

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD111920\
 Data File : VD067703.D
 Acq On : 19 Nov 2020 12:54
 Operator : VA/SY
 Sample : L4778-06
 Misc : 4.07G/5.00ml/MSVOA D/SOIL
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 VNJ-223

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_D\METHOD\82D110320S.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	2.203	47	48	52	rVB2	1573	1567	1.41%	0.155%
2	2.627	118	120	128	rVB2	1770	3599	3.25%	0.356%
3	2.703	131	133	137	rVB3	1513	1556	1.40%	0.154%
4	2.791	142	148	150	rVV4	1313	2245	2.02%	0.222%
5	4.968	513	518	524	rVB5	1647	3438	3.10%	0.340%
6	6.885	839	844	845	rVV	910	1271	1.15%	0.126%
7	7.262	905	908	915	rVB2	616	1420	1.28%	0.140%
8	7.920	1011	1020	1025	rBV3	17010	43331	39.07%	4.283%
9	7.979	1025	1030	1044	rVB4	34439	74136	66.85%	7.327%
10	8.291	1078	1083	1084	rBV	3177	4401	3.97%	0.435%
11	8.332	1084	1090	1099	rVB4	19179	44452	40.08%	4.393%
12	8.867	1172	1181	1189	rBV3	41394	82616	74.49%	8.165%
13	10.344	1426	1432	1440	rBV	59612	106400	95.94%	10.516%
14	10.726	1490	1497	1499	rBV3	1465	2513	2.27%	0.248%
15	10.891	1520	1525	1529	rBV3	1495	2197	1.98%	0.217%
16	11.408	1609	1613	1616	rBV2	797	1409	1.27%	0.139%
17	11.650	1647	1654	1661	rVB	65787	110906	100.00%	10.961%
18	11.755	1667	1672	1677	rVB3	859	1372	1.24%	0.136%
19	11.861	1685	1690	1694	rBV3	1680	3261	2.94%	0.322%
20	11.979	1709	1710	1717	rVB2	1047	1663	1.50%	0.164%
21	12.191	1745	1746	1751	rVB2	1296	1733	1.56%	0.171%
22	12.497	1791	1798	1803	rVV6	7473	14433	13.01%	1.426%
23	12.638	1813	1822	1828	rVV3	47480	83594	75.37%	8.262%
24	12.826	1848	1854	1857	rBV3	1078	1911	1.72%	0.189%
25	12.908	1864	1868	1872	rBV2	5068	5894	5.31%	0.583%
26	12.955	1874	1876	1880	rVV4	5523	6602	5.95%	0.653%
27	13.008	1880	1885	1888	rVV2	1263	2311	2.08%	0.228%
28	13.044	1888	1891	1893	rVB2	1798	1416	1.28%	0.140%
29	13.150	1906	1909	1910	rVV	2445	2682	2.42%	0.265%
30	13.185	1913	1915	1919	rVV3	2486	2923	2.64%	0.289%
31	13.232	1919	1923	1925	rVV3	1734	2376	2.14%	0.235%
32	13.332	1935	1940	1942	rBV2	4072	5611	5.06%	0.555%
33	13.402	1948	1952	1955	rBV5	2049	3295	2.97%	0.326%
34	13.444	1957	1959	1962	rVV2	3329	3807	3.43%	0.376%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD111920\
 Data File : VD067703.D
 Acq On : 19 Nov 2020 12:54
 Operator : VA/SY
 Sample : L4778-06
 Misc : 4.07G/5.00ml/MSVOA D/SOIL
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 VNJ-223

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_D\METHOD\82D110320S.M
 Title : SW846 8260

35	13.467	1962	1963	1966	rVV3	3467	3563	3.21%	0.352%
36	13.491	1966	1967	1968	rVV	3939	2051	1.85%	0.203%
37	13.508	1968	1970	1972	rVV2	4698	4959	4.47%	0.490%
38	13.526	1972	1973	1975	rVV2	3357	3035	2.74%	0.300%
39	13.579	1976	1982	1987	rVV3	57304	91997	82.95%	9.092%
40	13.638	1991	1992	1998	rVB4	4550	6505	5.87%	0.643%
41	13.779	2011	2016	2017	rBV2	2125	2614	2.36%	0.258%
42	13.820	2017	2023	2025	rBV5	3457	5019	4.53%	0.496%
43	13.867	2027	2031	2032	rVV3	5135	6979	6.29%	0.690%
44	13.897	2032	2036	2041	rVV4	15870	22715	20.48%	2.245%
45	13.944	2041	2044	2047	rVV3	3137	4008	3.61%	0.396%
46	13.997	2051	2053	2054	rVV2	2470	2339	2.11%	0.231%
47	14.049	2057	2062	2068	rBV6	2858	5810	5.24%	0.574%
48	14.114	2068	2073	2079	rBV8	5594	10333	9.32%	1.021%
49	14.232	2088	2093	2098	rBV7	4246	9638	8.69%	0.953%
50	14.285	2098	2102	2105	rVV6	3902	6849	6.18%	0.677%
51	14.320	2105	2108	2110	rVV4	2351	2703	2.44%	0.267%
52	14.344	2110	2112	2114	rVV2	3197	3033	2.73%	0.300%
53	14.391	2114	2120	2121	rVV3	5210	10164	9.16%	1.005%
54	14.420	2121	2125	2128	rVV4	6913	12852	11.59%	1.270%
55	14.455	2128	2131	2135	rVB5	4693	6554	5.91%	0.648%
56	14.532	2140	2144	2146	rVB5	4313	5579	5.03%	0.551%
57	14.567	2148	2150	2156	rVB4	4009	4918	4.43%	0.486%
58	14.614	2156	2158	2159	rBV2	2964	1698	1.53%	0.168%
59	14.632	2159	2161	2164	rVB2	4732	5314	4.79%	0.525%
60	14.655	2164	2165	2166	rBV	2646	1560	1.41%	0.154%
61	14.673	2166	2168	2169	rBV2	2146	1533	1.38%	0.152%
62	14.726	2172	2177	2179	rBV3	11288	15266	13.76%	1.509%
63	14.749	2179	2181	2187	rVB5	12515	15639	14.10%	1.546%
64	14.826	2187	2194	2198	rBV7	8221	16441	14.82%	1.625%
65	14.867	2199	2201	2202	rBV2	3515	2448	2.21%	0.242%
66	14.973	2218	2219	2220	rBV	3511	2066	1.86%	0.204%
67	15.038	2227	2230	2233	rBV4	3948	5180	4.67%	0.512%
68	15.161	2249	2251	2253	rBV3	3188	3572	3.22%	0.353%
69	15.202	2256	2258	2260	rVV3	4228	3902	3.52%	0.386%
70	15.249	2263	2266	2267	rVV3	2960	2959	2.67%	0.292%
71	15.285	2271	2272	2274	rVV2	3409	2699	2.43%	0.267%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD111920\
 Data File : VD067703.D
 Acq On : 19 Nov 2020 12:54
 Operator : VA/SY
 Sample : L4778-06
 Misc : 4.07G/5.00ml/MSVOA D/SOIL
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 VNJ-223

