	2.06%	0.167%
19	12.621	1753	1758	1763	rVB3	1148926	1970115	13.34%	1.082%
20	12.707	1763	1772	1776	rBV6	1316419	3422827	23.17%	1.881%
21	12.786	1780	1785	1791	rVB2	4853784	8712682	58.99%	4.787%
22	12.871	1791	1799	1804	rBV4	1804666	4669030	31.61%	2.565%
23	12.951	1805	1812	1825	rVB9	2187577	8231463	55.73%	4.523%
24	13.048	1825	1828	1830	rBV3	410886	529864	3.59%	0.291%
25	13.085	1830	1834	1838	rVV2	1808070	2767579	18.74%	1.521%
26	13.134	1838	1842	1845	rVV4	1583317	2770098	18.76%	1.522%
27	13.176	1845	1849	1862	rVV2	6309498	13283261	89.94%	7.298%
28	13.274	1862	1865	1866	rVV2	828040	1029694	6.97%	0.566%
29	13.304	1866	1870	1874	rVV	2192552	4276390	28.95%	2.350%
30	13.353	1874	1878	1881	rVV3	2163089	3802134	25.74%	2.089%
31	13.384	1881	1883	1887	rVV2	1179055	1922244	13.01%	1.056%
32	13.444	1887	1893	1897	rVV4	1919056	4938848	33.44%	2.714%
33	13.493	1897	1901	1905	rVB3	1528702	2260830	15.31%	1.242%
34	13.633	1921	1924	1925	rBV	510677	568053	3.85%	0.312%

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35	13.707	1931	1936	1940	rVB3	2102367	3620755	24.51%	1.989%
36	13.798	1946	1951	1954	rBV	1075122	1920446	13.00%	1.055%
37	13.890	1960	1966	1971	rBV2	8665260	14769712	100.00%	8.115%
38	13.993	1979	1983	1985	rBV3	1443928	2297360	15.55%	1.262%
39	14.103	1998	2001	2005	rVB	2207583	2825019	19.13%	1.552%
40	14.139	2005	2007	2012	rVB2	2048764	2897942	19.62%	1.592%
41	14.207	2013	2018	2023	rBV4	2757054	5291424	35.83%	2.907%
42	14.274	2024	2029	2035	rVB6	2005685	4303726	29.14%	2.365%
43	14.353	2035	2042	2045	rBV3	1896733	3764123	25.49%	2.068%
44	14.396	2045	2049	2052	rBV2	3845501	5607892	37.97%	3.081%
45	14.426	2052	2054	2058	rVV2	2028521	3196373	21.64%	1.756%
46	14.469	2058	2061	2065	rVB	3051755	4092949	27.71%	2.249%
47	14.511	2065	2068	2071	rBV	1274959	1670323	11.31%	0.918%
48	14.560	2071	2076	2083	rVB3	3337245	8072254	54.65%	4.435%
49	14.633	2084	2088	2096	rVB2	2015641	4782519	32.38%	2.628%
50	14.713	2096	2101	2103	rBV2	596163	1015562	6.88%	0.558%
51	14.743	2103	2106	2111	rVB3	1493049	2241366	15.18%	1.231%
52	14.804	2111	2116	2123	rBV2	2075636	4445496	30.10%	2.442%
53	14.926	2133	2136	2140	rVB	514037	720237	4.88%	0.396%
54	14.969	2140	2143	2145	rBV	443205	583238	3.95%	0.320%
55	15.078	2154	2161	2164	rBV3	2380940	4836218	32.74%	2.657%
56	15.182	2174	2178	2185	rVB3	2382494	5083406	34.42%	2.793%
57	15.328	2198	2202	2212	rVB3	717054	1629974	11.04%	0.896%
58	15.432	2213	2219	2221	rBV5	218919	452078	3.06%	0.248%
59	15.511	2229	2232	2239	rVB5	319181	712633	4.82%	0.392%
60	15.645	2250	2254	2264	rBV8	100191	311227	2.11%	0.171%
61	15.749	2265	2271	2275	rBV3	152413	337477	2.28%	0.185%
62	15.962	2301	2306	2311	rBV3	114370	213800	1.45%	0.117%
63	16.035	2314	2318	2323	rVV5	76817	158954	1.08%	0.087%
64	16.188	2334	2343	2355	rVB5	171319	492794	3.34%	0.271%
65	16.505	2383	2395	2409	rVB10	83309	463724	3.14%	0.255%
66	16.645	2411	2418	2427	rVV9	75497	210463	1.42%	0.116%

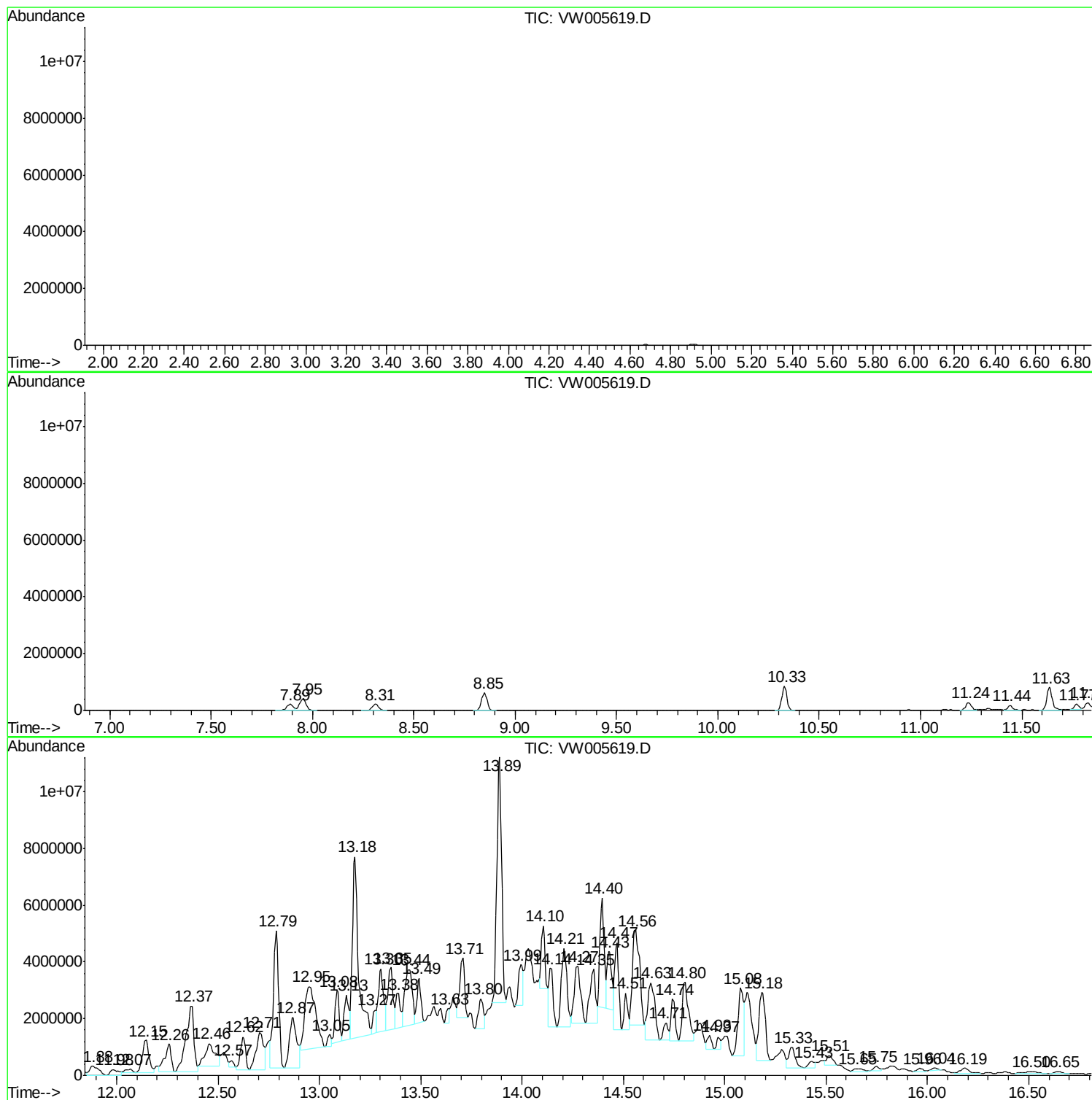
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W092118S.M
Quant Title : SW846 8260

