

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW020719\
 Data File : VW008551.D
 Acq On : 07 Feb 2019 14:35
 Operator : SY/VA
 Sample : K1390-01
 Misc : 5.55G/5ML/MSVOA W/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 SB-01(12-13)

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\82W012419S.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	3.032	175	185	203	rBV	2962994	8073885	1.39%	0.058%
2	3.392	233	244	263	rBV	2628883	7635397	1.31%	0.055%
3	4.111	348	362	392	rVB	2842842	12172094	2.10%	0.088%
4	4.958	473	501	549	rBV5	67276341	367431433	63.24%	2.661%
5	5.440	561	580	622	rBV3	55704730	264518622	45.53%	1.916%
6	5.940	646	662	686	rBV	25884780	96466981	16.60%	0.699%
7	6.732	773	792	805	rBV	13464175	67101338	11.55%	0.486%
8	6.891	805	818	852	rVV3	58863020	313314626	53.93%	2.269%
9	7.184	853	866	895	rVB	3794905	18104001	3.12%	0.131%
10	7.720	943	954	972	rVB2	16766057	78740229	13.55%	0.570%
11	8.067	1003	1011	1021	rVB4	104819202	405038620	69.72%	2.934%
12	8.183	1021	1030	1033	rBV8	132385778	405611162	69.81%	2.938%
13	8.452	1063	1074	1079	rBV4	120040999	418040546	71.95%	3.028%
14	8.592	1093	1097	1107	rVB3	68035976	179320054	30.86%	1.299%
15	8.708	1107	1116	1129	rVB4	102314834	273855733	47.14%	1.984%
16	8.836	1129	1137	1149	rBV2	34691310	110592791	19.04%	0.801%
17	8.939	1150	1154	1159	rVV5	6157018	16220135	2.79%	0.117%
18	9.000	1159	1164	1172	rVV	16137673	36214470	6.23%	0.262%
19	9.079	1172	1177	1181	rVV	3619992	8717993	1.50%	0.063%
20	9.165	1182	1191	1205	rVB2	27084043	107852625	18.56%	0.781%
21	9.531	1241	1251	1257	rBV2	62823746	155849870	26.82%	1.129%
22	9.628	1257	1267	1279	rVB6	126837489	458894851	78.99%	3.324%
23	9.781	1282	1292	1305	rBV7	134786069	564652255	97.19%	4.090%
24	9.921	1306	1315	1319	rBV6	112958586	375580677	64.65%	2.720%
25	10.122	1341	1348	1352	rBV6	125335214	381579705	65.68%	2.764%
26	10.281	1366	1374	1380	rBV5	126740369	322355369	55.48%	2.335%
27	10.354	1380	1386	1391	rBV4	157032998	462414019	79.59%	3.349%
28	10.470	1397	1405	1408	rBV6	94707108	237355590	40.85%	1.719%
29	10.524	1409	1414	1421	rVB8	161329722	458770611	78.96%	3.323%
30	10.585	1421	1424	1429	rVB3	31712636	50122624	8.63%	0.363%
31	10.640	1430	1433	1436	rVB3	14394020	17973316	3.09%	0.130%
32	10.689	1436	1441	1448	rVB4	79830760	186211433	32.05%	1.349%
33	10.780	1448	1456	1464	rBV10	164531424	580989021	100.00%	4.208%
34	10.866	1465	1470	1479	rVB6	123547700	355005943	61.10%	2.571%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW020719\
 Data File : VW008551.D
 Acq On : 07 Feb 2019 14:35
 Operator : SY/VA
 Sample : K1390-01
 Misc : 5.55G/5ML/MSVOA W/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 SB-01(12-13)

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\82W012419S.M
 Title : SW846 8260

35	10.969	1479	1487	1495	rBV9	114939012	456963612	78.65%	3.310%
36	11.360	1547	1551	1556	rBV6	54679024	138414309	23.82%	1.003%
37	11.536	1575	1580	1583	rBV5	94941896	195737055	33.69%	1.418%
38	11.756	1612	1616	1619	rBV3	48023544	84122402	14.48%	0.609%
39	11.799	1620	1623	1625	rBV2	31648760	45569061	7.84%	0.330%
40	12.036	1657	1662	1664	rBV4	31738592	61353429	10.56%	0.444%
41	12.085	1665	1670	1675	rVB7	86634360	188811237	32.50%	1.368%
42	12.353	1707	1714	1717	rBV5	118564488	284449640	48.96%	2.060%
43	12.683	1763	1768	1771	rVV5	67927635	143828360	24.76%	1.042%
44	12.719	1771	1774	1780	rVB5	76524216	133600711	23.00%	0.968%
45	12.811	1785	1789	1794	rVB5	81752144	161848415	27.86%	1.172%
46	12.878	1795	1800	1804	rBV5	65640448	129388715	22.27%	0.937%
47	13.006	1818	1821	1827	rVB5	76104136	129150559	22.23%	0.935%
48	13.091	1830	1835	1839	rBV4	82243384	151951563	26.15%	1.101%
49	13.146	1839	1844	1846	rBV5	73528032	120156058	20.68%	0.870%
50	13.182	1847	1850	1856	rVB3	87883552	149759869	25.78%	1.085%
51	13.243	1856	1860	1866	rVB7	55859776	109122199	18.78%	0.790%
52	13.304	1866	1870	1875	rVB4	71100056	117690893	20.26%	0.852%
53	13.365	1875	1880	1887	rBV8	63490472	178692444	30.76%	1.294%
54	13.457	1891	1895	1899	rVB5	70599048	108345087	18.65%	0.785%
55	13.499	1899	1902	1905	rBV4	65532648	86436384	14.88%	0.626%
56	13.573	1911	1914	1918	rVB5	79039656	111946482	19.27%	0.811%
57	13.640	1921	1925	1931	rVB4	108025168	189686962	32.65%	1.374%
58	13.725	1931	1939	1941	rBV5	68744192	146811610	25.27%	1.063%
59	13.762	1941	1945	1949	rVB3	93649632	164654962	28.34%	1.193%
60	13.823	1949	1955	1961	rVB7	84781088	232947091	40.09%	1.687%
61	13.890	1961	1966	1970	rBV4	160601088	293427425	50.50%	2.125%
62	14.024	1985	1988	1990	rBV2	36515848	53004525	9.12%	0.384%
63	14.103	1997	2001	2006	rVB3	84199744	135098834	23.25%	0.979%
64	14.207	2013	2018	2024	rVB4	94374892	173849198	29.92%	1.259%
65	14.298	2024	2033	2038	rBV6	63531036	170532787	29.35%	1.235%
66	14.359	2039	2043	2045	rVV3	21017253	37294270	6.42%	0.270%
67	14.396	2045	2049	2052	rVV3	92812133	151609263	26.10%	1.098%
68	14.432	2052	2055	2058	rVV2	62432081	89735718	15.45%	0.650%
69	14.463	2058	2060	2063	rVB2	30957040	34462050	5.93%	0.250%
70	14.505	2063	2067	2071	rBV2	77151016	108648175	18.70%	0.787%
71	14.554	2071	2075	2080	rVB4	44425696	87805001	15.11%	0.636%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW020719\
 Data File : VW008551.D
 Acq On : 07 Feb 2019 14:35
 Operator : SY/VA
 Sample : K1390-01
 Misc : 5.55G/5ML/MSVOA W/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 SB-01(12-13)

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W012419S.M
 Title : SW846 8260