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_D\METHOD\82D110320S.M
 Title : SW846 8260

72	15.302	2274	2275	2277	rVV2	3217	1462	1.32%	0.144%
73	15.320	2277	2278	2281	rVB3	3231	1520	1.37%	0.150%
74	15.355	2281	2284	2285	rBV3	3725	3185	2.87%	0.315%
75	15.373	2285	2287	2290	rVB4	2188	1519	1.37%	0.150%
76	15.461	2298	2302	2303	rBV4	2737	2758	2.49%	0.273%
77	15.514	2308	2311	2313	rVB3	2445	2789	2.51%	0.276%
78	15.532	2313	2314	2316	rBV2	1786	1636	1.48%	0.162%
79	15.602	2323	2326	2330	rVB5	4165	4909	4.43%	0.485%
80	15.679	2338	2339	2341	rBV2	2095	1383	1.25%	0.137%
81	15.708	2342	2344	2346	rVV3	3065	2905	2.62%	0.287%
82	15.726	2346	2347	2350	rVV3	3279	2026	1.83%	0.200%
83	15.755	2350	2352	2354	rVV2	2112	2188	1.97%	0.216%
84	15.779	2354	2356	2363	rVB6	3014	3949	3.56%	0.390%
85	15.896	2372	2376	2377	rBV4	1632	2120	1.91%	0.210%
86	15.926	2377	2381	2382	rVV3	2629	2956	2.67%	0.292%
87	15.943	2382	2384	2386	rVB2	2870	1934	1.74%	0.191%
88	15.991	2390	2392	2395	rVB3	1503	1263	1.14%	0.125%
89	16.132	2414	2416	2418	rVB2	1963	1730	1.56%	0.171%
90	16.149	2418	2419	2425	rVB4	1382	1994	1.80%	0.197%
91	16.220	2429	2431	2433	rVV3	3021	2657	2.40%	0.263%
92	16.238	2433	2434	2437	rVB3	2265	2009	1.81%	0.199%
93	16.314	2445	2447	2449	rVB3	2551	1866	1.68%	0.184%
94	16.420	2462	2465	2468	rVB4	2189	2016	1.82%	0.199%
95	16.479	2471	2475	2476	rBV2	2398	2642	2.38%	0.261%
96	16.508	2479	2480	2482	rBV2	1528	1510	1.36%	0.149%
97	16.579	2489	2492	2497	rVB3	2136	3495	3.15%	0.345%
98	16.620	2497	2499	2501	rBV2	2273	1823	1.64%	0.180%
99	16.691	2509	2511	2513	rVV2	2012	1267	1.14%	0.125%
100	16.773	2521	2525	2527	rVB3	1492	1454	1.31%	0.144%

