

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW100419\
 Data File : VW013443.D
 Acq On : 04 Oct 2019 17:03
 Operator : SY/VA
 Sample : K5213-07
 Misc : 5.02G/5ML/MSVOA W/SOIL
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 S-4(0-5)

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\82W100119S.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	2.154	35	41	46	rBV2	9045	14853	1.25%	0.071%
2	2.355	65	74	81	rBV	9149	15960	1.35%	0.077%
3	2.635	113	120	125	rBV6	6961	14102	1.19%	0.068%
4	4.379	397	406	419	rBV3	7859	25624	2.16%	0.123%
5	5.007	500	509	520	rVB2	6559	22730	1.92%	0.109%
6	7.165	847	863	876	rBV2	169964	528938	44.62%	2.544%
7	7.878	974	980	985	rBV3	23659	56633	4.78%	0.272%
8	7.945	985	991	998	rVB	53761	120868	10.20%	0.581%
9	8.256	1034	1042	1047	rBV2	53108	134488	11.34%	0.647%
10	8.842	1131	1138	1147	rBV	70907	150343	12.68%	0.723%
11	9.092	1173	1179	1186	rBV	36510	70540	5.95%	0.339%
12	9.750	1283	1287	1292	rBV8	7427	15869	1.34%	0.076%
13	10.323	1375	1381	1390	rBV	121638	226451	19.10%	1.089%
14	10.707	1440	1444	1448	rVB	14737	24400	2.06%	0.117%
15	10.860	1464	1469	1475	rBV	61418	98282	8.29%	0.473%
16	10.933	1478	1481	1485	rBV4	18958	27793	2.34%	0.134%
17	11.225	1525	1529	1536	rVB	64729	122869	10.36%	0.591%
18	11.280	1536	1538	1542	rBV5	13259	14300	1.21%	0.069%
19	11.335	1544	1547	1551	rBV5	22044	31776	2.68%	0.153%
20	11.433	1560	1563	1571	rVB2	62407	110268	9.30%	0.530%
21	11.512	1573	1576	1584	rVB	31645	55060	4.64%	0.265%
22	11.634	1591	1596	1607	rVB	94768	189157	15.96%	0.910%
23	11.762	1614	1617	1620	rBV4	22952	27989	2.36%	0.135%
24	11.872	1631	1635	1639	rBV	48277	92508	7.80%	0.445%
25	11.975	1649	1652	1655	rBV4	15941	24570	2.07%	0.118%
26	12.134	1672	1678	1684	rBV3	122841	272397	22.98%	1.310%
27	12.250	1693	1697	1701	rVB	155379	225069	18.99%	1.083%
28	12.359	1702	1715	1722	rBV5	181793	674200	56.87%	3.243%
29	12.615	1751	1757	1763	rBV4	162715	333153	28.10%	1.602%
30	12.780	1779	1784	1789	rVB2	225570	384995	32.48%	1.852%
31	12.859	1790	1797	1803	rBV6	175655	535685	45.19%	2.577%
32	12.938	1804	1810	1824	rVB10	241212	1021409	86.16%	4.913%
33	13.079	1829	1833	1837	rVB3	155626	209398	17.66%	1.007%
34	13.127	1837	1841	1843	rBV3	92199	145979	12.31%	0.702%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW100419\
Data File : VW013443.D
Acq On : 04 Oct 2019 17:03
Operator : SY/VA
Sample : K5213-07
Misc : 5.02G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
S-4(0-5)

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\82W100119S.M
Title : SW846 8260

35	13.164	1843	1847	1860	rVB2	339111	731641	61.72%	3.519%
36	13.268	1860	1864	1865	rBV2	114646	170464	14.38%	0.820%
37	13.292	1865	1868	1871	rBV3	120953	153446	12.94%	0.738%
38	13.377	1880	1882	1885	rVB2	77030	80557	6.80%	0.387%
39	13.444	1886	1893	1897	rBV6	284810	680011	57.36%	3.271%
40	13.481	1897	1899	1904	rVB4	137532	179008	15.10%	0.861%
41	13.627	1919	1923	1928	rBV7	75970	166069	14.01%	0.799%
42	13.694	1929	1934	1944	rVV5	234702	658053	55.51%	3.165%
43	13.877	1958	1964	1968	rBV4	640204	1185441	100.00%	5.702%
44	13.981	1977	1981	1984	rBV4	213609	398519	33.62%	1.917%
45	14.091	1996	1999	2003	rVB	226672	304439	25.68%	1.464%
46	14.194	2012	2016	2022	rBV4	197965	411723	34.73%	1.980%
47	14.261	2023	2027	2033	rVB4	262142	461793	38.96%	2.221%
48	14.322	2033	2037	2041	rBV5	128079	269156	22.71%	1.295%
49	14.383	2042	2047	2051	rBV2	595200	969950	81.82%	4.665%
50	14.493	2062	2065	2068	rBV4	159724	250688	21.15%	1.206%
51	14.548	2069	2074	2079	rVV5	590746	988341	83.37%	4.754%
52	14.737	2102	2105	2109	rVB3	180422	239582	20.21%	1.152%
53	14.798	2110	2115	2119	rVV4	422700	915652	77.24%	4.404%
54	14.847	2119	2123	2130	rVB5	497414	1127179	95.09%	5.421%
55	14.975	2140	2144	2146	rBV3	154574	231342	19.52%	1.113%
56	15.060	2154	2158	2163	rVB6	289696	484877	40.90%	2.332%
57	15.176	2172	2177	2183	rVB4	244700	514351	43.39%	2.474%
58	15.328	2197	2202	2209	rVB2	393673	692613	58.43%	3.331%
59	15.420	2210	2217	2222	rBV5	333382	911390	76.88%	4.384%
60	15.737	2264	2269	2275	rVB7	237063	509678	42.99%	2.451%
61	15.865	2287	2290	2300	rVB6	198351	383961	32.39%	1.847%
62	16.371	2368	2373	2379	rVB3	177110	365624	30.84%	1.759%
63	16.633	2411	2416	2425	rVB8	127266	336821	28.41%	1.620%