72	14.609	2080	2084	2086	rBV2	24896740	34062222	5.86%	0.247%
73	14.646	2087	2090	2095	rVB2	61129306	91153146	15.69%	0.660%
74	14.713	2095	2101	2110	rVB5	56209284	191846297	33.02%	1.390%
75	14.798	2110	2115	2117	rBV3	49294436	87270777	15.02%	0.632%
76	14.859	2122	2125	2134	rVB4	34797266	61171774	10.53%	0.443%
77	15.030	2148	2153	2155	rBV3	13065254	23337993	4.02%	0.169%
78	15.066	2155	2159	2164	rVB3	27651310	50010851	8.61%	0.362%
79	15.182	2172	2178	2184	rVB3	24896252	58490006	10.07%	0.424%
80	15.334	2198	2203	2211	rVB6	6065667	14396699	2.48%	0.104%
81	15.426	2212	2218	2222	rBV2	6992342	14036978	2.42%	0.102%
82	15.475	2222	2226	2230	rVV2	4657632	6970950	1.20%	0.050%
83	15.658	2249	2256	2263	rVB3	3760334	8319783	1.43%	0.060%
84	15.865	2286	2290	2300	rVB	2776037	5888647	1.01%	0.043%

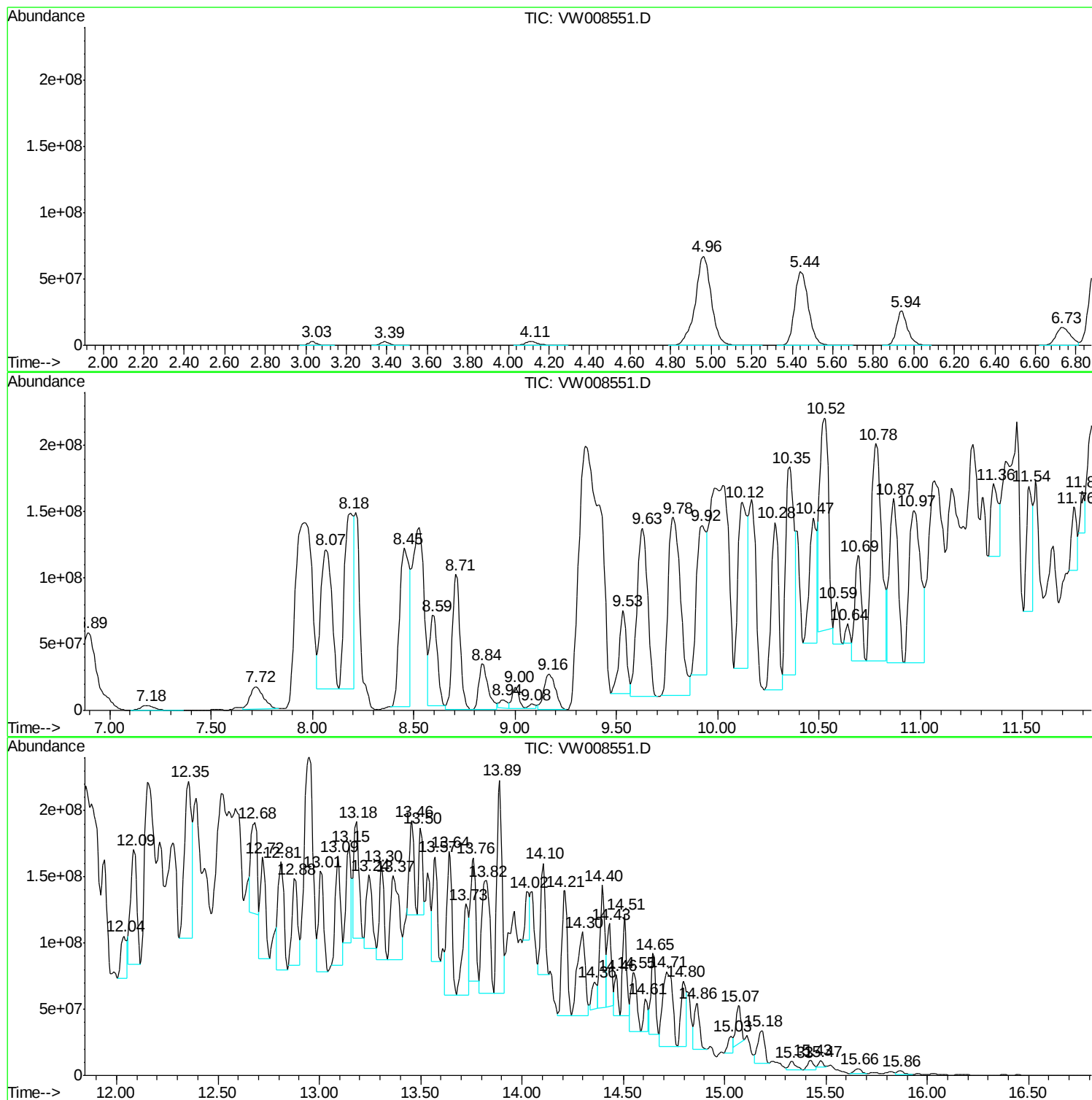
Sum of corrected areas: 13806640497

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW020719\
Data File : VW008551.D
Acq On : 07 Feb 2019 14:35
Operator : SY/VA
Sample : K1390-01
Misc : 5.55G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
SB-01(12-13)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W012419S.M
Quant Title : SW846 8260

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW020719\
Data File : VW008551.D
Acq On : 07 Feb 2019 14:35
Operator : SY/VA
Sample : K1390-01
Misc : 5.55G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
SB-01(12-13)