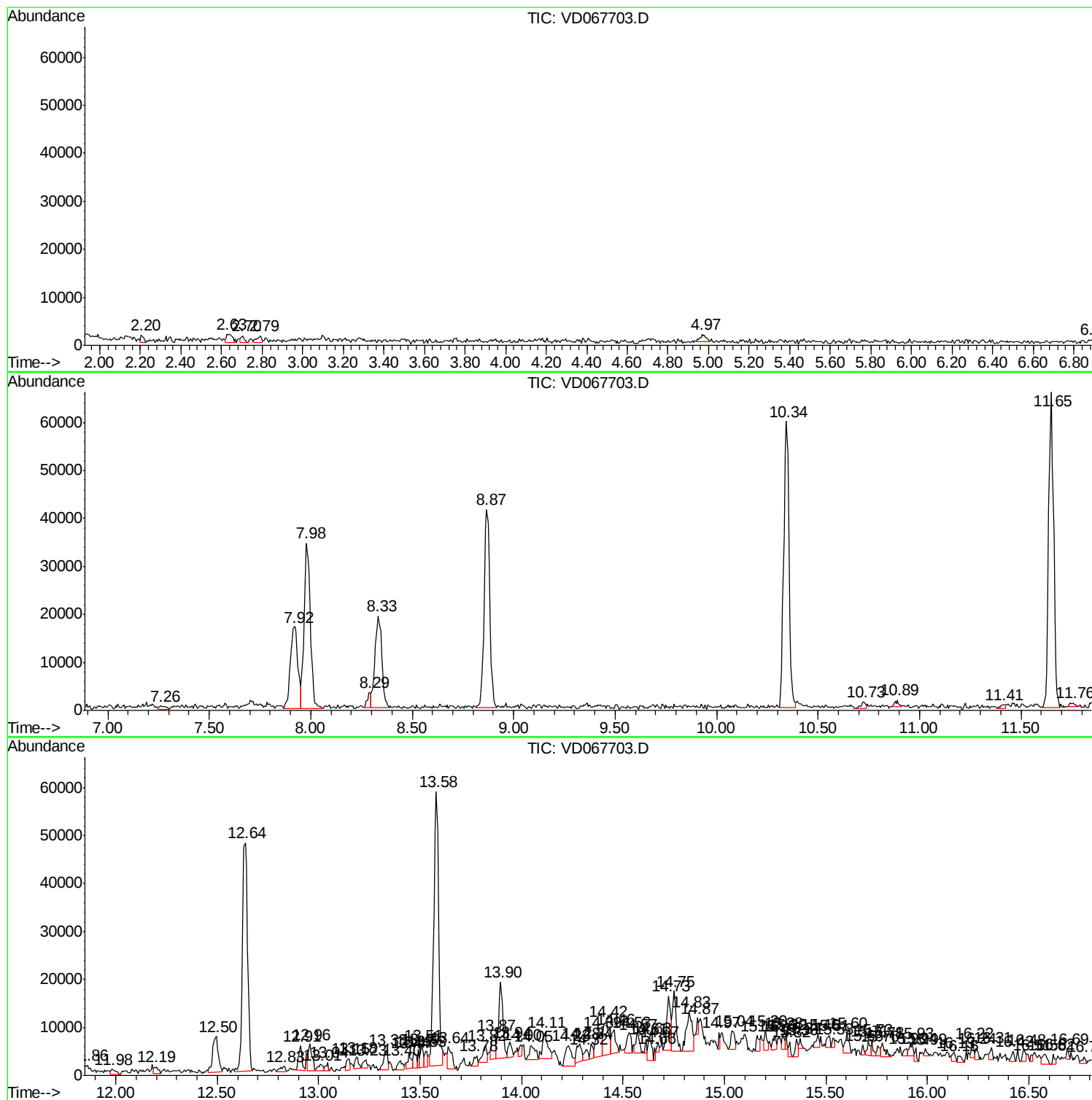
Sum of corrected areas: 1011800

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111920\
Data File : VD067703.D
Acq On : 19 Nov 2020 12:54
Operator : VA/SY
Sample : L4778-06
Misc : 4.07G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
VNJ-223

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.M
Quant Title : SW846 8260

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111920\
Data File : VD067703.D
Acq On : 19 Nov 2020 12:54
Operator : VA/SY
Sample : L4778-06
Misc : 4.07G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

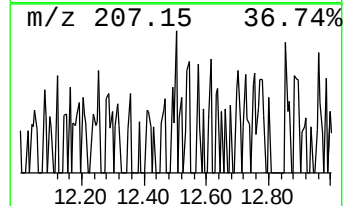
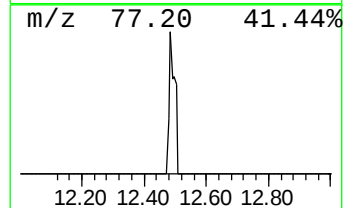
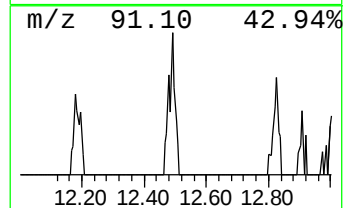
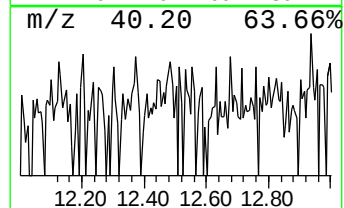
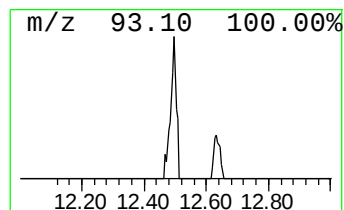
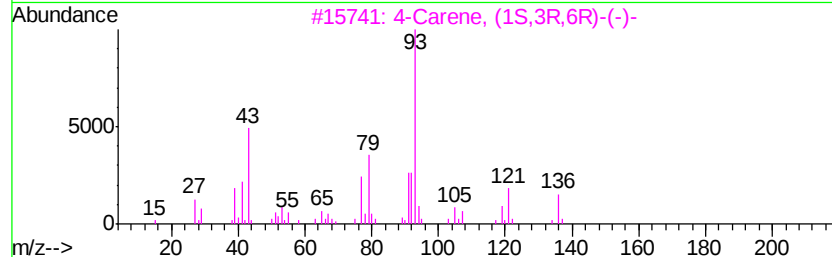
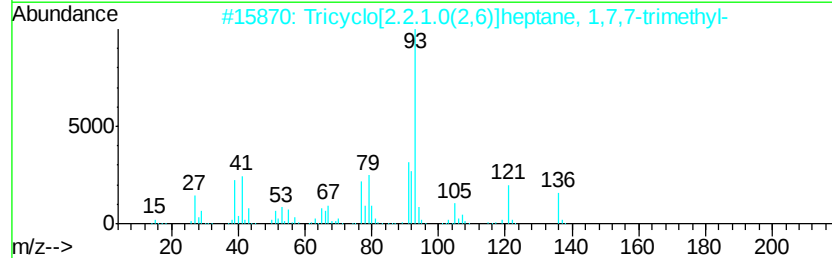
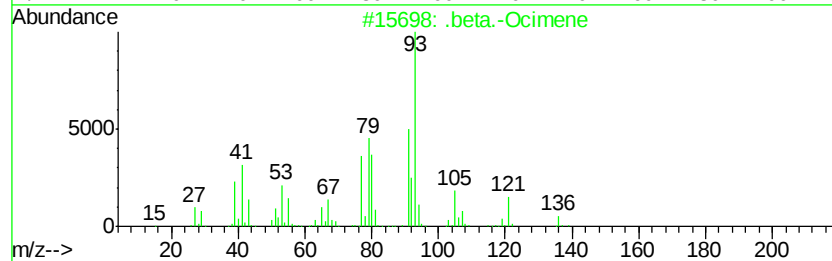
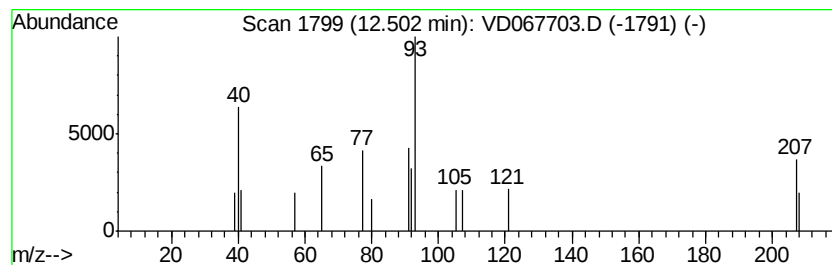
Instrument :
MSVOA_D
ClientSampleId :
VNJ-223

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.M
Quant Title : SW846 8260

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

Peak Number 1 .beta.-Ocimene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.50	6.51 ug/l	14433	Chlorobenzene-d5	11.65	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-Ocimene	136	C10H16	013877-91-3	86
2	Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	80
3	4-Carene, (1S,3R,6R)-(-)-	136	C10H16	005208-49-1	80
4	3-Carene	136	C10H16	013466-78-9	80
5	Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000488-97-1	80



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111920\
Data File : VD067703.D
Acq On : 19 Nov 2020 12:54
Operator : VA/SY
Sample : L4778-06
Misc : 4.07G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