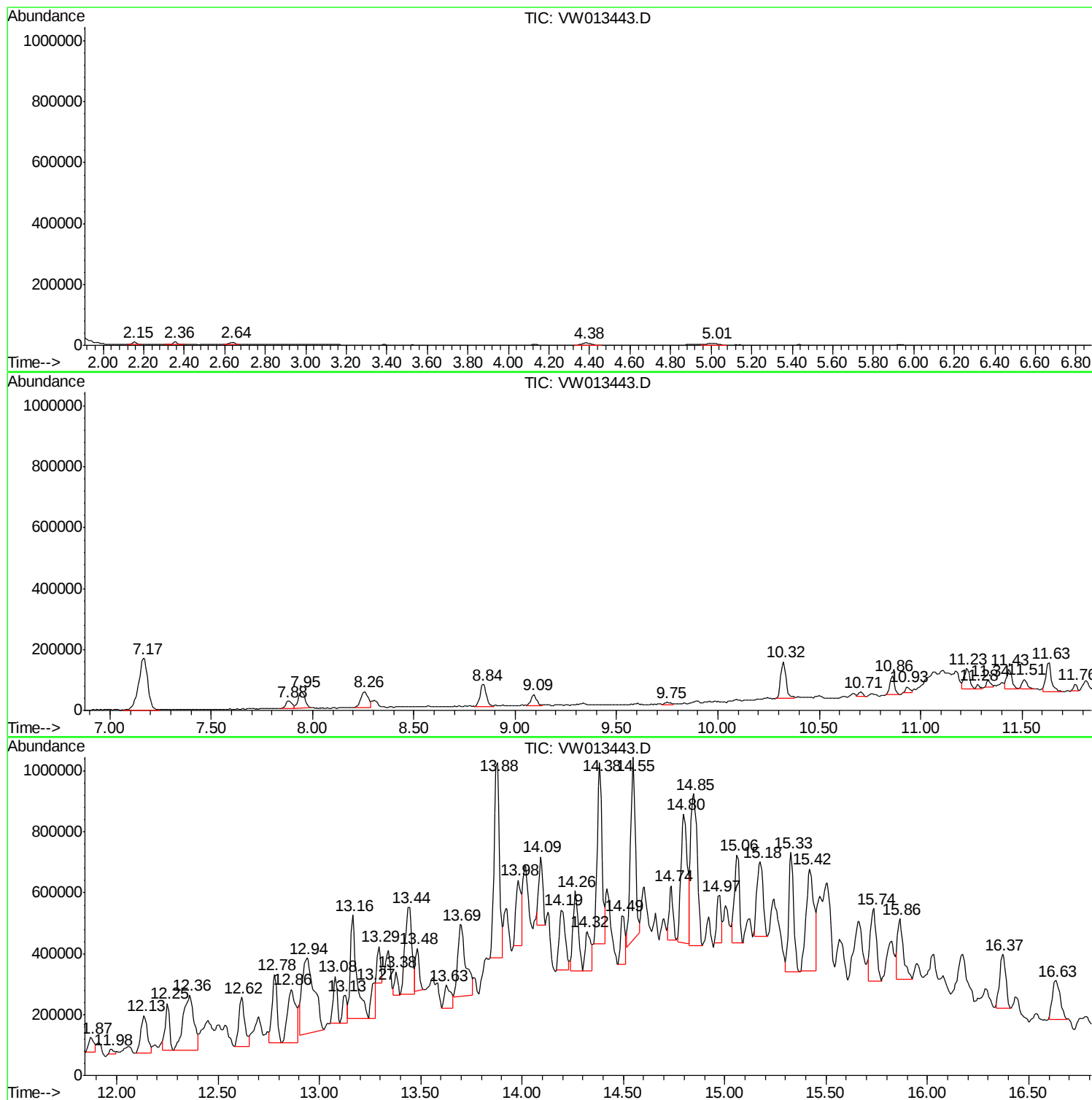
Sum of corrected areas: 20791025

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW100419\
Data File : VW013443.D
Acq On : 04 Oct 2019 17:03
Operator : SY/VA
Sample : K5213-07
Misc : 5.02G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
S-4(0-5)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W100119S.M
Quant Title : SW846 8260

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW100419\
Data File : VW013443.D
Acq On : 04 Oct 2019 17:03
Operator : SY/VA
Sample : K5213-07
Misc : 5.02G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
S-4(0-5)

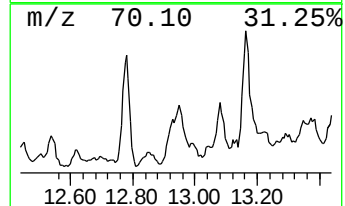
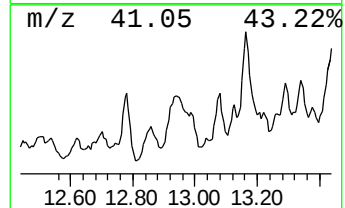
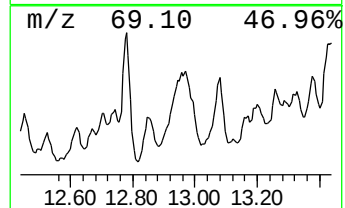
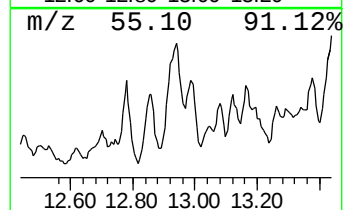
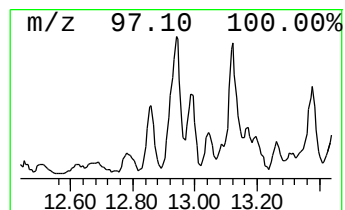
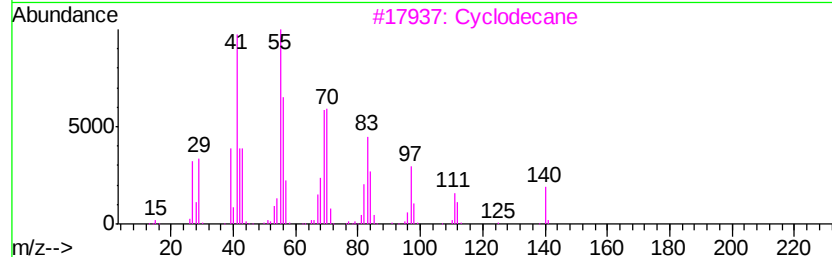
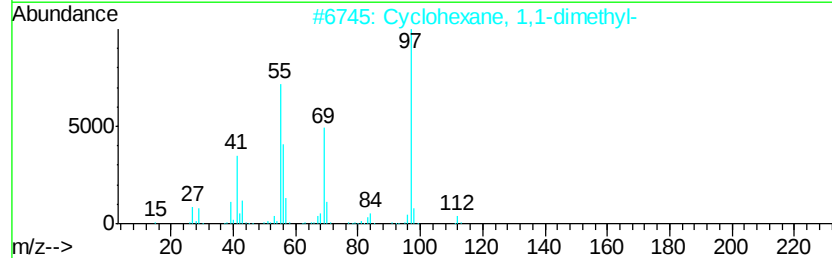
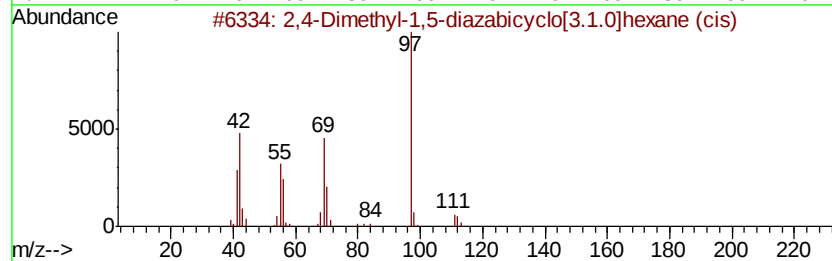
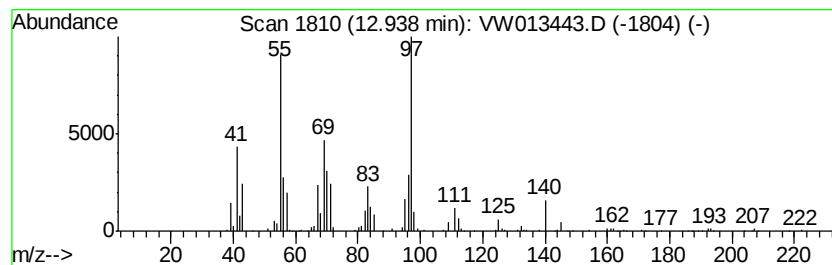
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W100119S.M
Quant Title : SW846 8260