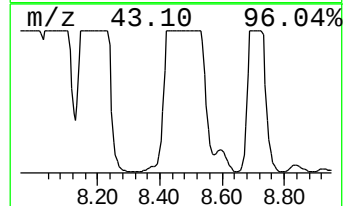
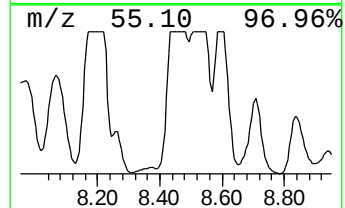
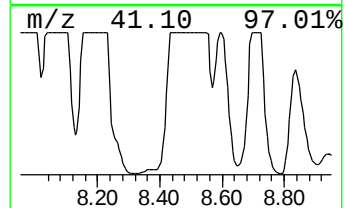
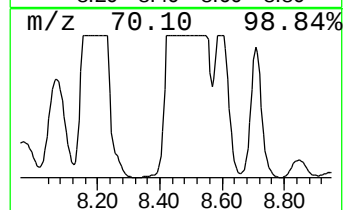
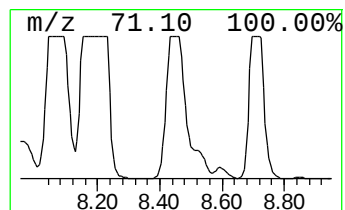
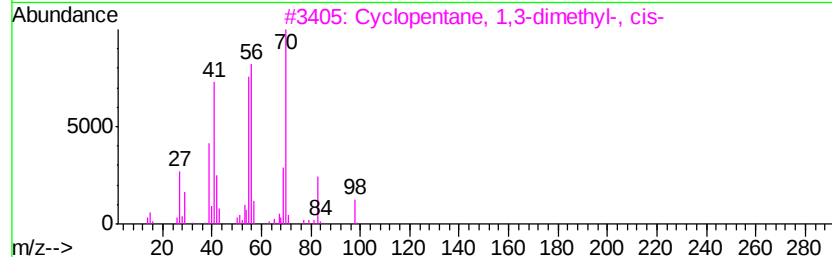
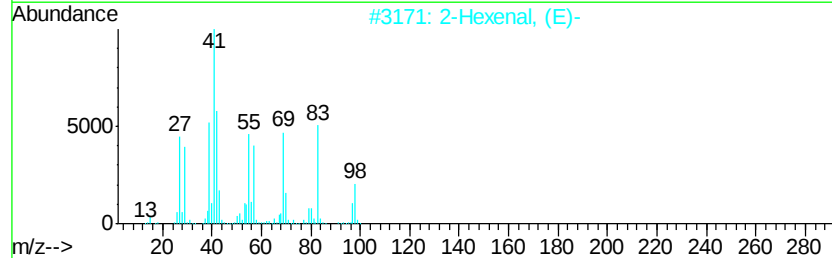
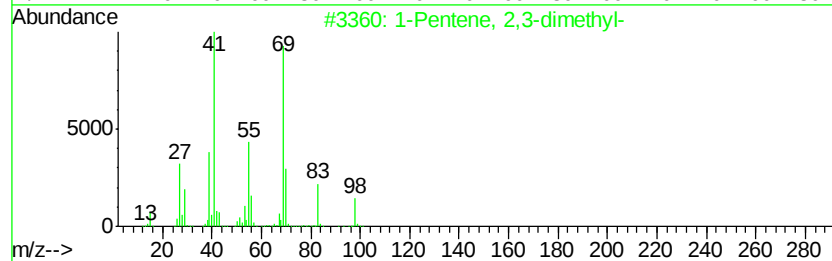
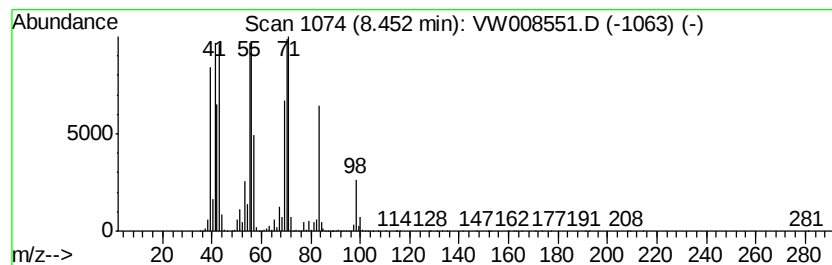
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W012419S.M
Quant Title : SW846 8260

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

Peak Number 1 1-Pentene, 2,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.45	189.00 ug/l	418041000	1,4-Difluorobenzene	8.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Pentene, 2,3-dimethyl-	98	C7H14	003404-72-6	58
2		2-Hexenal, (E)-	98	C6H10O	006728-26-3	52
3		Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	49
4		Cycloheptane	98	C7H14	000291-64-5	38
5		Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	25



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW020719\
Data File : VW008551.D
Acq On : 07 Feb 2019 14:35
Operator : SY/VA
Sample : K1390-01
Misc : 5.55G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
SB-01(12-13)

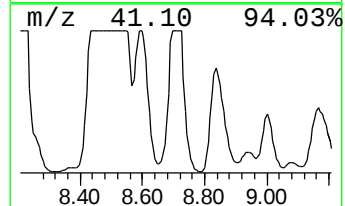
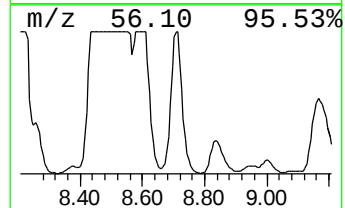
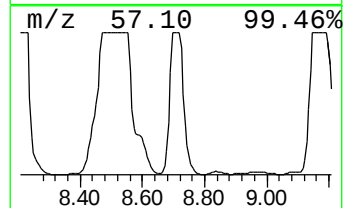
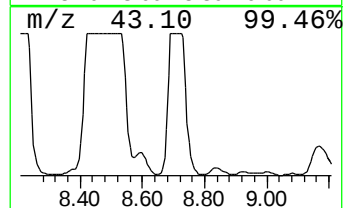
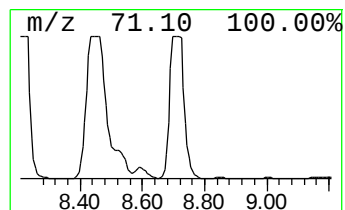
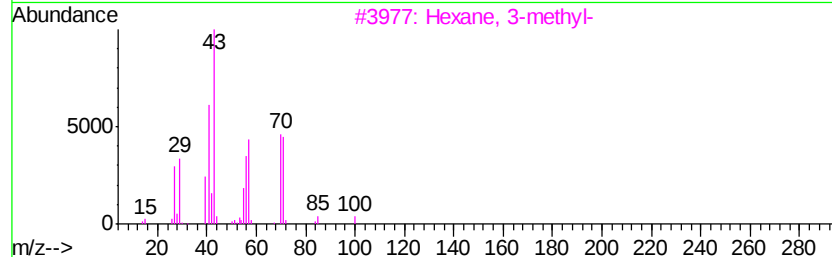
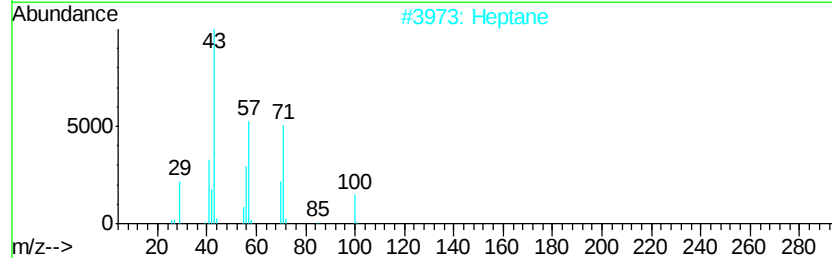
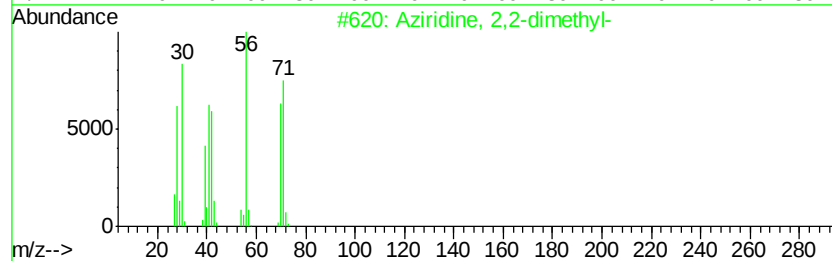
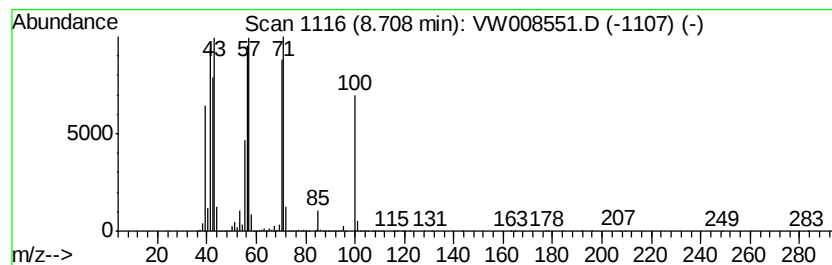
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W012419S.M
Quant Title : SW846 8260