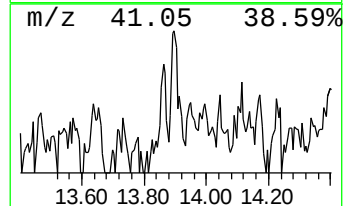
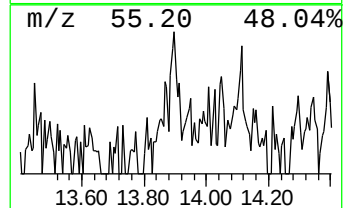
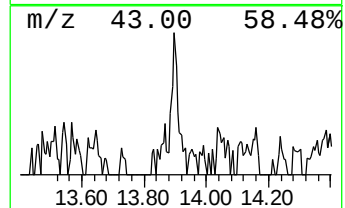
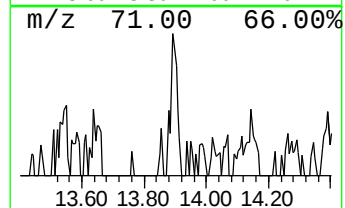
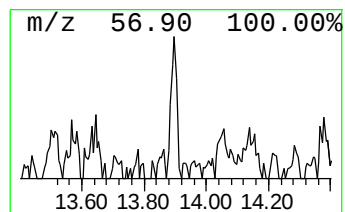
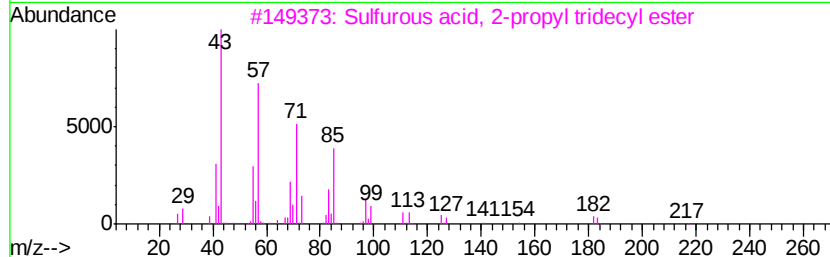
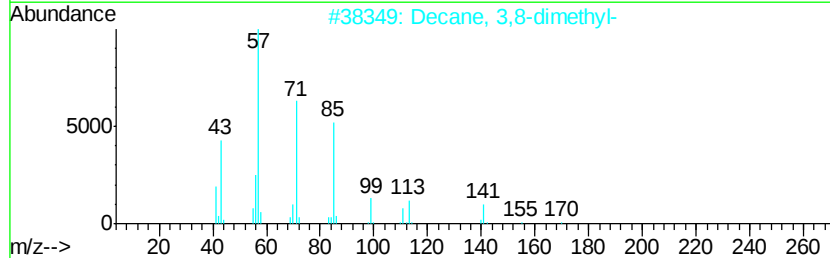
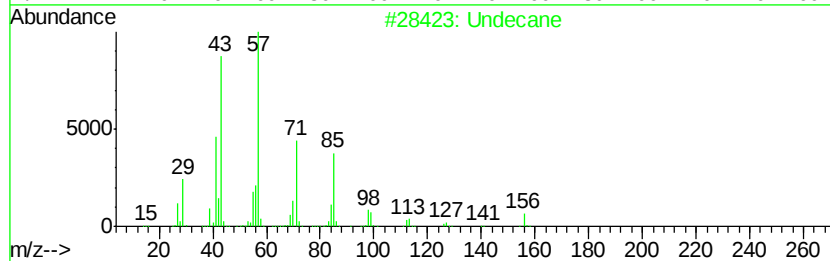
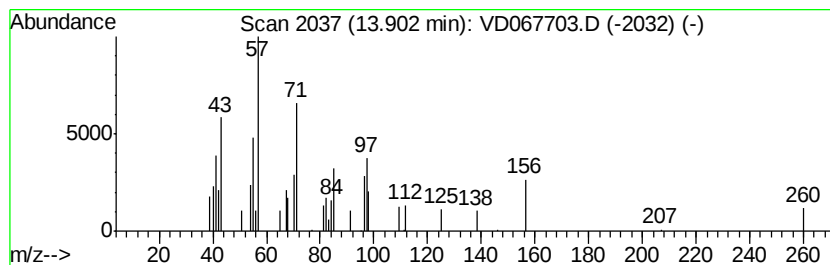
Instrument :
MSVOA_D
ClientSampleId :
VNJ-223

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.M
Quant Title : SW846 8260

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

Peak Number 2 unknown13.90 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	12.35 ug/l	22715	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Undecane		156 C11H24	001120-21-4 46
2	Decane, 3,8-dimethyl-		170 C12H26	017312-55-9 43
3	Sulfurous acid, 2-propyl tridecyl...		306 C16H34O3S	1000309-12-4 43
4	Heptacosane		380 C27H56	000593-49-7 43
5	Eicosane, 10-methyl-		296 C21H44	054833-23-7 43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111920\
Data File : VD067703.D
Acq On : 19 Nov 2020 12:54
Operator : VA/SY
Sample : L4778-06
Misc : 4.07G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