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

Peak Number 1 unknown12.94 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.94	307.53 ug/l	1021410	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4-Dimethyl-1,5-diazabicyclo[3.1.0]hexane (cis)	112	C6H12N2	100463-01-2	46
2		Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	43
3		Cyclodecane	140	C10H20	000293-96-9	42
4		2,4-Dimethyl-1,5-diazabicyclo[3.1.0]hexane (cis)	112	C6H12N2	1000283-20-9	38
5		1H-Pyrazole, 4,5-dihydro-3,5,5-t...	112	C6H12N2	003975-85-7	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW100419\
Data File : VW013443.D
Acq On : 04 Oct 2019 17:03
Operator : SY/VA
Sample : K5213-07
Misc : 5.02G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

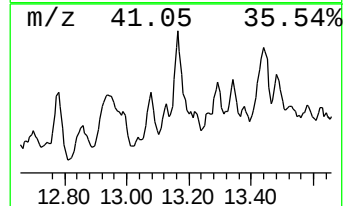
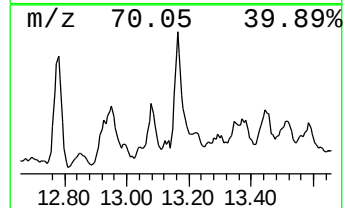
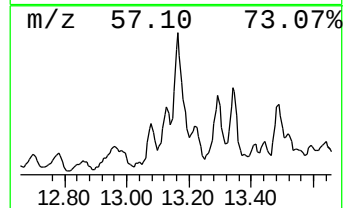
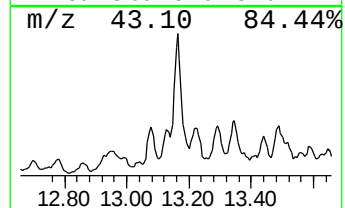
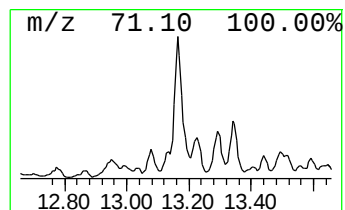
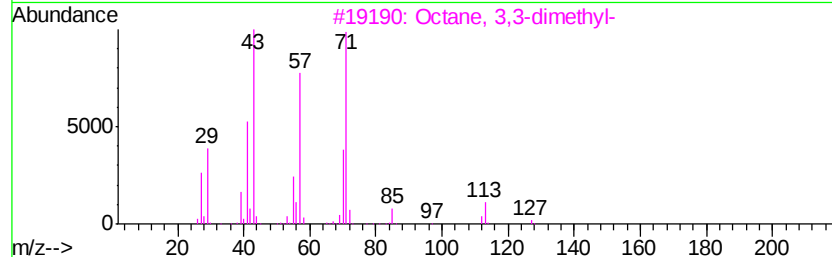
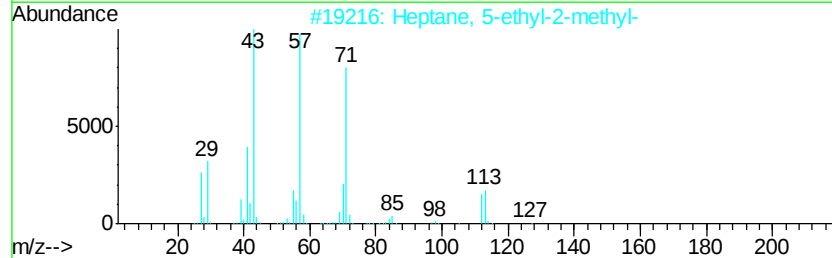
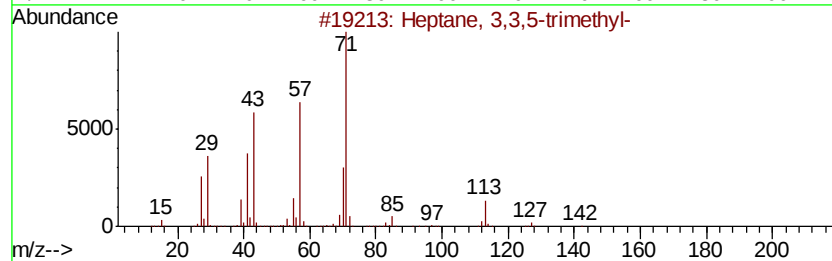
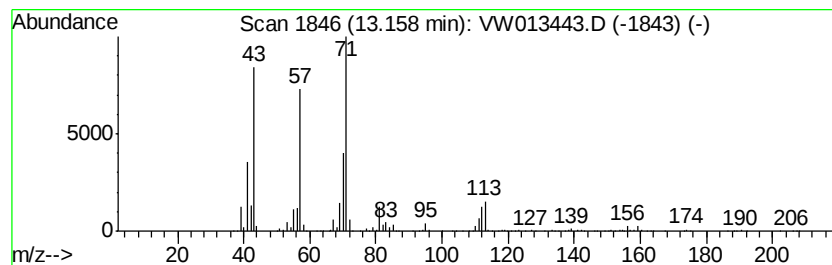
Instrument :
MSVOA_W
ClientSampled :
S-4(0-5)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W100119S.M
Quant Title : SW846 8260

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

Peak Number 2 Heptane, 3,3,5-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.16	220.28 ug/l	731641	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptane, 3,3,5-trimethyl-	142 C10H22	007154-80-5	72
2	Heptane, 5-ethyl-2-methyl-	142 C10H22	013475-78-0	59
3	Octane, 3,3-dimethyl-	142 C10H22	004110-44-5	53
4	Undecane, 4,5-dimethyl-	184 C13H28	017312-79-7	50
5	Oxalic acid, hexyl neopentyl ester	244 C13H24O4	1000309-73-1	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW100419\
Data File : VW013443.D
Acq On : 04 Oct 2019 17:03
Operator : SY/VA
Sample : K5213-07
Misc : 5.02G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
S-4(0-5)