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



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EP1-Q1-H4

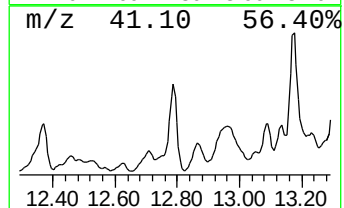
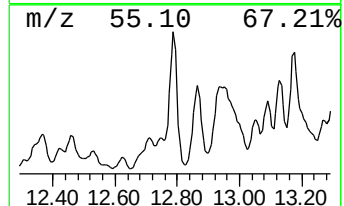
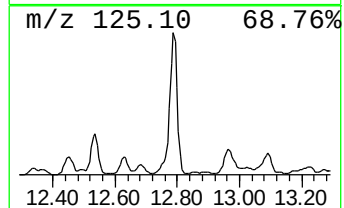
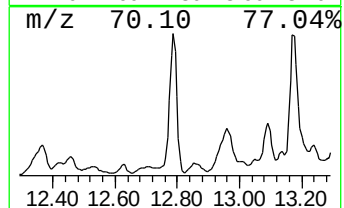
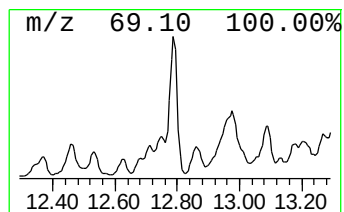
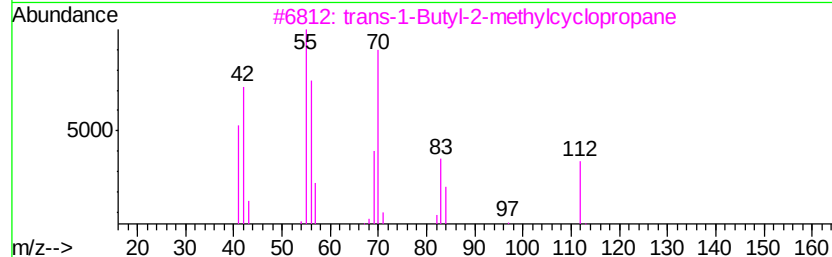
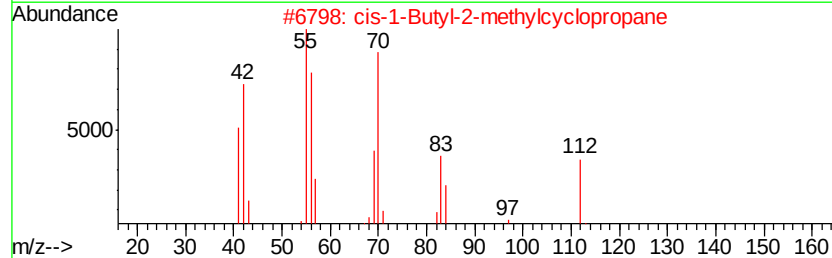
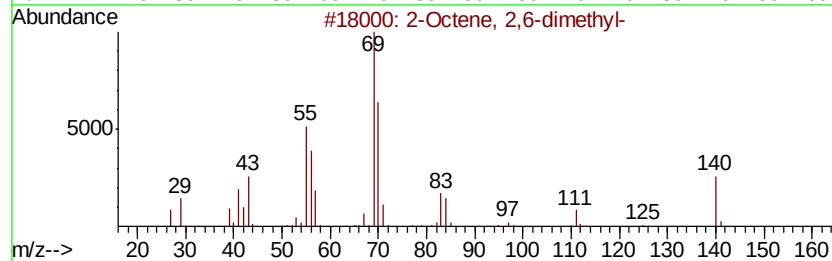
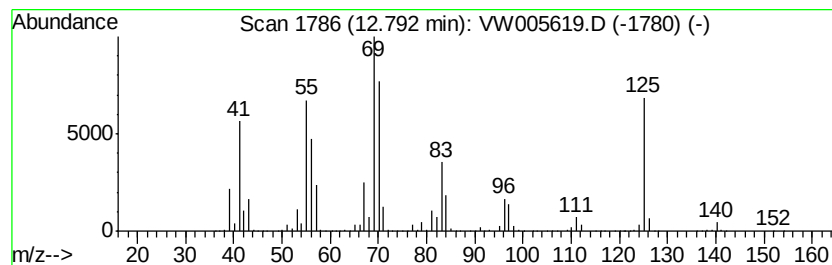
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W092118S.M
Quant Title : SW846 8260

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

Peak Number 1 2-Octene, 2,6-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.79	766.89 ug/l	8712680	1,4-Dichlorobenzene-d4	13.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	70
2			cis-1-Butyl-2-methylcyclopropane	112	C8H16	038851-69-3	50
3			trans-1-Butyl-2-methylcyclopropane	112	C8H16	038851-70-6	45
4			Cyclohexane, 1,2,3-trimethyl-, (...)	126	C9H18	001678-81-5	43
5			2-Methyl-2-octene	126	C9H18	016993-86-5	38