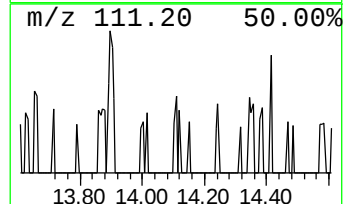
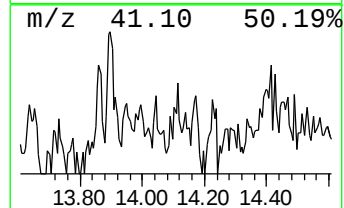
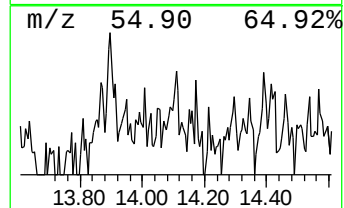
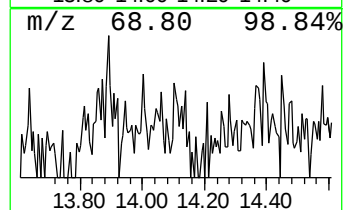
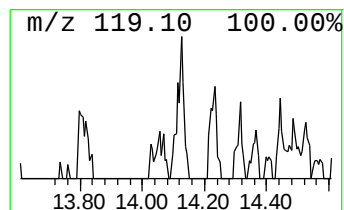
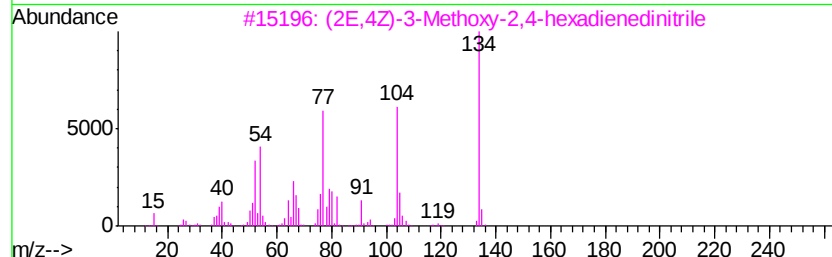
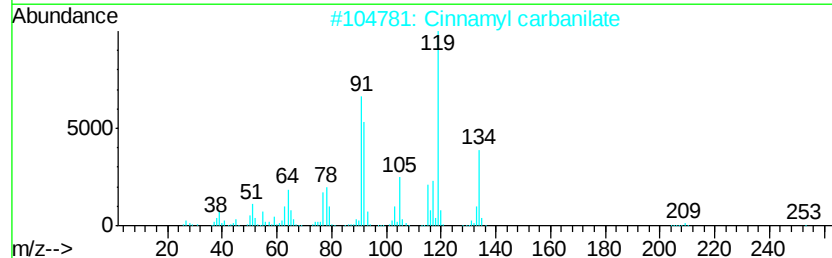
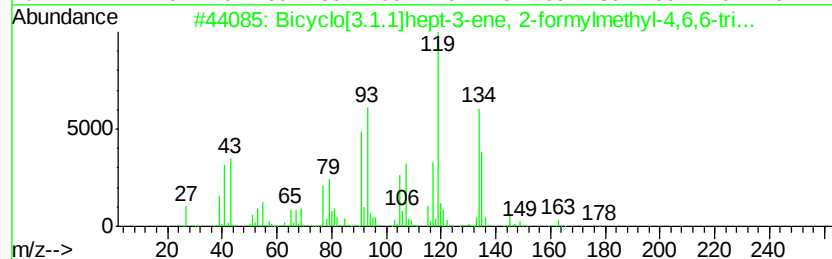
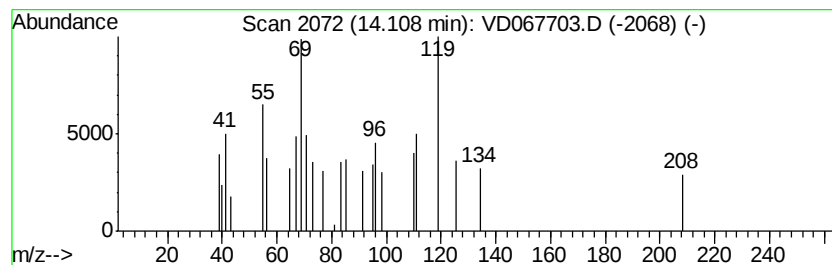
Instrument :
MSVOA_D
ClientSampleId :
VNJ-223

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.M
Quant Title : SW846 8260

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

Peak Number 3 unknown14.11 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.11	5.62 ug/l	10333	1,4-Dichlorobenzene-d4	13.58	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicvclo[3.1.1]hept-3-ene, 2-form...	178	C12H18O	135004-95-4	28
2	Cinnamyl carbanilate	253	C16H15NO2	025076-44-2	28
3	(2E,4Z)-3-Methoxy-2,4-hexadiened...	134	C7H6N2O	001789-46-4	9
4	1,3,8-p-Menthatriene	134	C10H14	018368-95-1	9
5	1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	9



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111920\
Data File : VD067703.D
Acq On : 19 Nov 2020 12:54
Operator : VA/SY
Sample : L4778-06
Misc : 4.07G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
VNJ-223

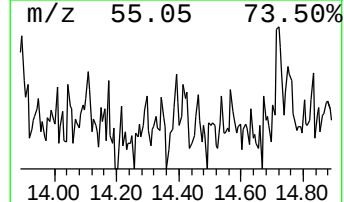
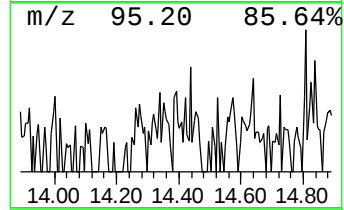
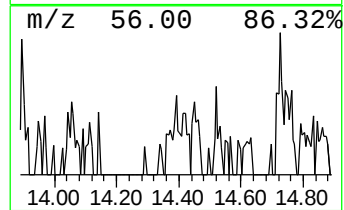
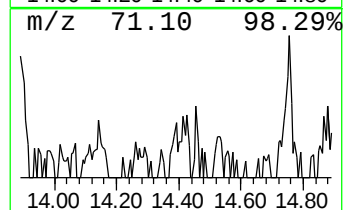
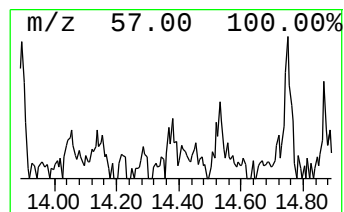
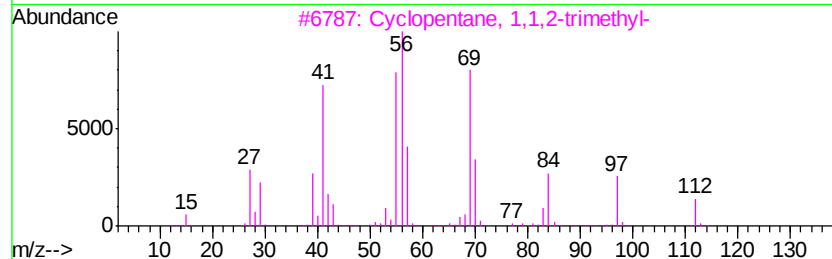
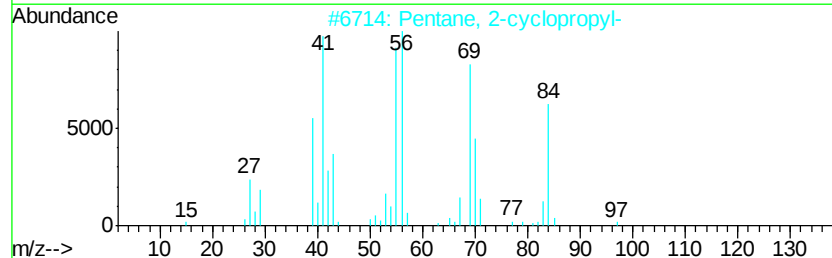
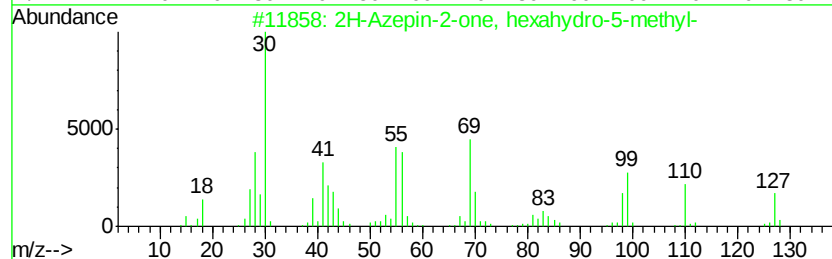
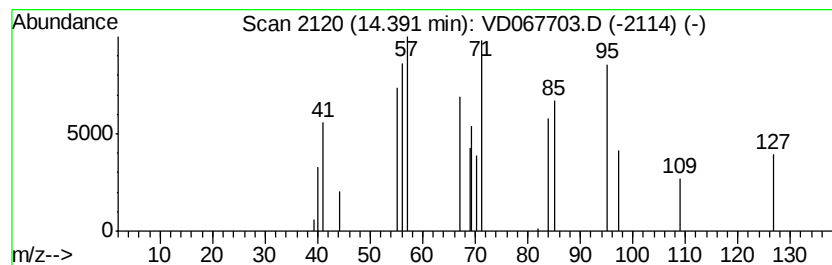
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.M
Quant Title : SW846 8260