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

Peak Number 2 Aziridine, 2,2-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.71	123.81 ug/l	273856000	1,4-Difluorobenzene	8.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Aziridine, 2,2-dimethyl-	71	C4H9N	002658-24-4	52
2		Heptane	100	C7H16	000142-82-5	50
3		Hexane, 3-methyl-	100	C7H16	000589-34-4	50
4		Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	38
5		Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW020719\
Data File : VW008551.D
Acq On : 07 Feb 2019 14:35
Operator : SY/VA
Sample : K1390-01
Misc : 5.55G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
SB-01(12-13)

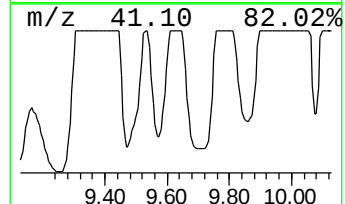
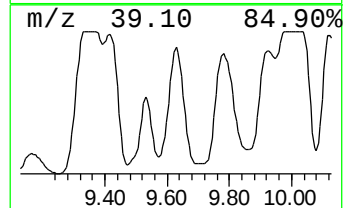
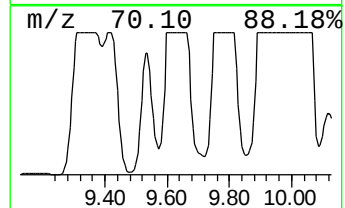
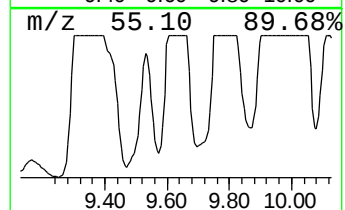
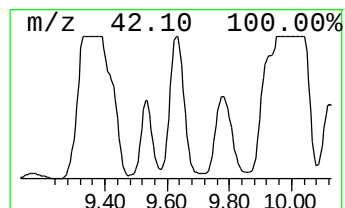
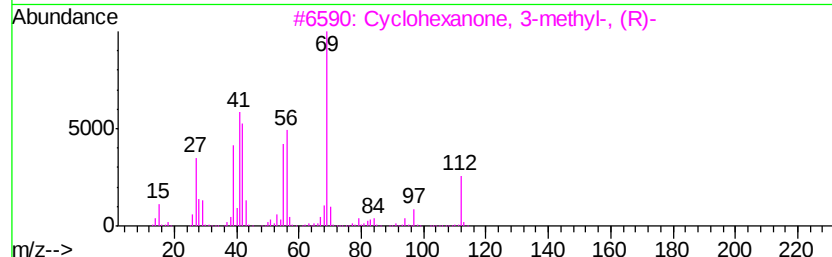
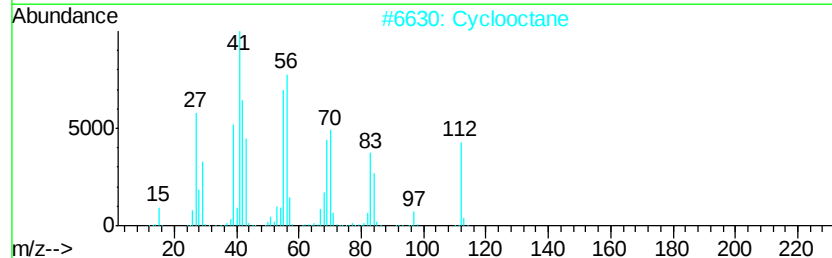
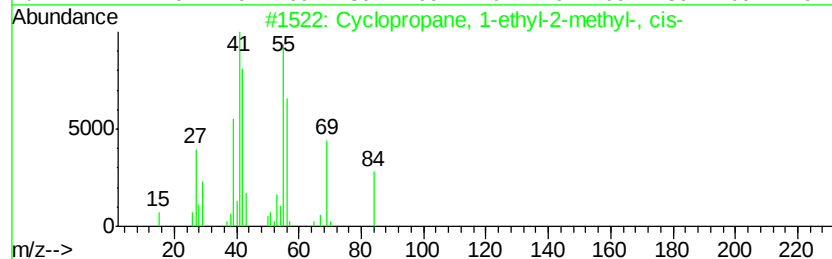
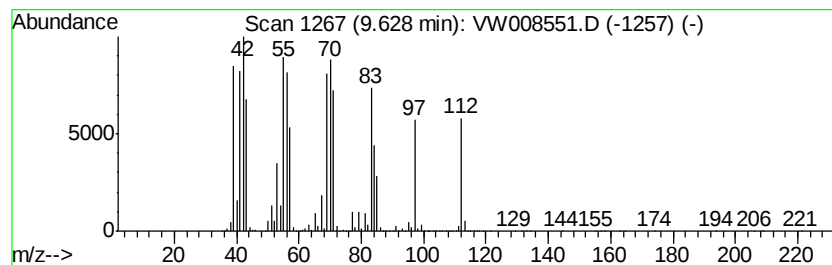
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W012419S.M
Quant Title : SW846 8260

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

Peak Number 3 Cyclopropane, 1-ethyl-2-met... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.63	207.47 ug/l	458895000	1,4-Difluorobenzene	8.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopropane, 1-ethyl-2-methyl-, ...	84	C6H12	019781-68-1	58
2		Cyclooctane	112	C8H16	000292-64-8	50
3		Cyclohexanone, 3-methyl-, (R)-	112	C7H12O	013368-65-5	49
4		1-Pentene, 3-methyl-	84	C6H12	000760-20-3	46
5		Cyclohexane, 1,2-dimethyl-, cis-	112	C8H16	002207-01-4	27



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW020719\
Data File : VW008551.D
Acq On : 07 Feb 2019 14:35
Operator : SY/VA
Sample : K1390-01
Misc : 5.55G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
SB-01(12-13)