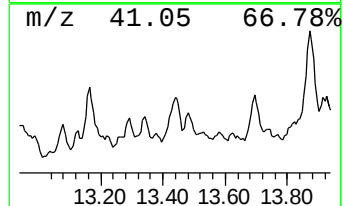
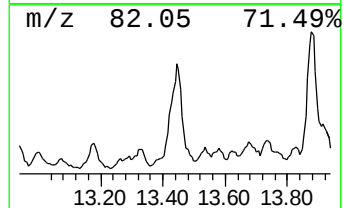
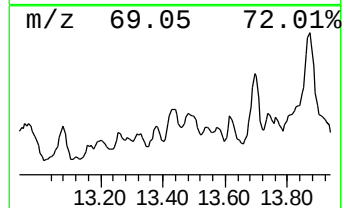
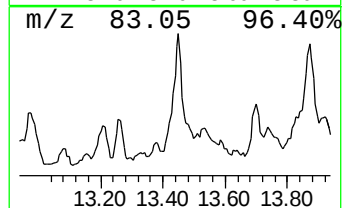
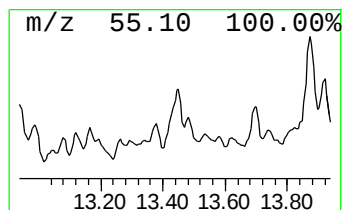
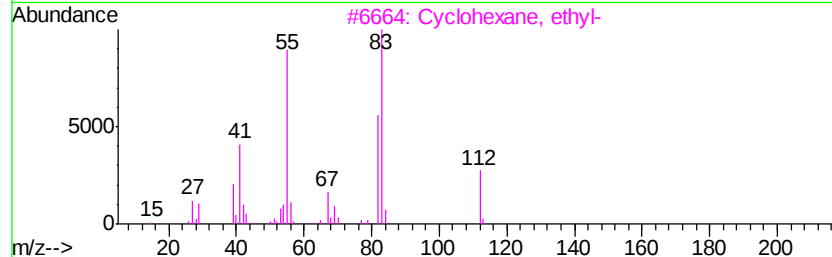
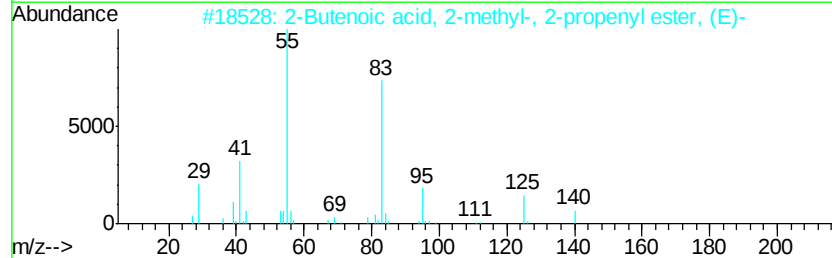
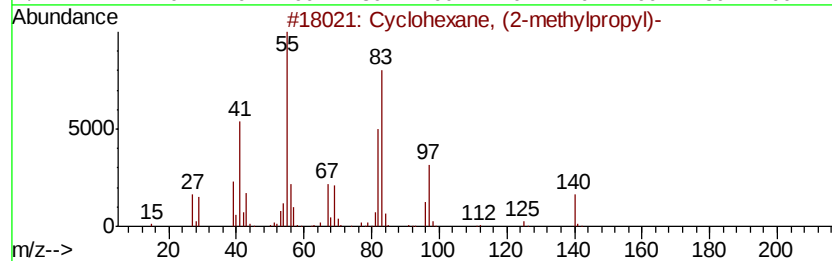
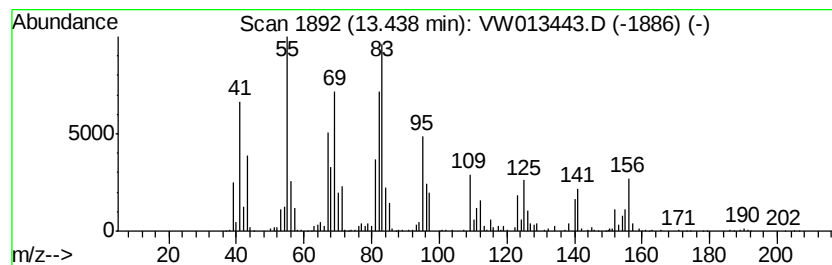
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W100119S.M
Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, (2-methylpropyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.44	204.74 ug/l	680011	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	53
2		2-Butenoic acid, 2-methyl-, 2-propenyl ester, (E)-	140	C8H12O2	007493-71-2	43
3		Cyclohexane, ethyl-	112	C8H16	001678-91-7	43
4		Cyclohexane, butyl-	140	C10H20	001678-93-9	43
5		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	41



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW100419\
Data File : VW013443.D
Acq On : 04 Oct 2019 17:03
Operator : SY/VA
Sample : K5213-07
Misc : 5.02G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
S-4(0-5)

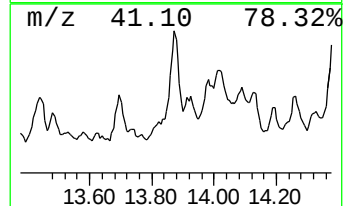
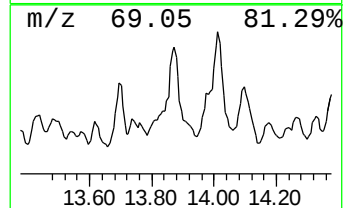
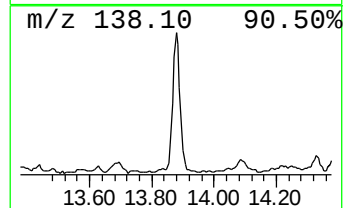
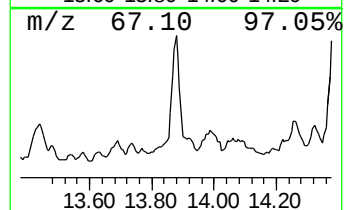
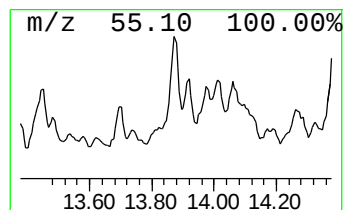
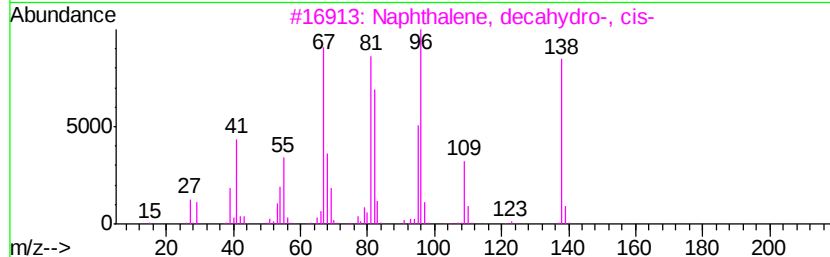
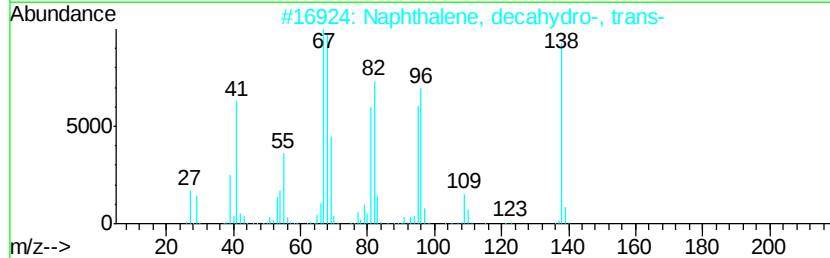
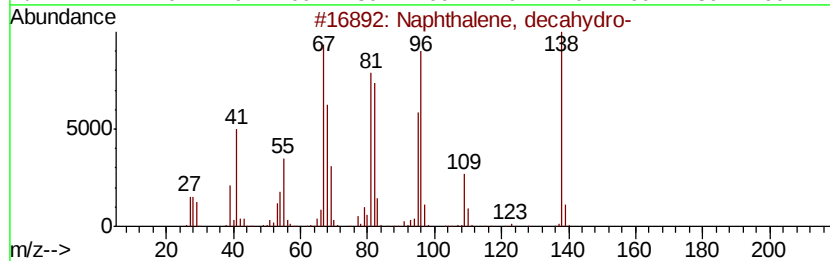
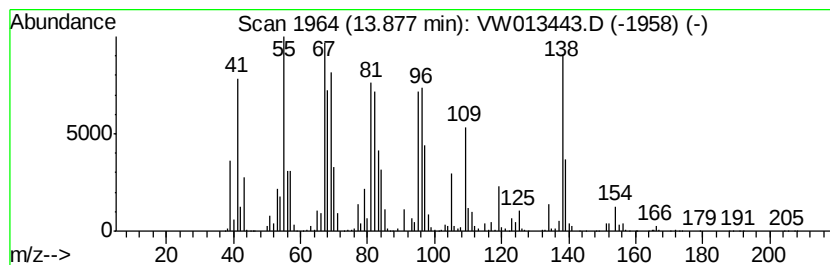
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W100119S.M
Quant Title : SW846 8260