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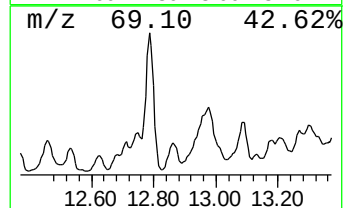
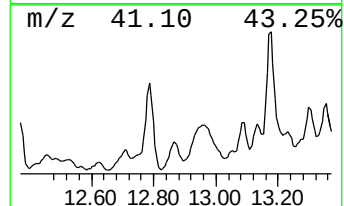
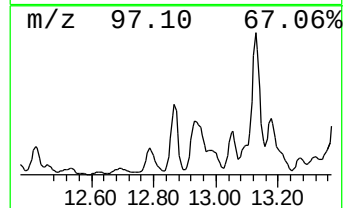
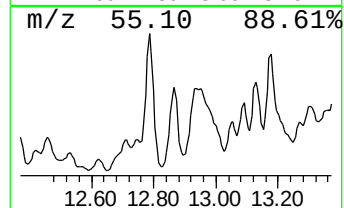
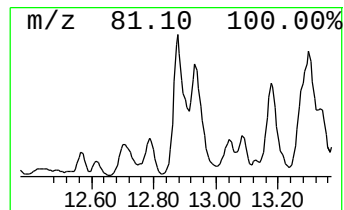
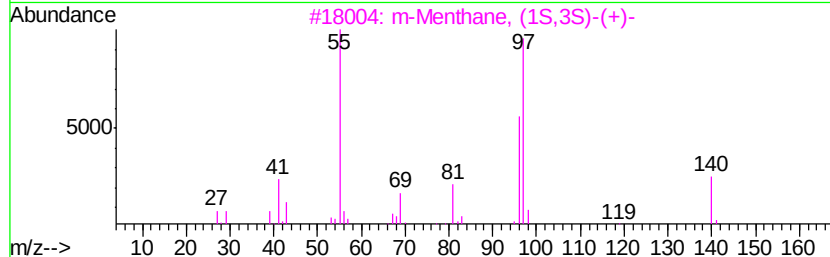
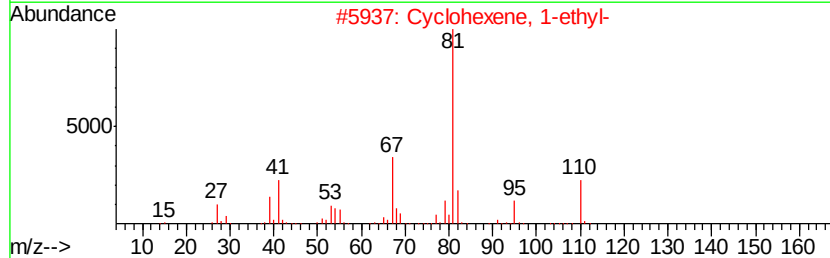
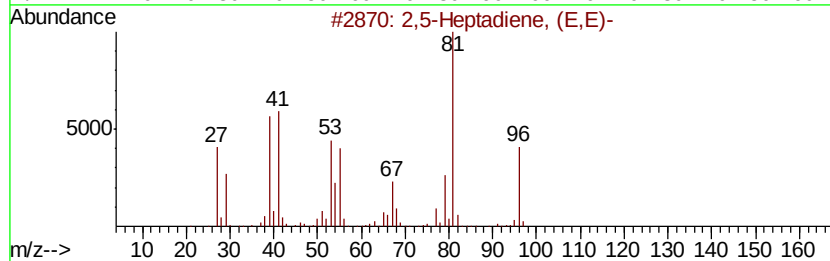
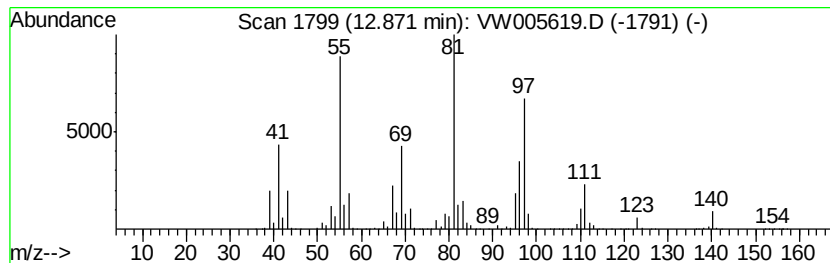
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W092118S.M
Quant Title : SW846 8260

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

Peak Number 2 unknown12.87 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.87	410.97 ug/l	4669030	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,5-Heptadiene, (E,E)-	96	C7H12	039619-60-8	46
2		Cyclohexene, 1-ethyl-	110	C8H14	001453-24-3	43
3		m-Menthane, (1S,3S)-(+)-	140	C10H20	013837-67-7	38
4		2,4-Heptadienal, (E,E)-	110	C7H10O	004313-03-5	35
5		Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	001678-82-6	30