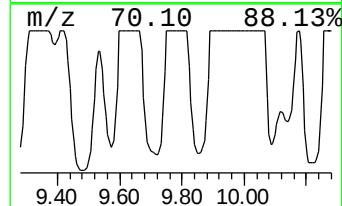
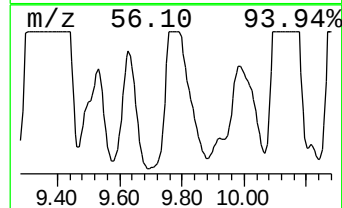
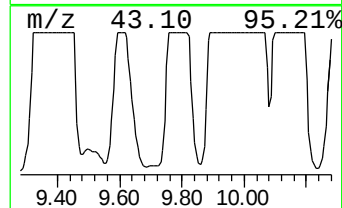
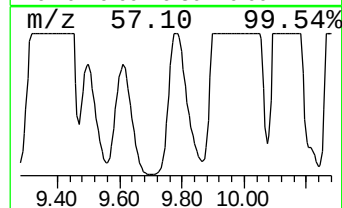
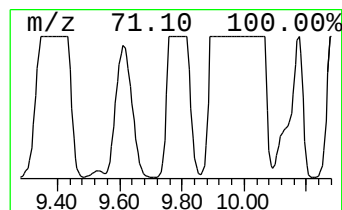
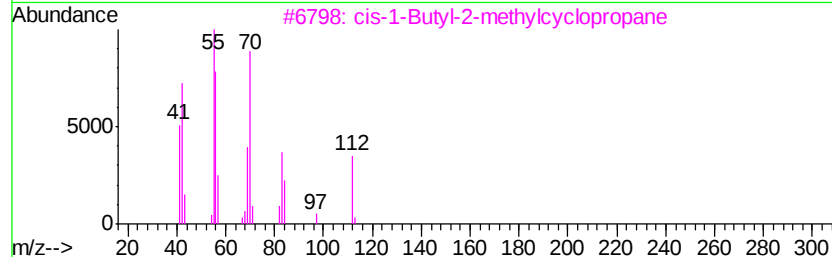
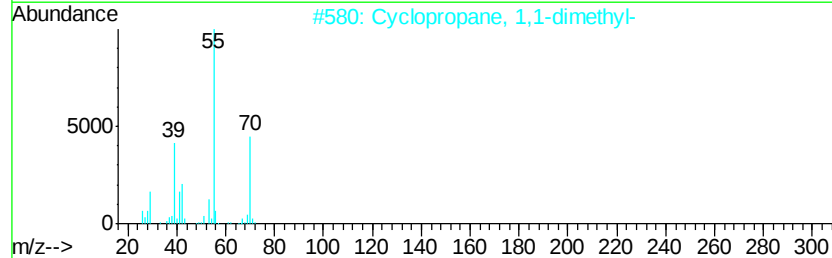
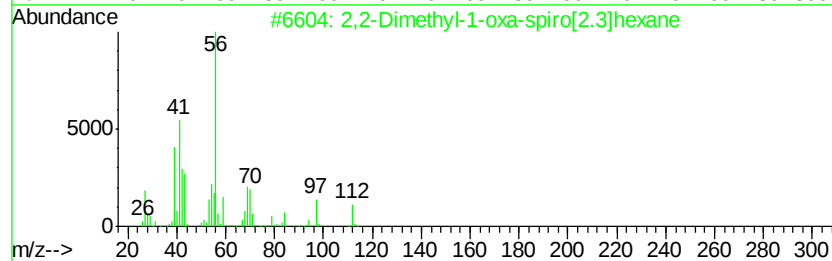
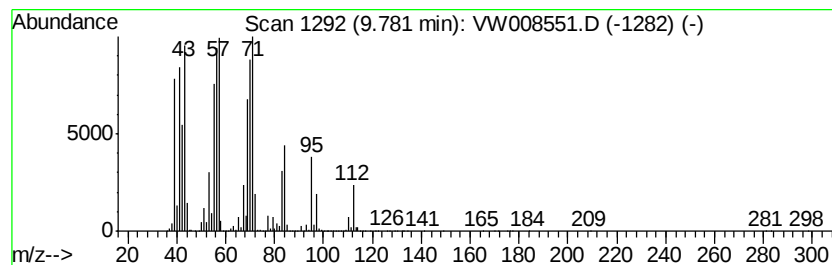
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W012419S.M
Quant Title : SW846 8260

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

Peak Number 4 2,2-Dimethyl-1-oxa-spiro[2.3]hexane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.78	255.28 ug/l	564652000	1,4-Difluorobenzene	8.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2-Dimethyl-1-oxa-spiro[2.3]hexane	112	C7H12O	1000194-04-4	50
2		Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	50
3		cis-1-Butyl-2-methylcyclopropane	112	C8H16	038851-69-3	46
4		Furan, 2,5-dihydro-3-methyl-	84	C5H8O	001708-31-2	46
5		Cyclooctane	112	C8H16	000292-64-8	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW020719\
Data File : VW008551.D
Acq On : 07 Feb 2019 14:35
Operator : SY/VA
Sample : K1390-01
Misc : 5.55G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
SB-01(12-13)

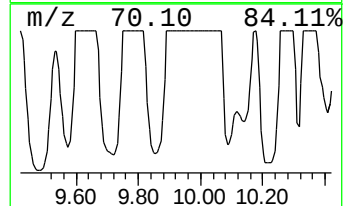
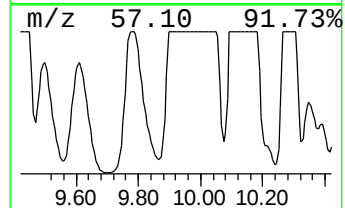
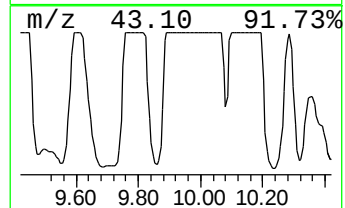
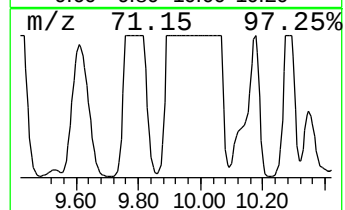
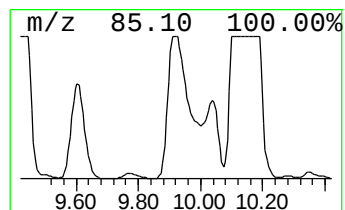
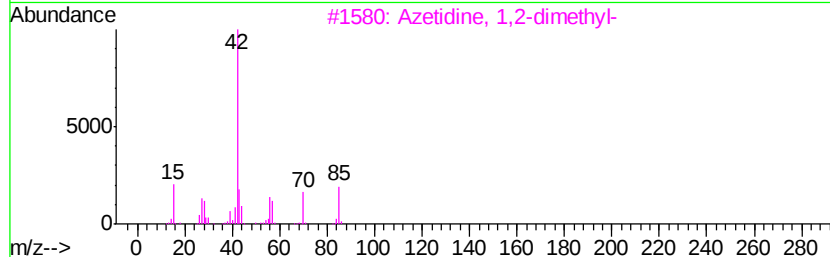
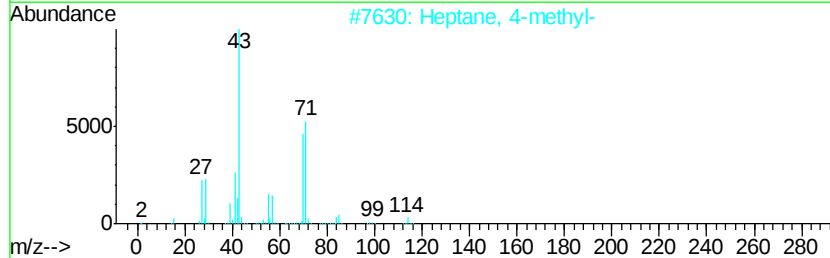
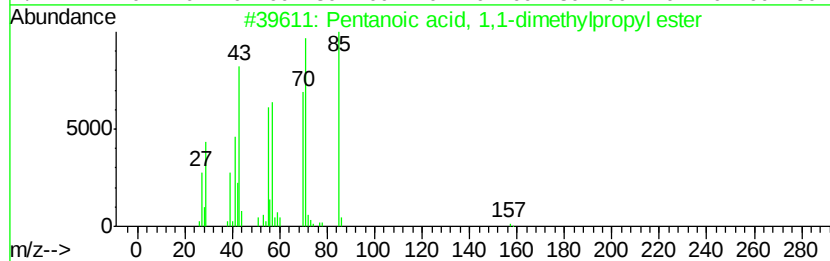
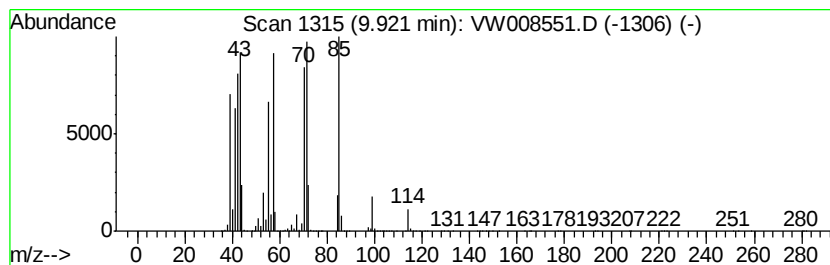
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W012419S.M
Quant Title : SW846 8260