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

Peak Number 4 Naphthalene, decahydro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.88	356.91 ug/l	1185440	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, decahydro-	138	C10H18	000091-17-8	93
2		Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	91
3		Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	64
4		Spiro[4.5]decane	138	C10H18	000176-63-6	60
5		Bicyclo[4.1.0]heptane, 3,7,7-tri...	138	C10H18	018968-23-5	55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW100419\
Data File : VW013443.D
Acq On : 04 Oct 2019 17:03
Operator : SY/VA
Sample : K5213-07
Misc : 5.02G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

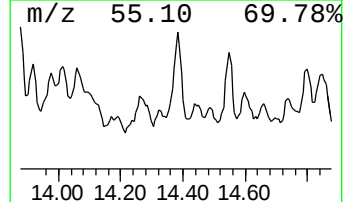
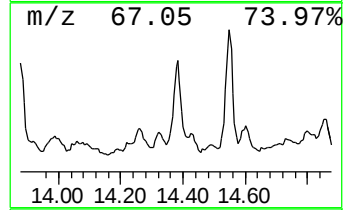
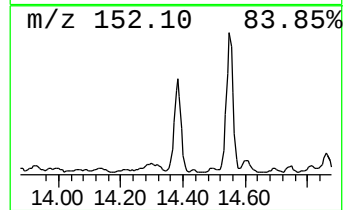
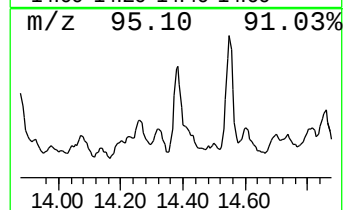
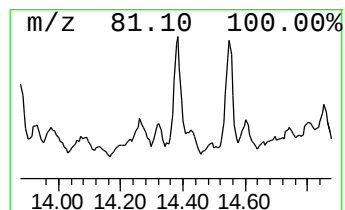
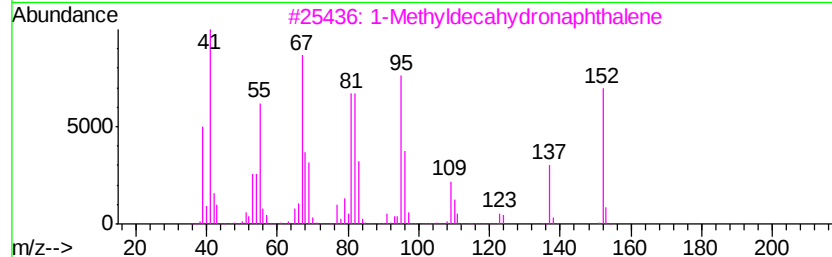
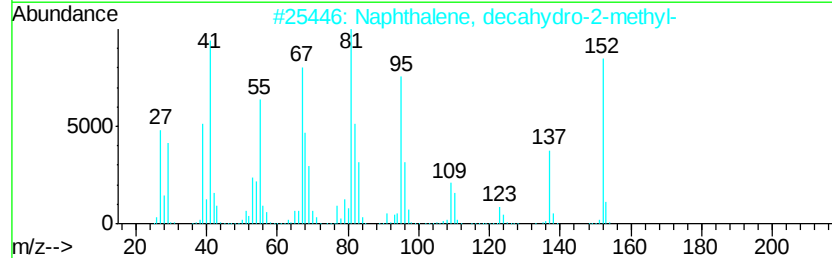
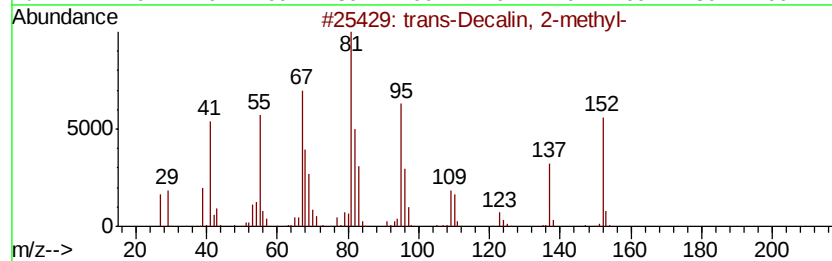
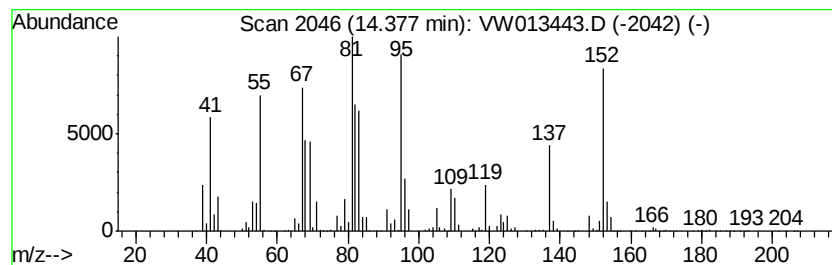
Instrument :
MSVOA_W
ClientSampleId :
S-4(0-5)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W100119S.M
Quant Title : SW846 8260

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

Peak Number 5 trans-Decalin, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.38	292.03 ug/l	969950	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	trans-Decalin, 2-methyl-	152 C11H20	1000152-47-3	86
2	Naphthalene, decahydro-2-methyl-	152 C11H20	002958-76-1	64
3	1-Methyldecahydronaphthalene	152 C11H20	002958-75-0	58
4	Cyclohexene, 1-pentyl-	152 C11H20	015232-85-6	50
5	6,6-Dimethyl-cyclooct-4-ene	152 C10H16O	1000194-10-0	49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW100419\
Data File : VW013443.D
Acq On : 04 Oct 2019 17:03
Operator : SY/VA
Sample : K5213-07
Misc : 5.02G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