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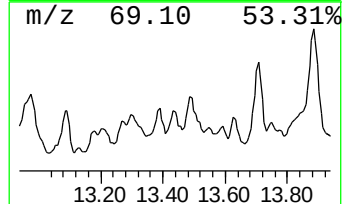
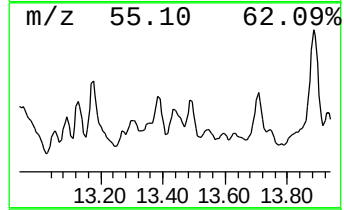
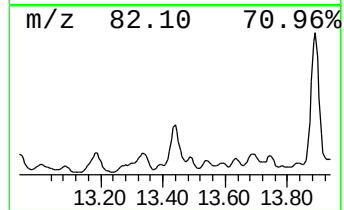
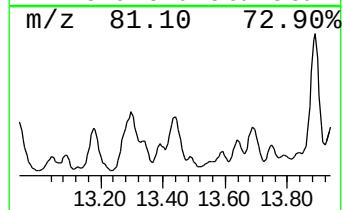
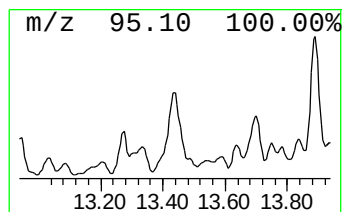
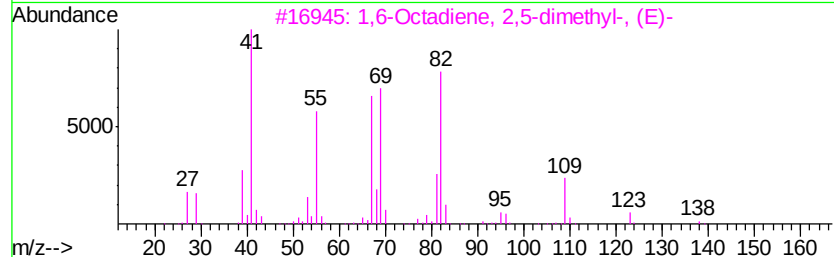
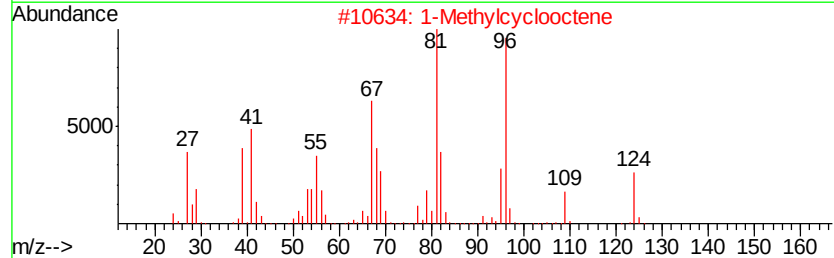
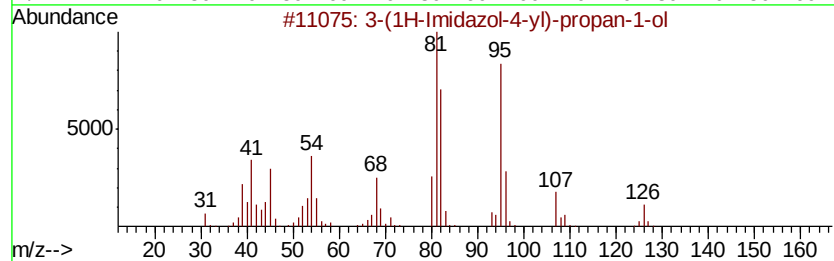
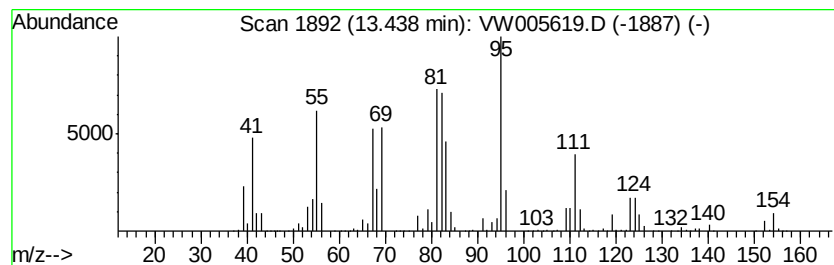
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Peak Number 3 3-(1H-Imidazol-4-yl)-propan... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.44	434.72 ug/l	4938850	1,4-Dichlorobenzene-d4	13.57		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-(1H-Imidazol-4-yl)-propan-1-ol	126	C6H10N2O	049549-75-9	58
2		1-Methylcyclooctene	124	C9H16	000933-11-9	38
3		1,6-Octadiene, 2,5-dimethyl-, (E)-	138	C10H18	068702-25-0	30
4		1H-Indene, octahydro-, cis-	124	C9H16	004551-51-3	30
5		Bicyclo[3.1.1]heptane, 2,6,6-tri...	138	C10H18	006876-13-7	30



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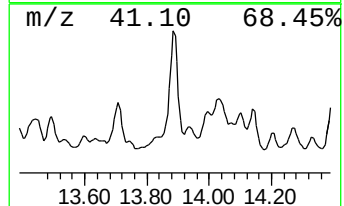
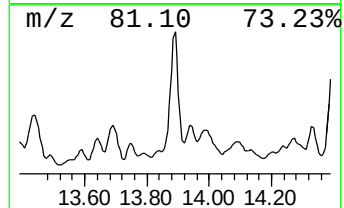
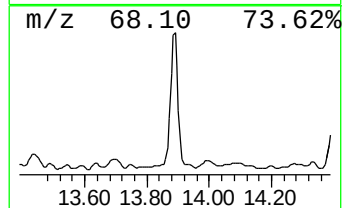
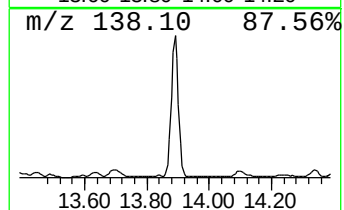
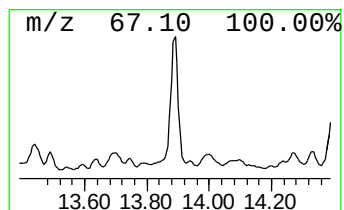
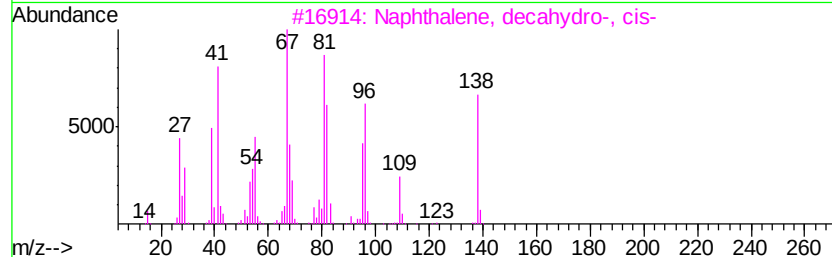
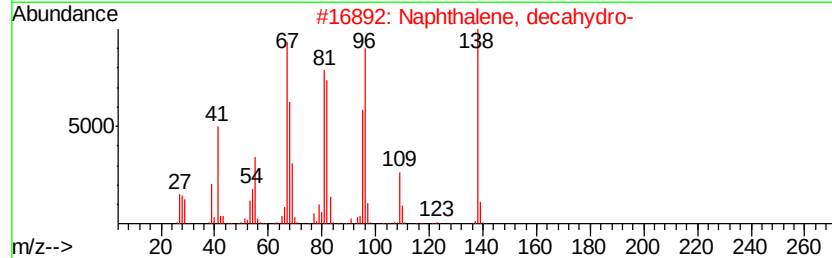
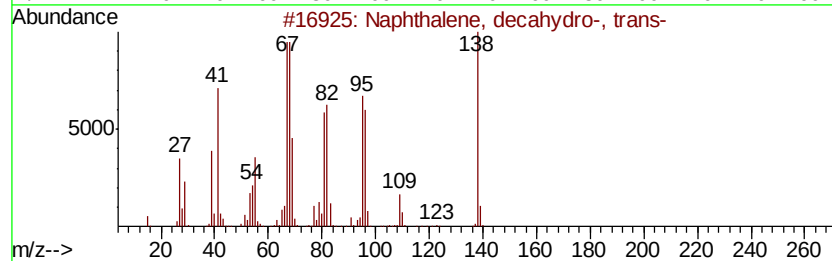
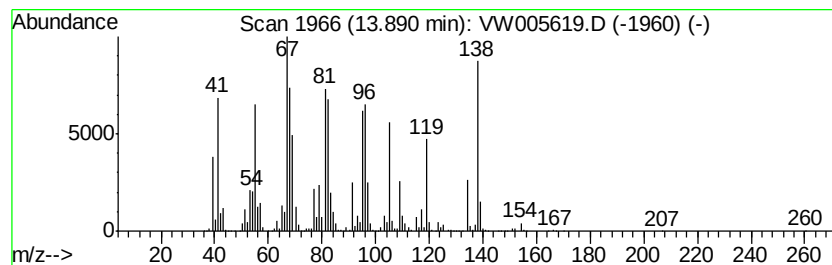
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Peak Number 4 Naphthalene, decahydro-, tr... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	1300.03 ug/l	14769700	1,4-Dichlorobenzene-d4	13.57
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, decahydro-, trans-	138 C10H18	000493-02-7 96
2		Naphthalene, decahydro-	138 C10H18	000091-17-8 96
3		Naphthalene, decahydro-, cis-	138 C10H18	000493-01-6 70
4		Cyclohexane, 1-methyl-3-(1-methy...	138 C10H18	024399-15-3 64
5		Spiro[4.5]decane	138 C10H18	000176-63-6 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW092518\
Data File : VW005619.D
Acq On : 25 Sep 2018 18:44
Operator : VA/AP
Sample : J5066-09
Misc : 4.83G/5ML/MSVOA W/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EP1-Q1-H4