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

Peak Number 5 Pentanoic acid, 1,1-dimethy... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.92	169.80 ug/l	375581000	1,4-Difluorobenzene	8.84

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentanoic acid, 1,1-dimethylprop...	172	C10H20O2	117421-32-6	53
2	Heptane, 4-methyl-	114	C8H18	000589-53-7	43
3	Azetidine, 1,2-dimethyl-	85	C5H11N	051764-32-0	38
4	2-Pentene	70	C5H10	000109-68-2	35
5	2-Pentene, (Z)-	70	C5H10	000627-20-3	35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW020719\
Data File : VW008551.D
Acq On : 07 Feb 2019 14:35
Operator : SY/VA
Sample : K1390-01
Misc : 5.55G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

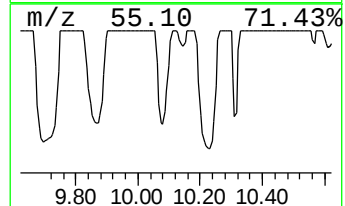
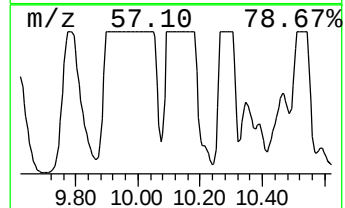
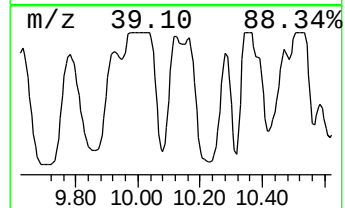
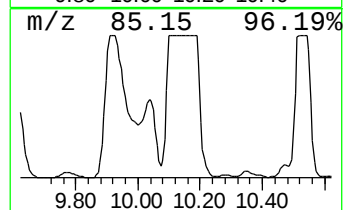
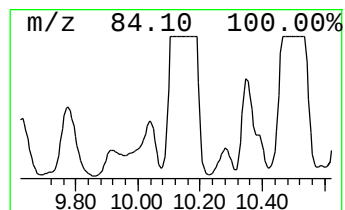
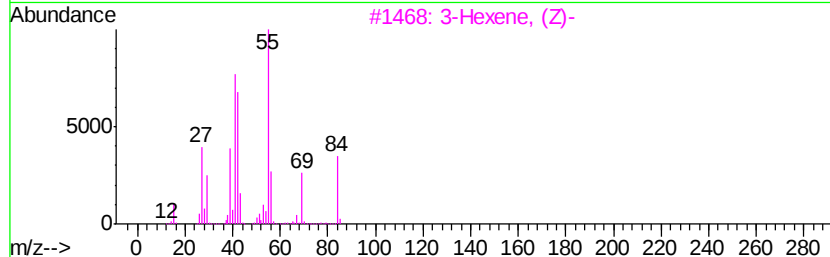
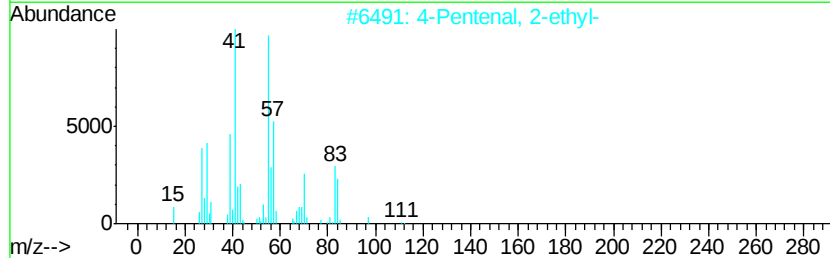
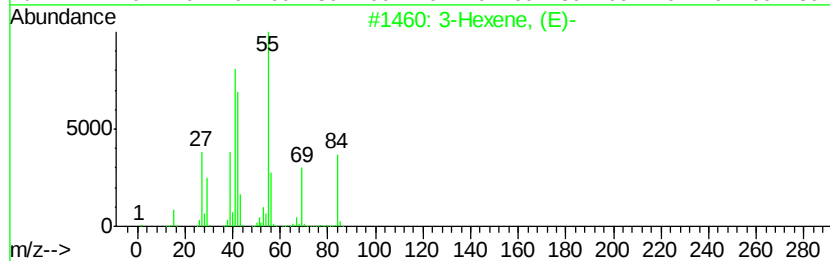
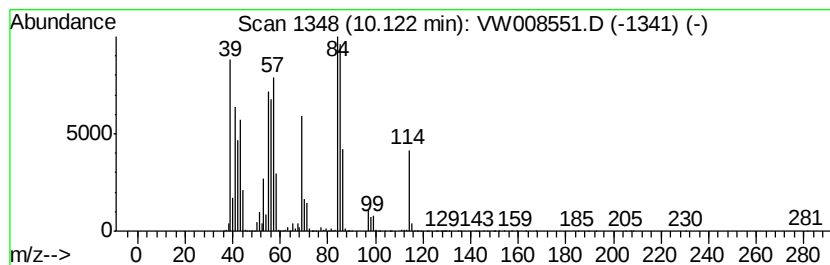
Instrument :
MSVOA_W
ClientSampleId :
SB-01(12-13)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W012419S.M
Quant Title : SW846 8260

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

Peak Number 6 3-Hexene, (E)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.12	172.52 ug/l	381580000	1,4-Difluorobenzene	8.84
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		3-Hexene, (E)-	84 C6H12	013269-52-8 50
2		4-Pentenal, 2-ethyl-	112 C7H12O	005204-80-8 50
3		3-Hexene, (Z)-	84 C6H12	007642-09-3 50
4		Pentane, 3-methylene-	84 C6H12	000760-21-4 40
5		Furan, 2,3-dihydro-3-methyl-	84 C5H8O	001708-27-6 38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW020719\
Data File : VW008551.D
Acq On : 07 Feb 2019 14:35
Operator : SY/VA
Sample : K1390-01
Misc : 5.55G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
SB-01(12-13)