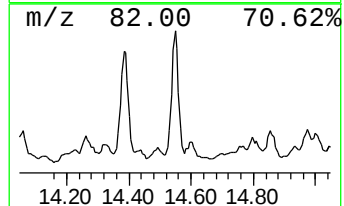
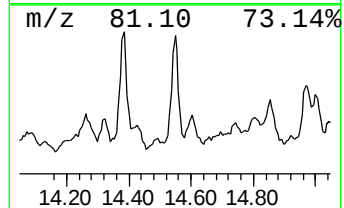
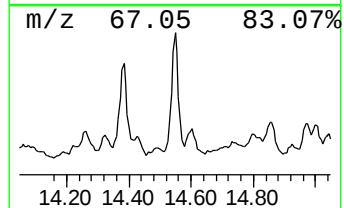
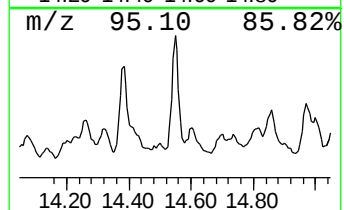
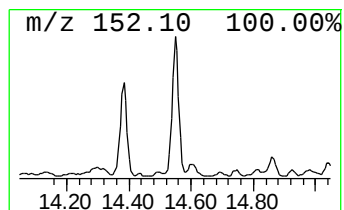
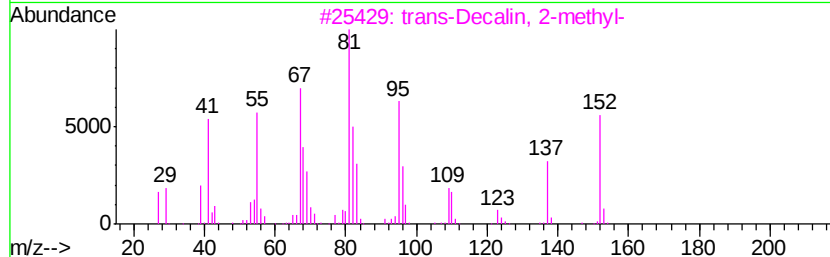
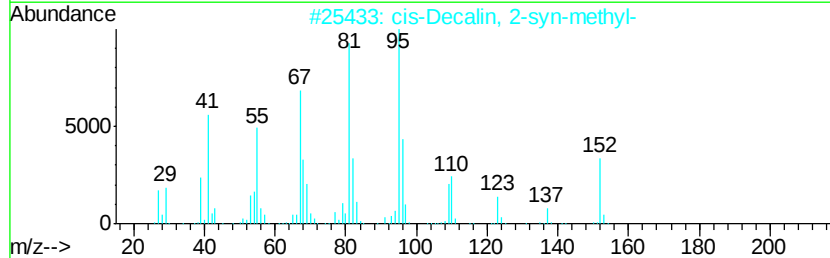
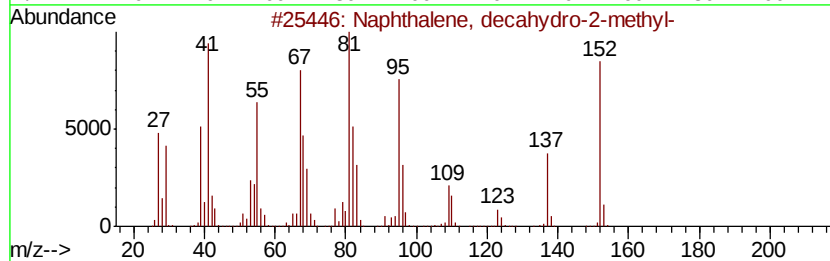
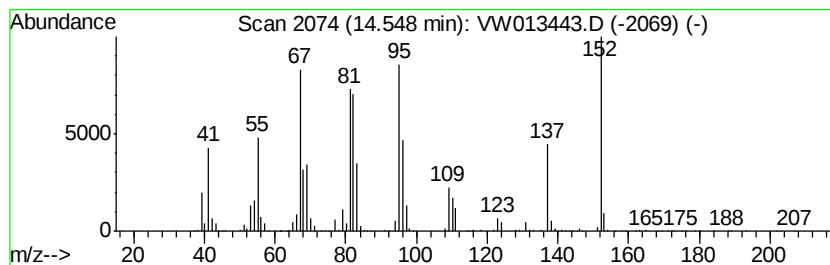
Instrument :
MSVOA_W
ClientSampleId :
S-4(0-5)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W100119S.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.
14.55	297.57 ug/l	988341	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, decahydro-2-methyl-	152 C11H20	002958-76-1 90
2		cis-Decalin, 2-syn-methyl-	152 C11H20	1000155-85-6 83
3		trans-Decalin, 2-methyl-	152 C11H20	1000152-47-3 64
4		Bicyclo[3.3.2]decan-9-one	152 C10H16O	028054-91-3 58
5		Bicyclo[5.4.0]undecane (trans)	152 C11H20	021394-30-9 52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW100419\
Data File : VW013443.D
Acq On : 04 Oct 2019 17:03
Operator : SY/VA
Sample : K5213-07
Misc : 5.02G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