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

Peak Number 5 unknown14.39 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.39	5.52 ug/l	10164	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-Azepin-2-one, hexahydro-5-met...	127	C7H13NO	002210-07-3	35
2		Pentane, 2-cyclopropyl-	112	C8H16	005458-16-2	35
3		Cyclopentane, 1,1,2-trimethyl-	112	C8H16	004259-00-1	35
4		Oxalic acid, cyclobutyl hexadecyl...	368	C22H40O4	1000309-70-6	28
5		1-Heptanol, 6-methyl-	130	C8H18O	001653-40-3	25



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111920\
Data File : VD067703.D
Acq On : 19 Nov 2020 12:54
Operator : VA/SY
Sample : L4778-06
Misc : 4.07G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
VNJ-223

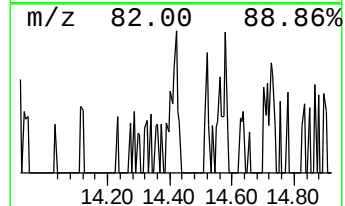
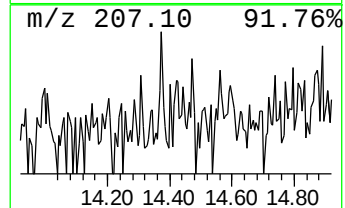
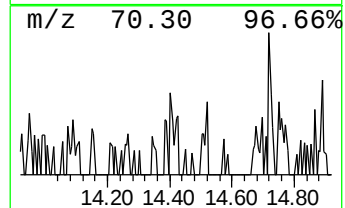
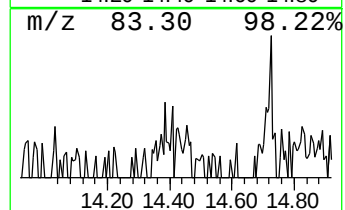
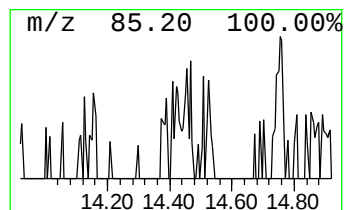
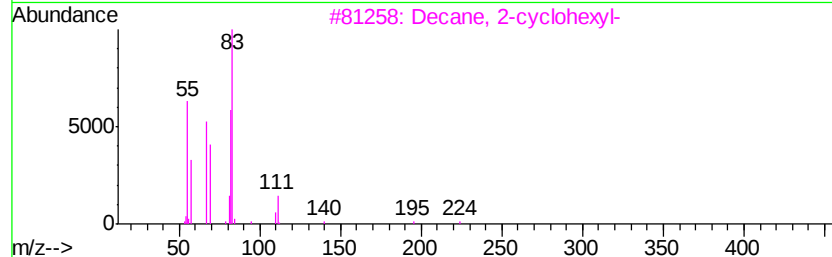
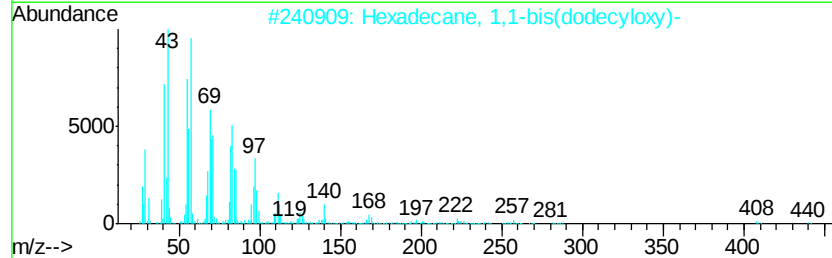
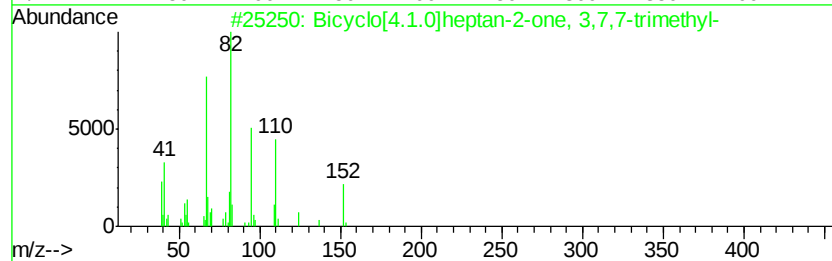
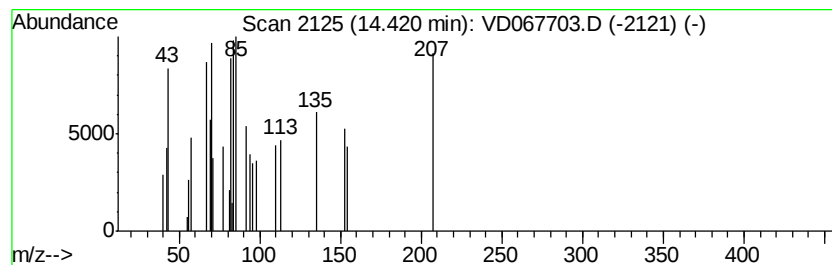
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.M
Quant Title : SW846 8260