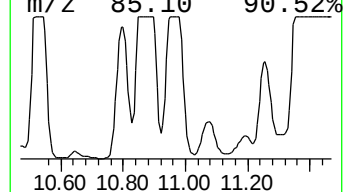
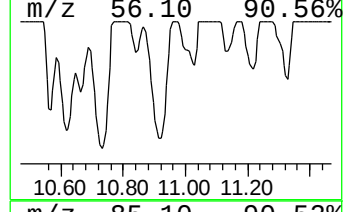
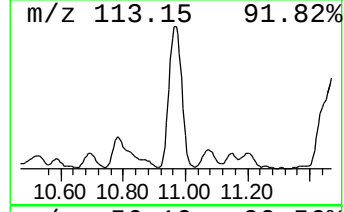
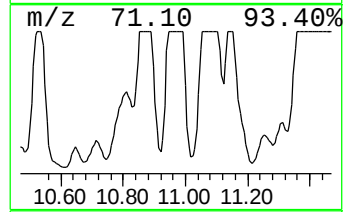
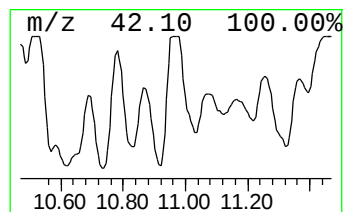
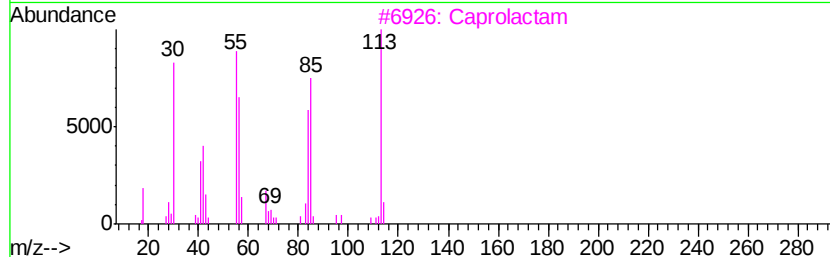
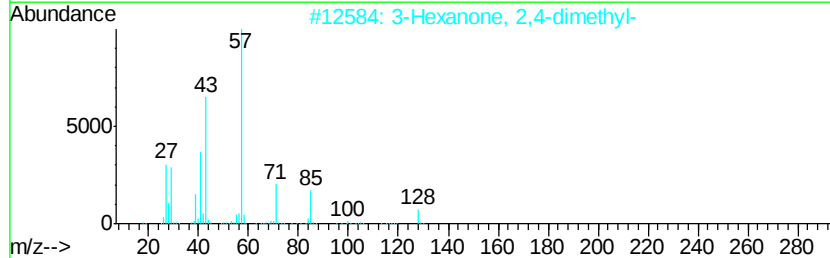
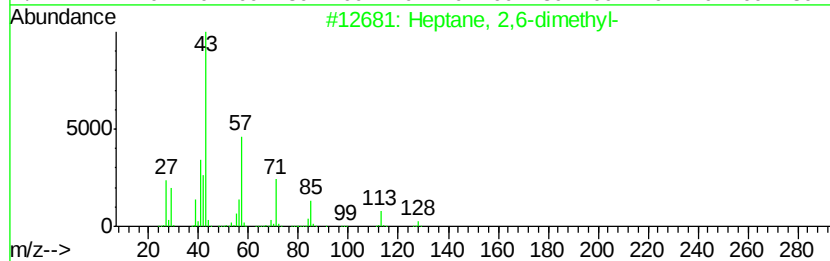
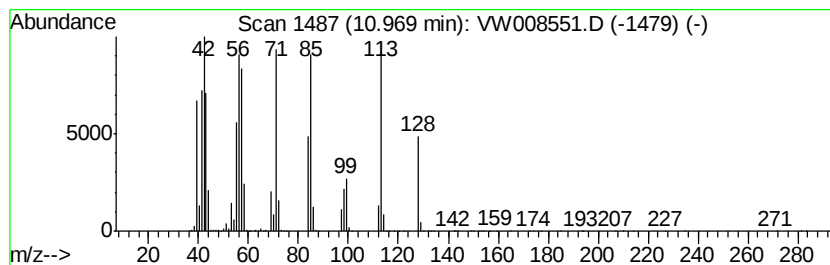
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W012419S.M
Quant Title : SW846 8260

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

Peak Number 8 unknown10.97 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.97	116.73 ug/l	456964000	Chlorobenzene-d5	11.63

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	49
2		3-Hexanone, 2,4-dimethyl-	128	C8H16O	018641-70-8	35
3		Caprolactam	113	C6H11NO	000105-60-2	27
4		Undecane, 2,7-dimethyl-	184	C13H28	017301-24-5	27
5		3,4-Dimethyl-isoxazol-5(4H)-one	113	C5H7NO2	015731-93-8	27



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW020719\
Data File : VW008551.D
Acq On : 07 Feb 2019 14:35
Operator : SY/VA
Sample : K1390-01
Misc : 5.55G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
SB-01(12-13)

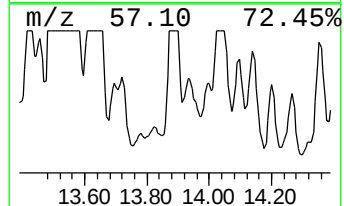
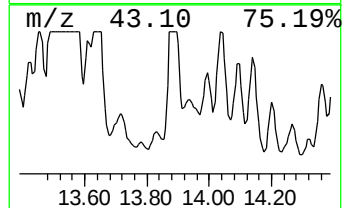
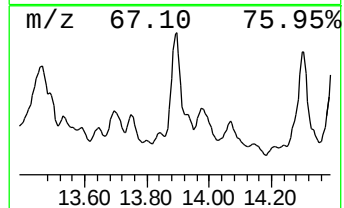
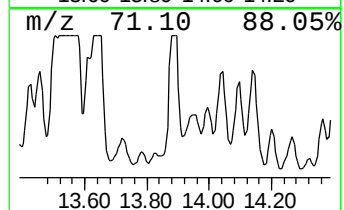
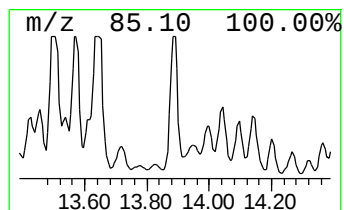
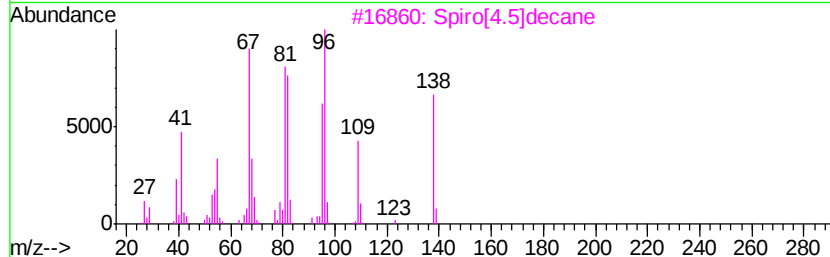
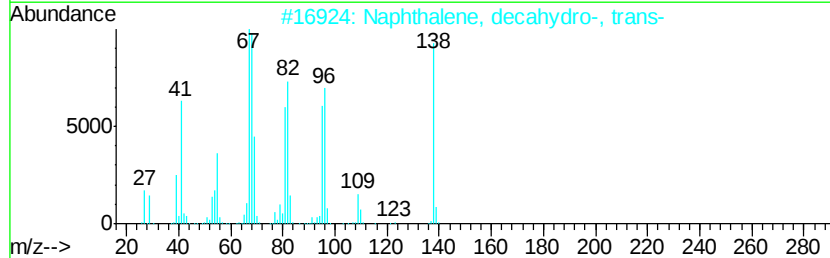
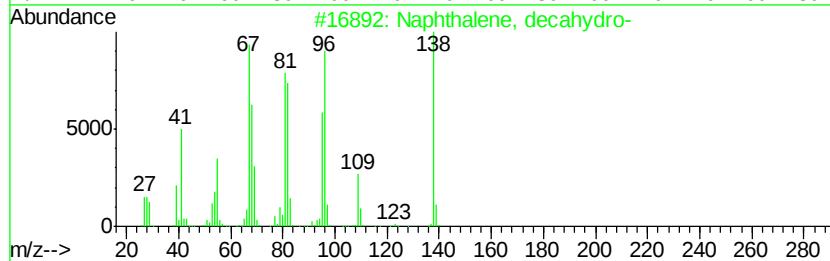
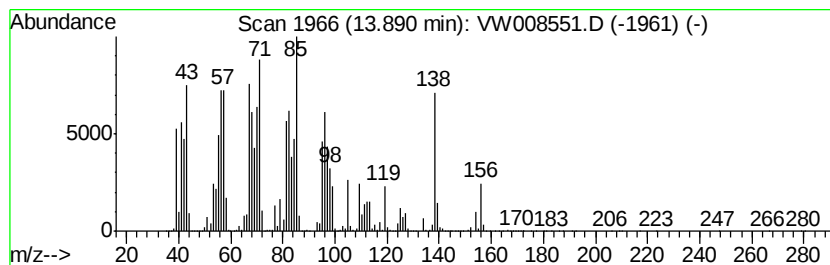
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W012419S.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, decahydro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	131.06 ug/l	293427000	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, decahydro-	138	C10H18	000091-17-8	94
2		Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	83
3		Spiro[4.5]decane	138	C10H18	000176-63-6	81
4		Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	80
5		trans-2-Dodecen-1-ol, trifluorooa...	280	C14H23F3O2	1000352-26-2	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW020719\
Data File : VW008551.D
Acq On : 07 Feb 2019 14:35
Operator : SY/VA
Sample : K1390-01
Misc : 5.55G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
SB-01(12-13)