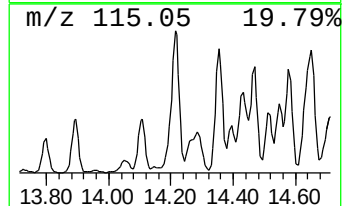
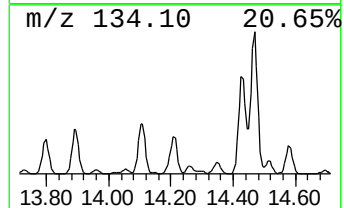
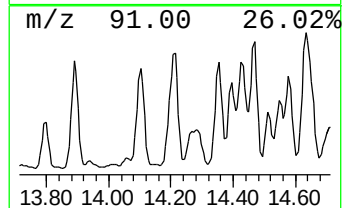
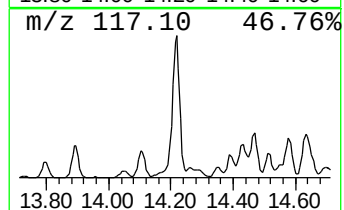
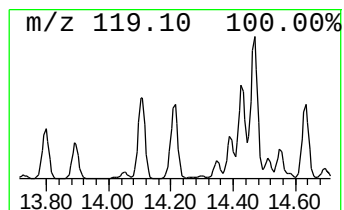
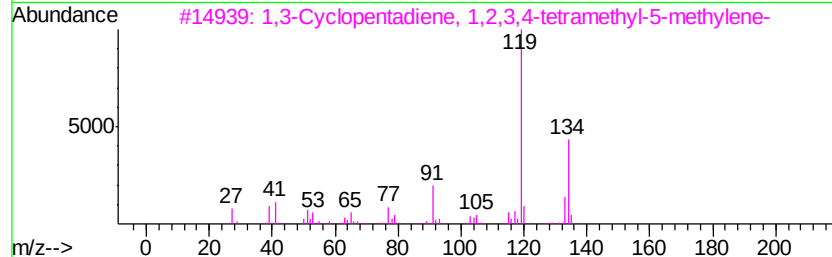
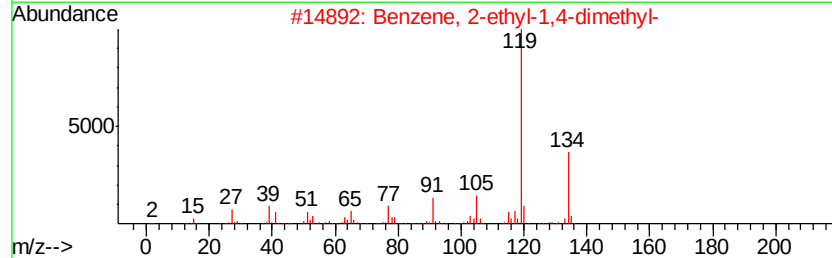
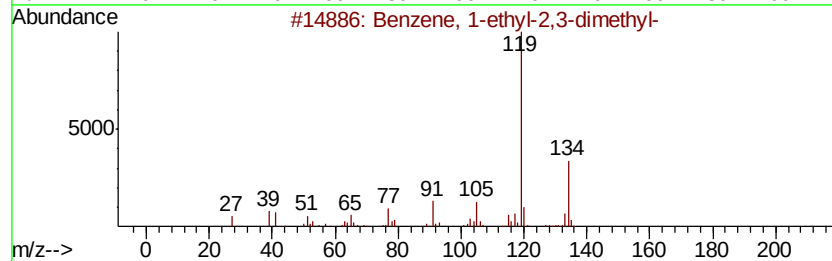
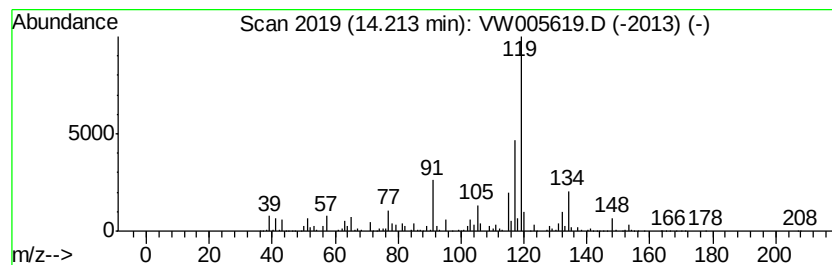
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W092118S.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	465.75 ug/l	5291420	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	76
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	76
3		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	70
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	64
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW092518\
Data File : VW005619.D
Acq On : 25 Sep 2018 18:44
Operator : VA/AP
Sample : J5066-09
Misc : 4.83G/5ML/MSVOA W/SOIL
ALS Vial : 20 Sample Multiplier: 1