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

Peak Number 6 unknown14.42 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.42	6.99 ug/l	12852	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[4.1.0]heptan-2-one, 3,7,...	152	C10H16O	000497-62-1	14
2		Hexadecane, 1,1-bis(dodecyloxy)-	595	C40H82O2	056554-64-4	10
3		Decane, 2-cyclohexyl-	224	C16H32	013151-73-0	10
4		Azeleoneitrile	150	C9H14N2	001675-69-0	10
5		Tricyclo[4.2.1.0(3,7)]nonane-2,6...	154	C9H14O2	106435-10-3	10



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111920\
Data File : VD067703.D
Acq On : 19 Nov 2020 12:54
Operator : VA/SY
Sample : L4778-06
Misc : 4.07G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
VNJ-223

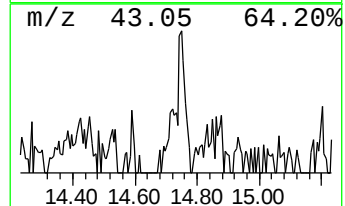
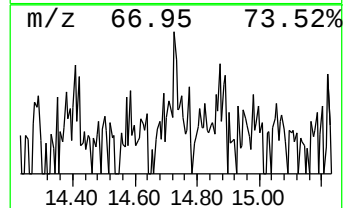
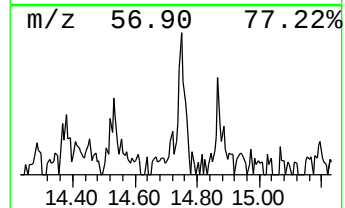
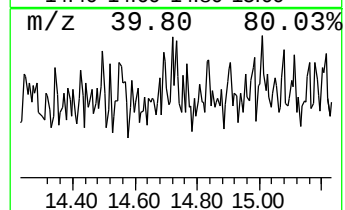
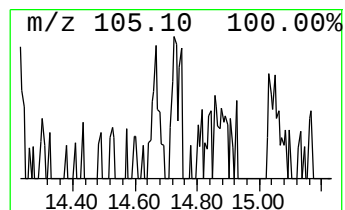
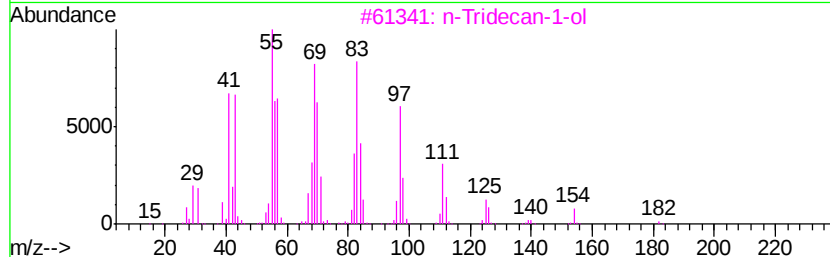
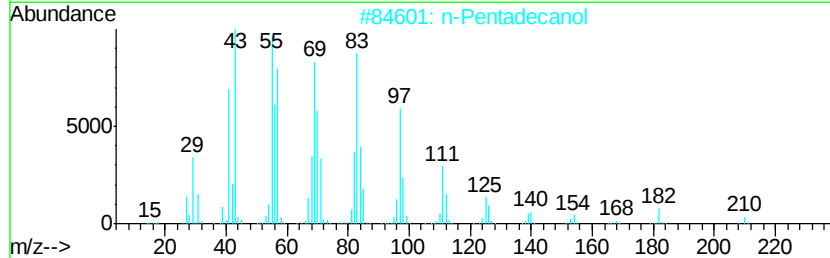
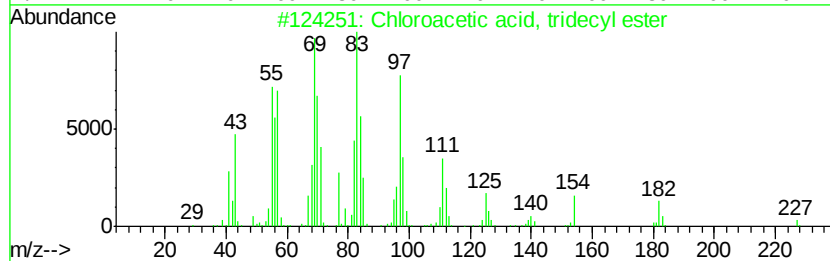
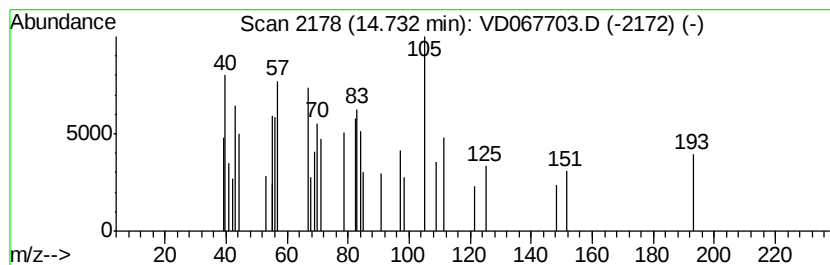
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.M
Quant Title : SW846 8260

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

Peak Number 7 Chloroacetic acid, tridecyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.73	8.30 ug/l	15266	1,4-Dichlorobenzene-d4	13.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Chloroacetic acid, tridecyl ester	276	C15H29ClO2	018277-85-5	72
2		n-Pentadecanol	228	C15H32O	000629-76-5	64
3		n-Tridecan-1-ol	200	C13H28O	000112-70-9	64
4		Carbonic acid, tridecyl 2,2,2-tr...	374	C16H29Cl3O3	1000314-55-9	64
5		1-Undecanol	172	C11H24O	000112-42-5	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111920\
Data File : VD067703.D
Acq On : 19 Nov 2020 12:54
Operator : VA/SY
Sample : L4778-06
Misc : 4.07G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
VNJ-223