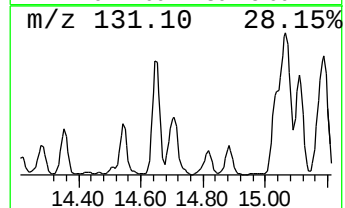
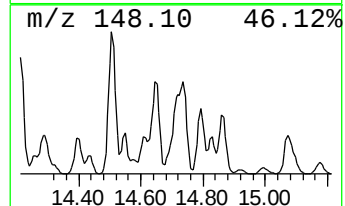
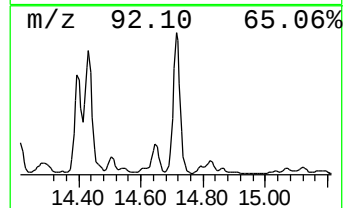
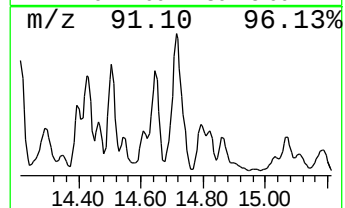
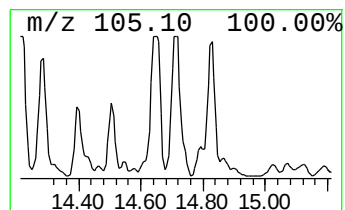
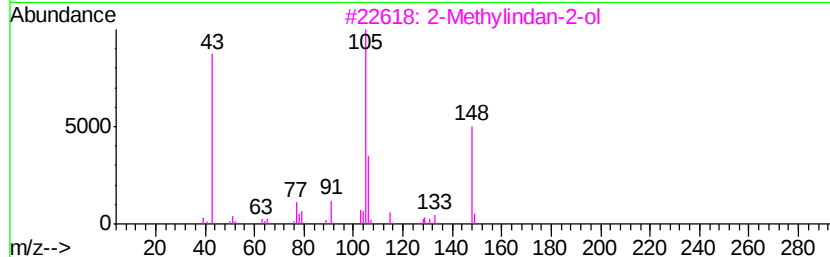
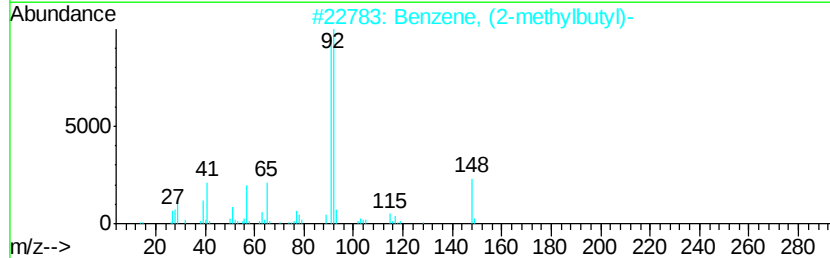
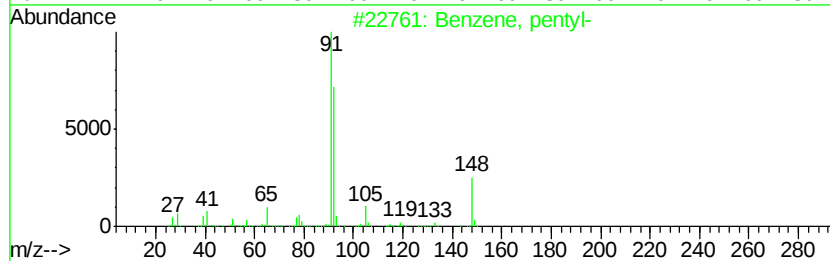
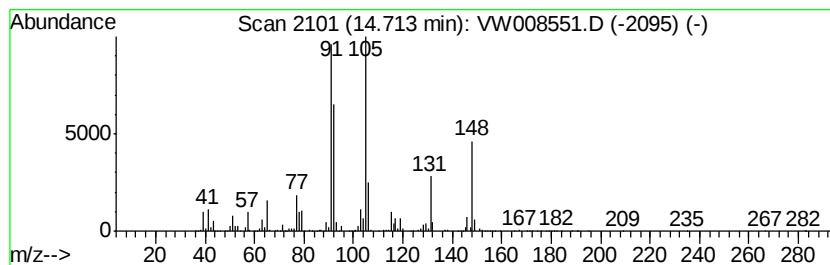
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W012419S.M
Quant Title : SW846 8260

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

Peak Number 10 unknown14.71 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.71	85.69 ug/l	191846000	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, pentvl-	148	C11H16	000538-68-1	49
2		Benzene, (2-methylbutyl)-	148	C11H16	003968-85-2	38
3		2-Methylindan-2-ol	148	C10H12O	033223-84-6	38
4		Benzene, (1-methylbutyl)-	148	C11H16	002719-52-0	35
5		Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW020719\
 Data File : VW008551.D
 Acq On : 07 Feb 2019 14:35
 Operator : SY/VA
 Sample : K1390-01
 Misc : 5.55G/5ML/MSVOA_W/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 SB-01(12-13)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W012419S.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
1-Pentene, 2,3-di...	8.45	189.0	ug/l	418041000	2	8.84	110593000	50.0
Aziridine, 2,2-di...	8.71	123.8	ug/l	273856000	2	8.84	110593000	50.0
Cyclopropane, 1-e...	9.63	207.5	ug/l	458895000	2	8.84	110593000	50.0
2,2-Dimethyl-1-ox...	9.78	255.3	ug/l	564652000	2	8.84	110593000	50.0
Pentanoic acid, 1...	9.92	169.8	ug/l	375581000	2	8.84	110593000	50.0
3-Hexene, (E)-	10.12	172.5	ug/l	381580000	2	8.84	110593000	50.0
unknown10.97	10.97	116.7	ug/l	456964000	3	11.63	195737000	50.0
Naphthalene, deca...	13.89	131.1	ug/l	293427000	4	13.55	111946000	50.0
unknown14.71	14.71	85.7	ug/l	191846000	4	13.55	111946000	50.0