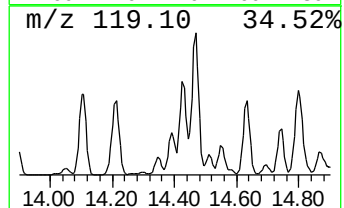
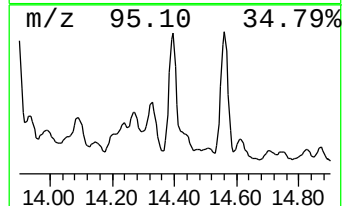
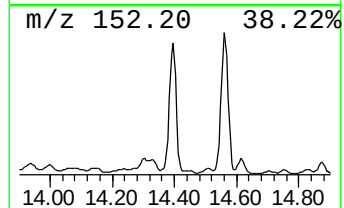
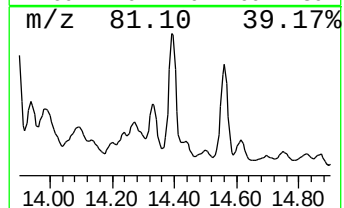
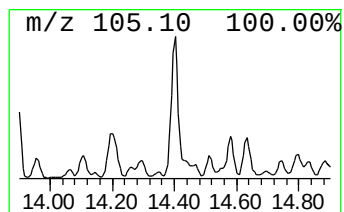
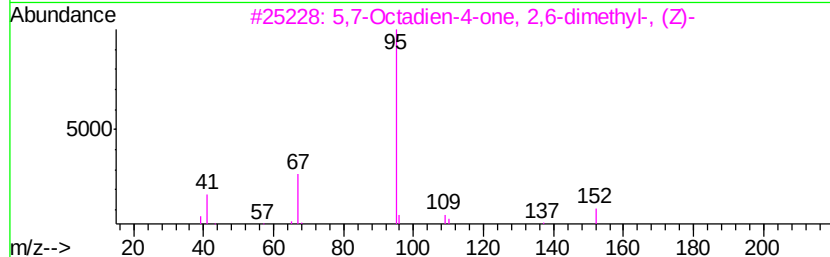
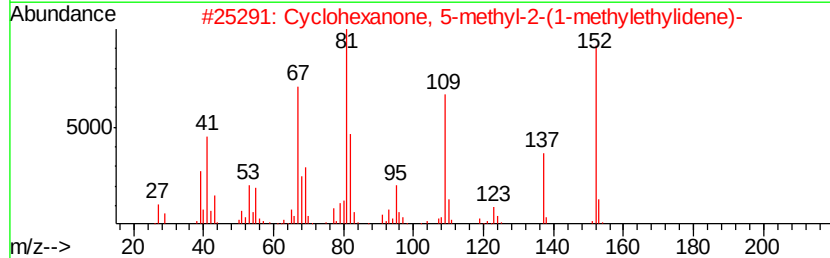
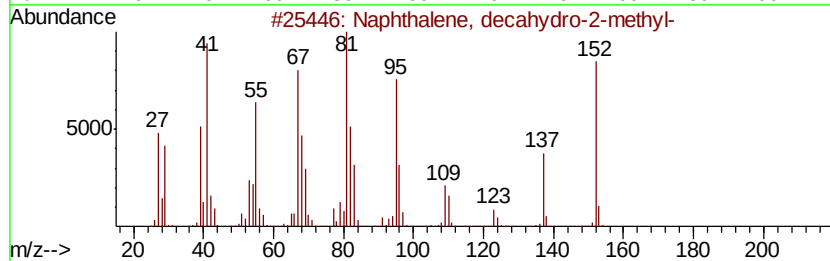
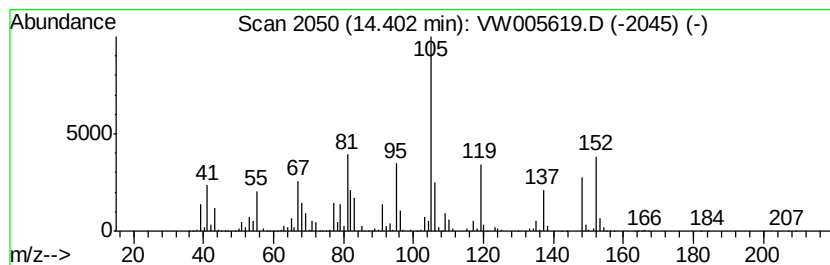
Instrument :
MSVOA_W
ClientSampleId :
EP1-Q1-H4

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W092118S.M
Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, decahydro-2-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.40	493.61 ug/l	5607890	1,4-Dichlorobenzene-d4	13.57
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, decahydro-2-methyl-	152 C11H20	002958-76-1 93
2		Cyclohexanone, 5-methyl-2-(1-met...	152 C10H16O	015932-80-6 50
3		5,7-Octadien-4-one, 2,6-dimethyl...	152 C10H16O	003588-18-9 43
4		2H-Inden-2-one, octahydro-3a-met...	152 C10H16O	013351-29-6 43
5		1-Adamantanol	152 C10H16O	000768-95-6 42



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW092518\
Data File : VW005619.D
Acq On : 25 Sep 2018 18:44
Operator : VA/AP
Sample : J5066-09
Misc : 4.83G/5ML/MSVOA W/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EP1-Q1-H4