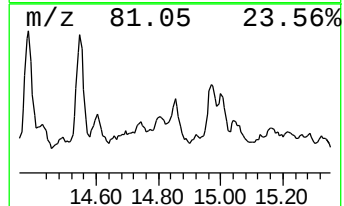
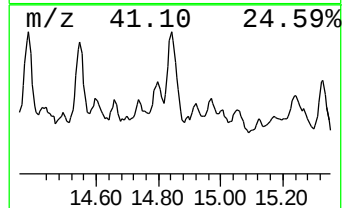
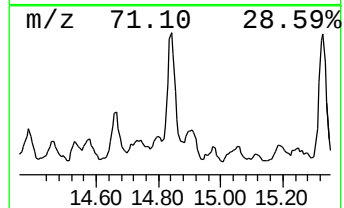
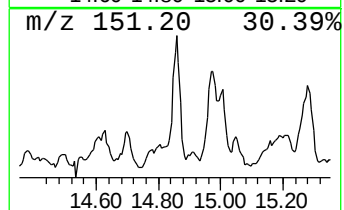
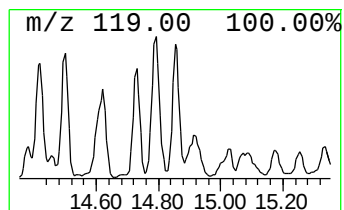
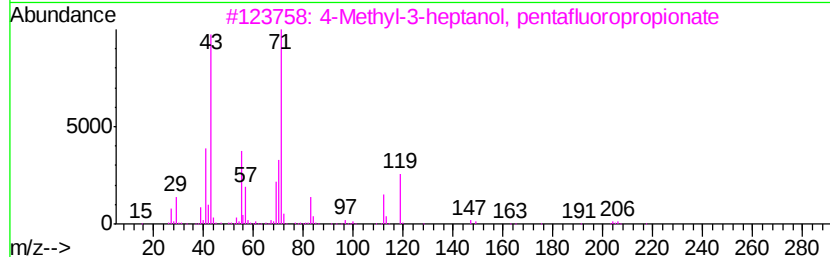
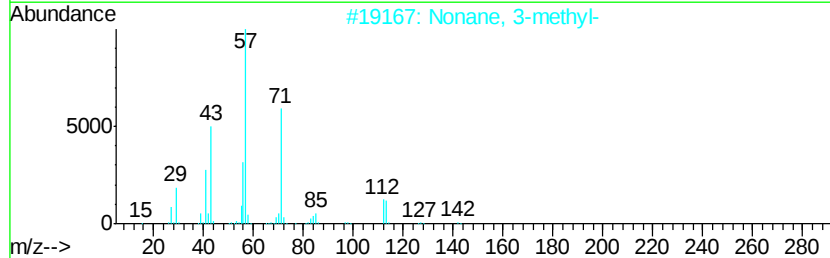
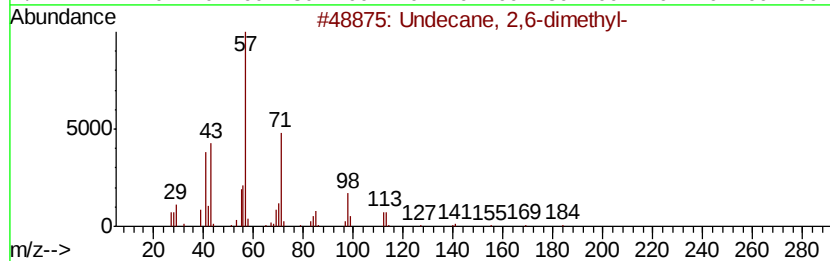
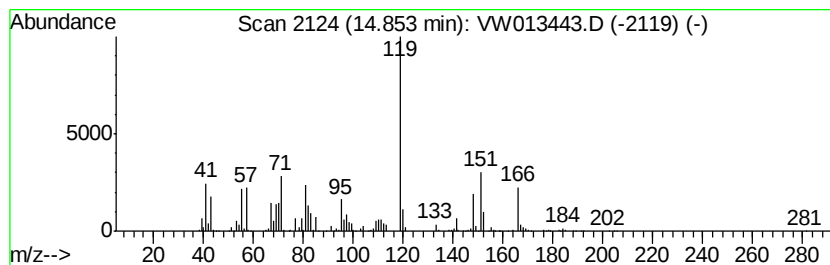
Instrument :
MSVOA_W
ClientSampleId :
S-4(0-5)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W100119S.M
Quant Title : SW846 8260

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

Peak Number 8 Undecane, 2,6-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	339.37 ug/l	1127180	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Undecane, 2,6-dimethyl-	184 C13H28	017301-23-4	38
2	Nonane, 3-methyl-	142 C10H22	005911-04-6	30
3	4-Methyl-3-heptanol, pentafluoro...	276 C11H17F5O2	1000365-51-7	27
4	4-Methyl-3-heptanol, heptafluoro...	326 C12H17F7O2	1000365-51-5	27
5	Oxalic acid, 2-ethylhexyl nonyl ...	328 C19H36O4	1000309-39-2	27



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW100419\
Data File : VW013443.D
Acq On : 04 Oct 2019 17:03
Operator : SY/VA
Sample : K5213-07
Misc : 5.02G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
S-4(0-5)

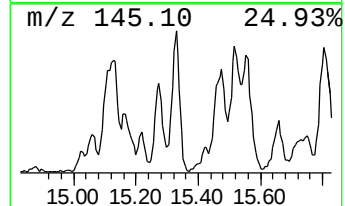
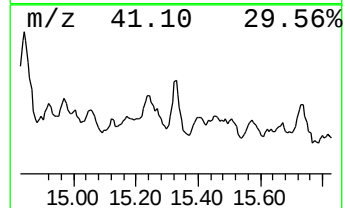
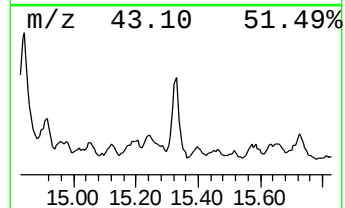
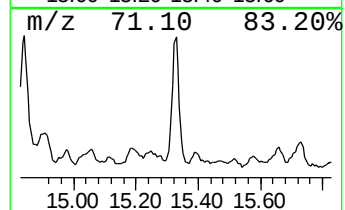
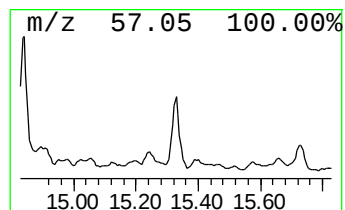
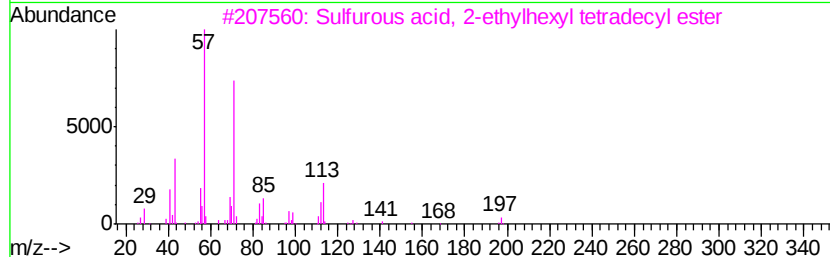
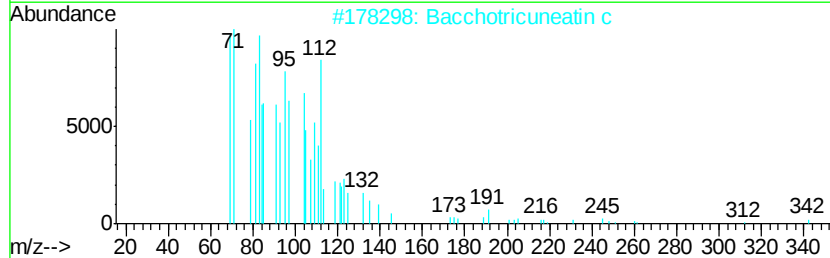
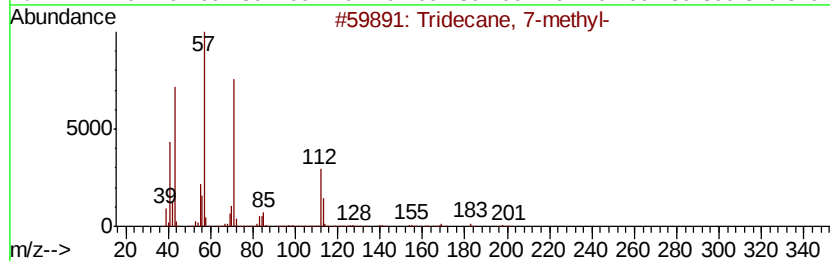
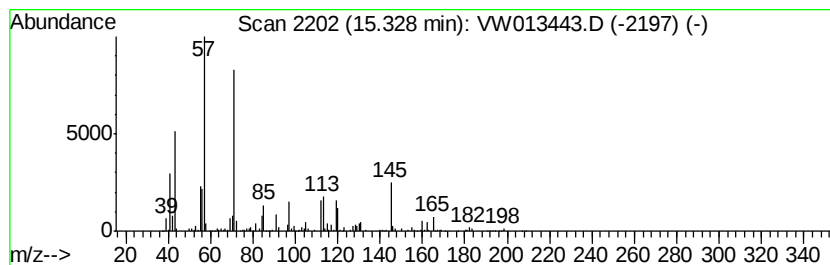
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W100119S.M
Quant Title : SW846 8260