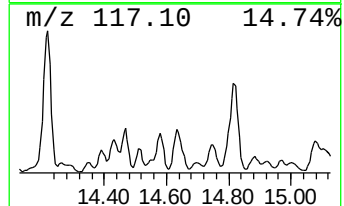
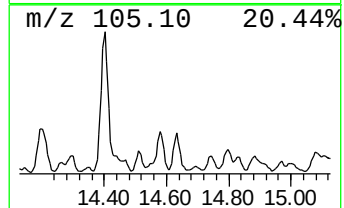
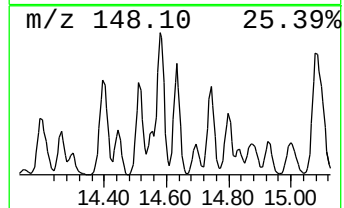
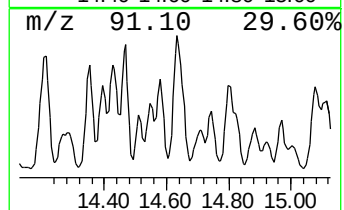
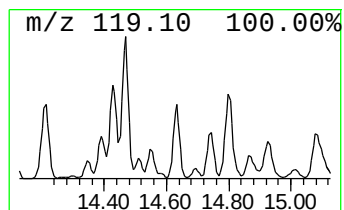
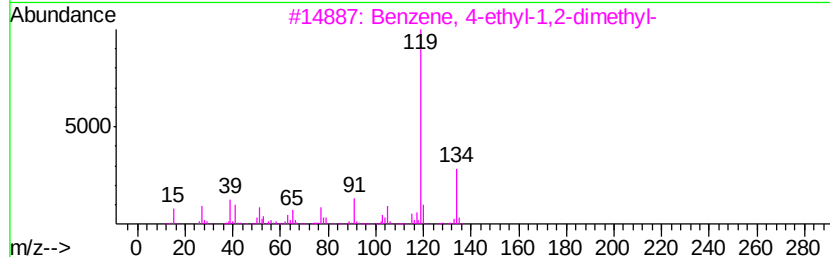
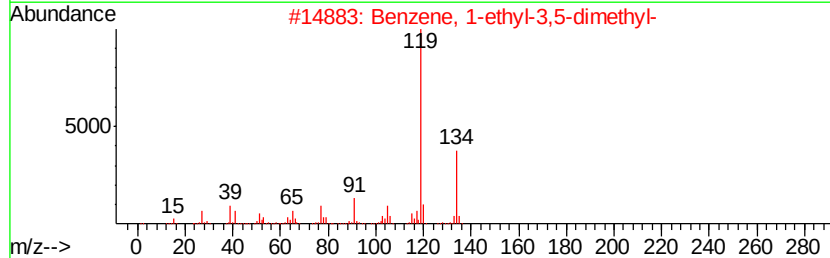
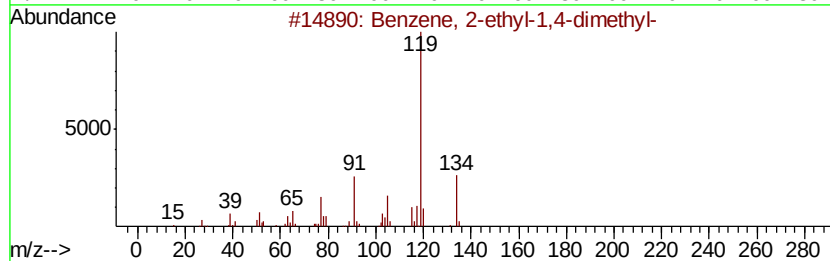
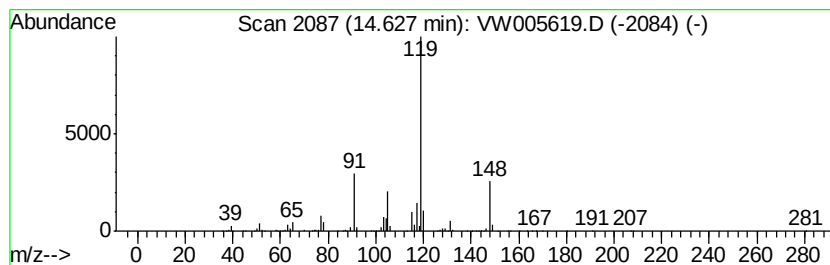
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W092118S.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.63	420.96 ug/l	4782520	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	59
2		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	59
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	59
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	59
5		Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW092518\
Data File : VW005619.D
Acq On : 25 Sep 2018 18:44
Operator : VA/AP
Sample : J5066-09
Misc : 4.83G/5ML/MSVOA W/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EP1-Q1-H4

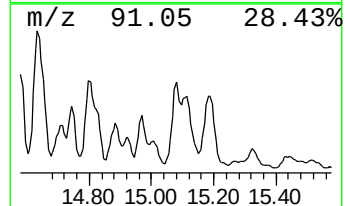
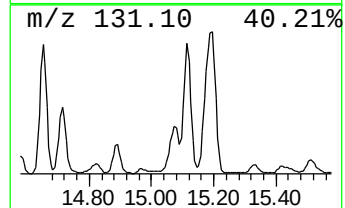
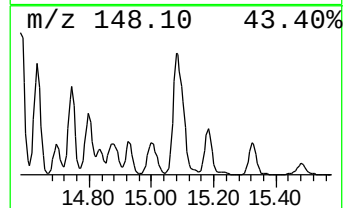
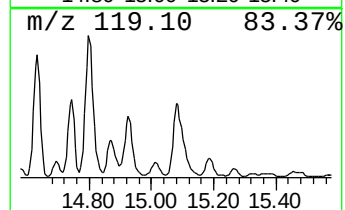
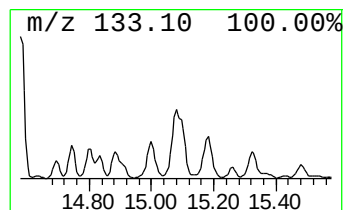
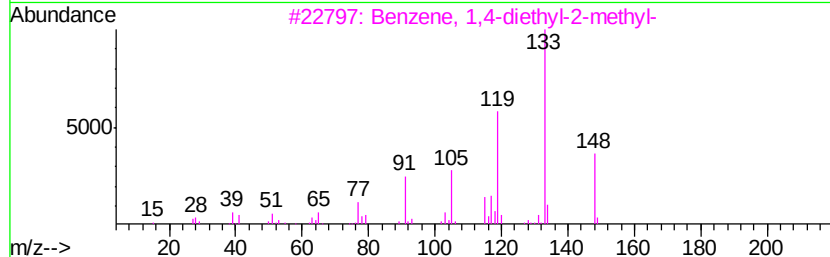
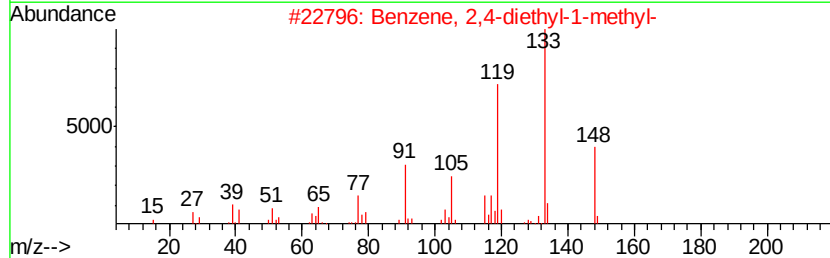
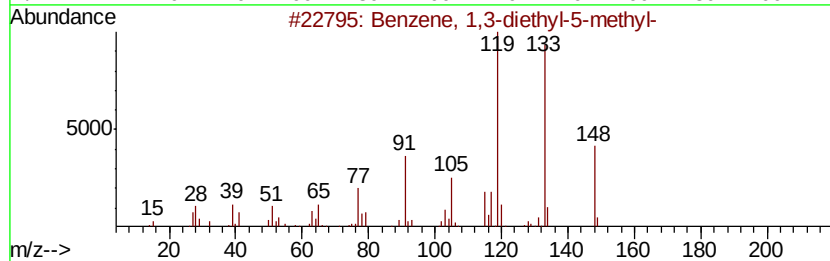
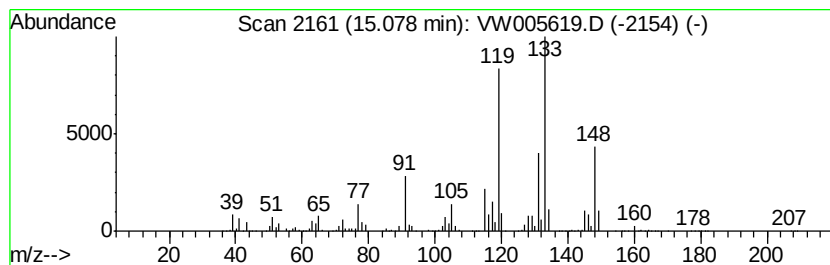
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W092118S.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.08	425.68 ug/l	4836220	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	93
2		Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	81
3		Benzene, 1,4-diethyl-2-methyl-	148	C11H16	013632-94-5	81
4		Benzene, pentamethyl-	148	C11H16	000700-12-9	46
5		3,4-Dimethylcumene	148	C11H16	1000370-34-1	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW092518\
Data File : VW005619.D
Acq On : 25 Sep 2018 18:44
Operator : VA/AP
Sample : J5066-09
Misc : 4.83G/5ML/MSVOA W/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EP1-Q1-H4

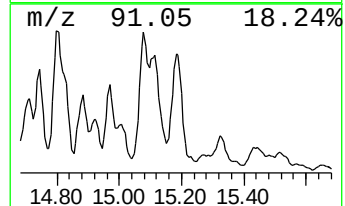
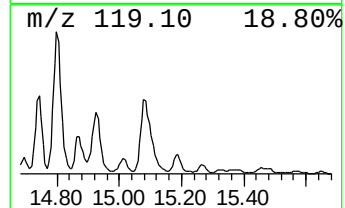
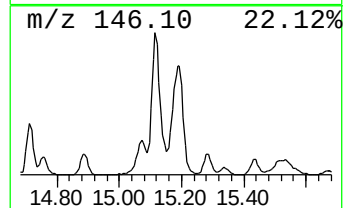
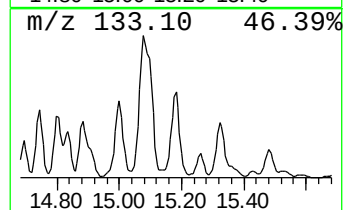
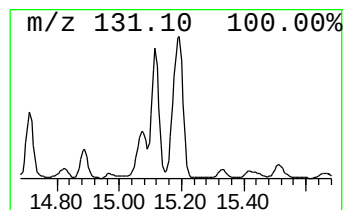
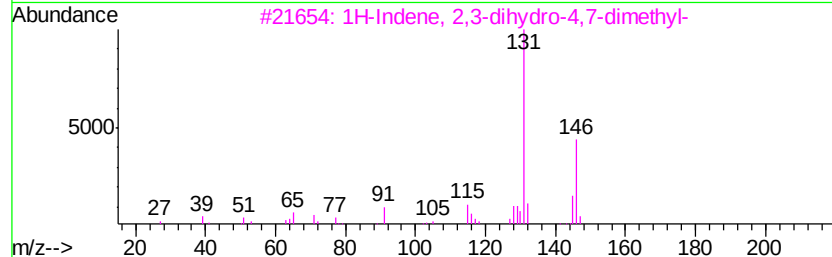
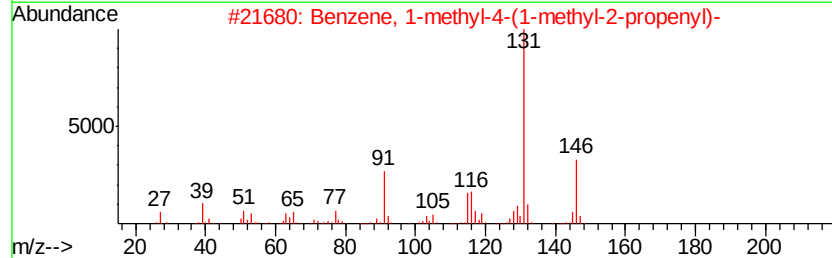
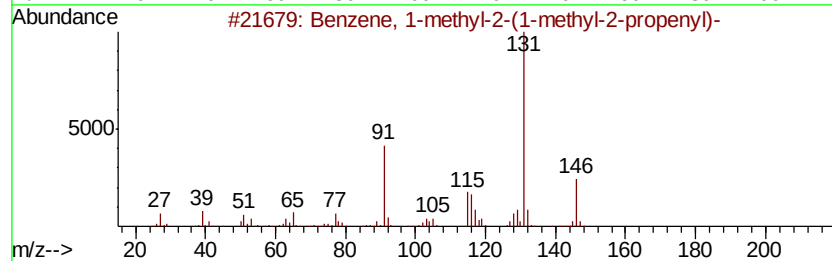
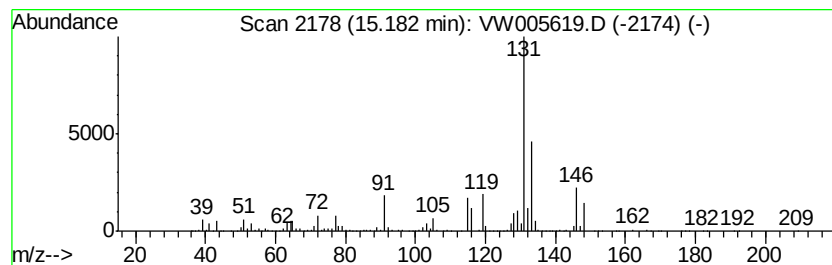
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W092118S.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-methyl-2-(1-meth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.18	447.44 ug/l	5083410	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	70
2		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	70
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	64
4		1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	64
5		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW092518\
Data File : VW005619.D
Acq On : 25 Sep 2018 18:44
Operator : VA/AP
Sample : J5066-09
Misc : 4.83G/5ML/MSVOA_W/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EP1-Q1-H4

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W092118S.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown12.87	12.87	411.0	ug/l	4669030	4	13.57	568053	50.0
3-(1H-Imidazol-4-...	13.44	434.7	ug/l	4938850	4	13.57	568053	50.0
Naphthalene, deca...	13.89	1300.0	ug/l	14769700	4	13.57	568053	50.0
Benzene, 1-ethyl-...	14.21	465.8	ug/l	5291420	4	13.57	568053	50.0
Naphthalene, deca...	14.40	493.6	ug/l	5607890	4	13.57	568053	50.0
Benzene, 2-ethyl-...	14.63	421.0	ug/l	4782520	4	13.57	568053	50.0
Benzene, 1,3-diet...	15.08	425.7	ug/l	4836220	4	13.57	568053	50.0
Benzene, 1-methyl...	15.18	447.4	ug/l	5083410	4	13.57	568053	50.0