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

Peak Number 9 Tridecane, 7-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.33	208.53 ug/l	692613	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 7-methyl-	198	C14H30	026730-14-3	70
2		Bacchotricuneatin c	342	C20H22O5	066563-30-2	68
3		Sulfurous acid, 2-ethylhexyl tet...	390	C22H46O3S	1000309-19-7	53
4		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	53
5		Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW100419\
Data File : VW013443.D
Acq On : 04 Oct 2019 17:03
Operator : SY/VA
Sample : K5213-07
Misc : 5.02G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
S-4(0-5)

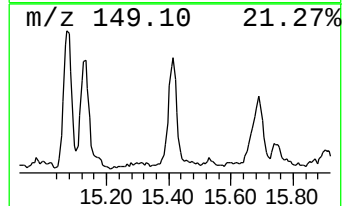
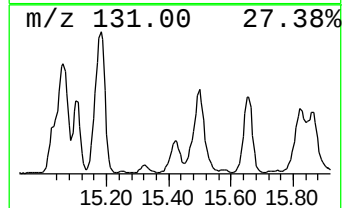
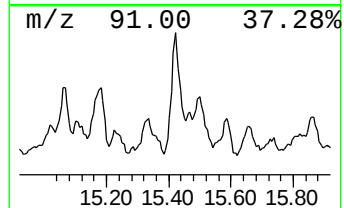
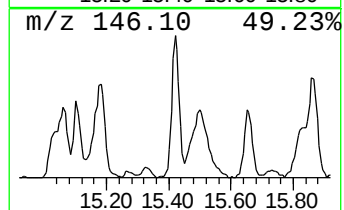
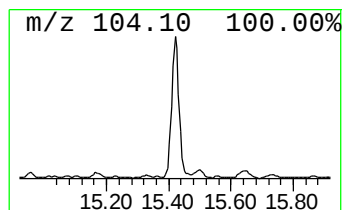
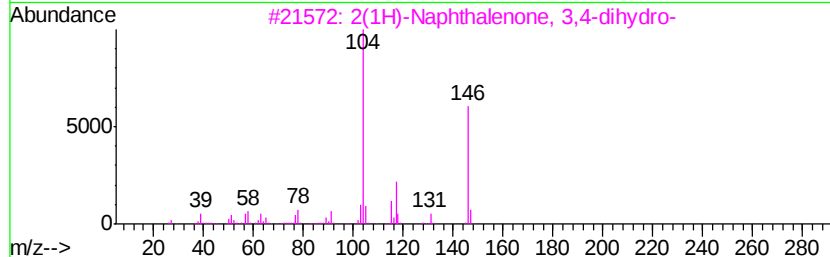
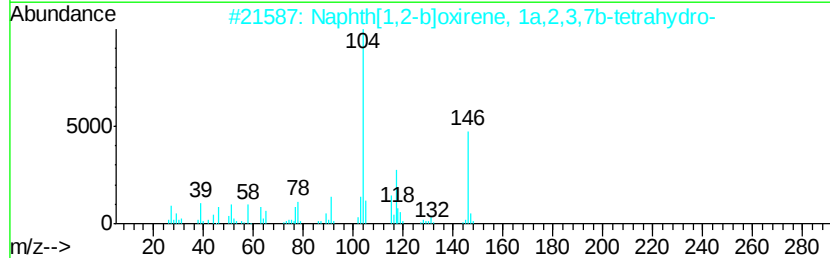
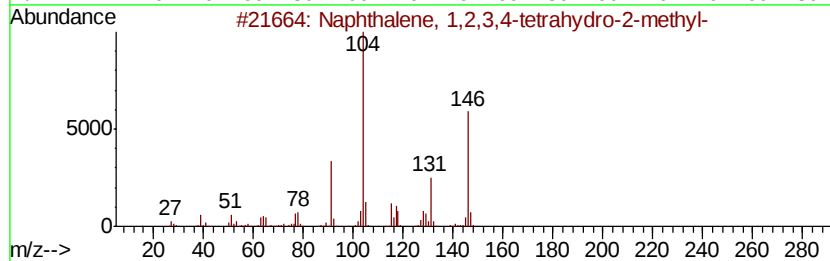
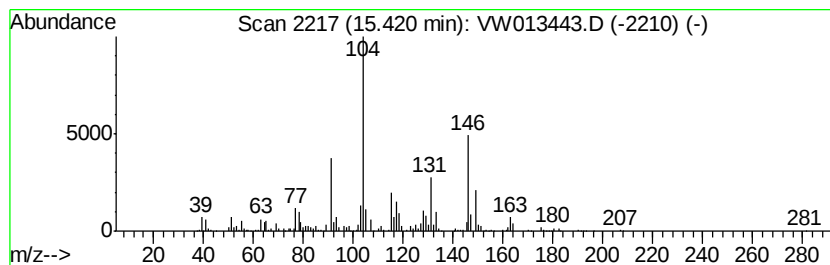
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W100119S.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.42	274.40 ug/l	911390	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	96
2		Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	58
3		2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	52
4		Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	49
5		Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW100419\
Data File : VW013443.D
Acq On : 04 Oct 2019 17:03
Operator : SY/VA
Sample : K5213-07
Misc : 5.02G/5ML/MSVOA_W/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
S-4(0-5)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W100119S.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
unknown12.94	12.94	307.5	ug/l	1021410	4	13.56	166069	50.0
Heptane, 3,3,5-tr...	13.16	220.3	ug/l	731641	4	13.56	166069	50.0
Cyclohexane, (2-m...	13.44	204.7	ug/l	680011	4	13.56	166069	50.0
Naphthalene, deca...	13.88	356.9	ug/l	1185440	4	13.56	166069	50.0
trans-Decalin, 2-...	14.38	292.0	ug/l	969950	4	13.56	166069	50.0
Naphthalene, deca...	14.55	297.6	ug/l	988341	4	13.56	166069	50.0
Undecane, 2,6-dim...	14.85	339.4	ug/l	1127180	4	13.56	166069	50.0
Tridecane, 7-methyl-	15.33	208.5	ug/l	692613	4	13.56	166069	50.0
Naphthalene, 1,2,...	15.42	274.4	ug/l	911390	4	13.56	166069	50.0