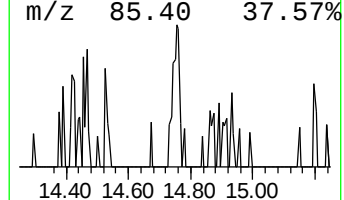
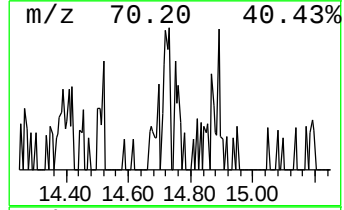
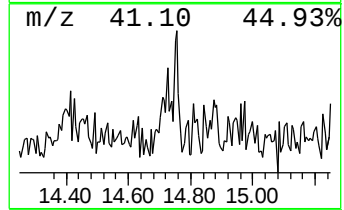
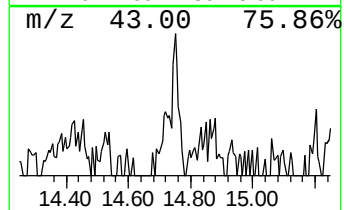
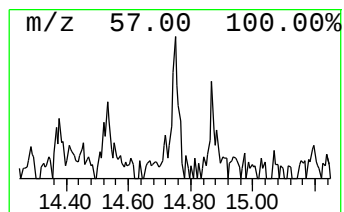
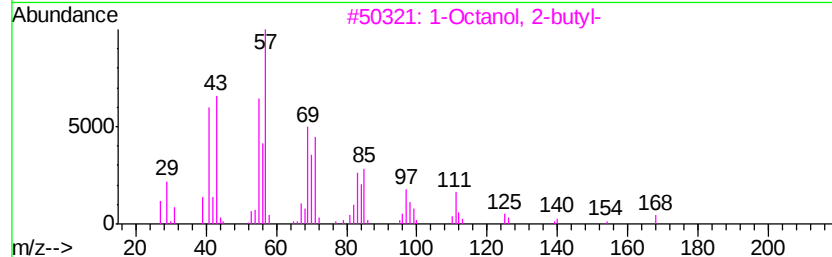
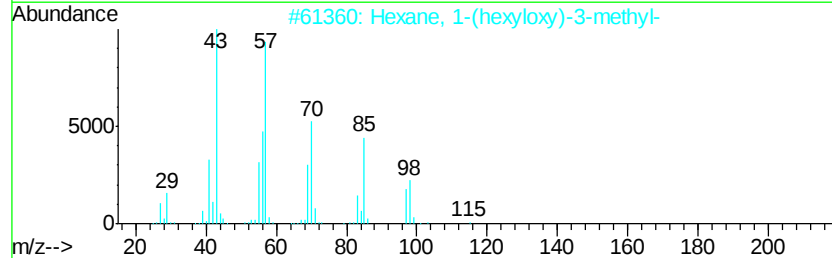
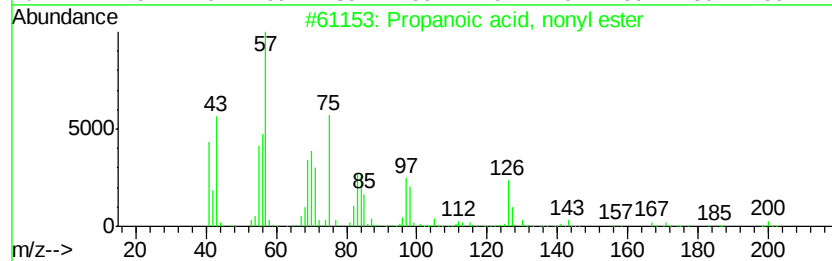
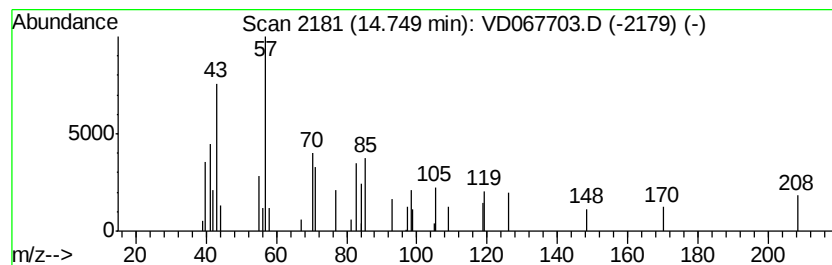
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.M
Quant Title : SW846 8260

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

Peak Number 8 unknown14.75 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.75	8.50 ug/l	15639	1,4-Dichlorobenzene-d4	13.58

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, nonyl ester	200	C12H24O2	053184-67-1	47
2			Hexane, 1-(hexyloxy)-3-methyl-	200	C13H28O	074421-18-4	35
3			1-Octanol, 2-butyl-	186	C12H26O	003913-02-8	32
4			1-Dodecene	168	C12H24	000112-41-4	25
5			Nonane, 4-methyl-	142	C10H22	017301-94-9	22



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111920\
Data File : VD067703.D
Acq On : 19 Nov 2020 12:54
Operator : VA/SY
Sample : L4778-06
Misc : 4.07G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

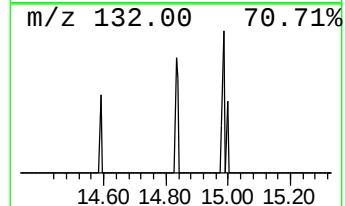
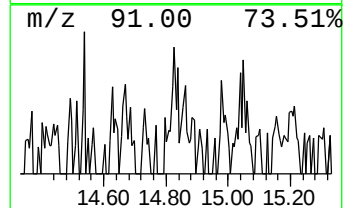
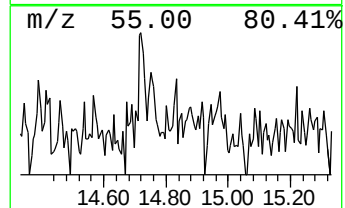
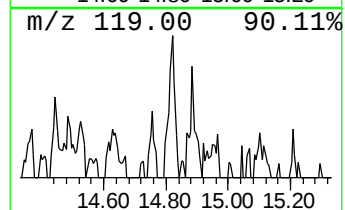
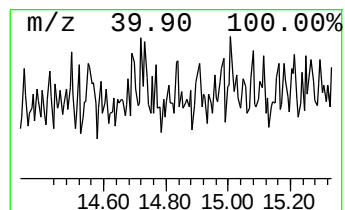
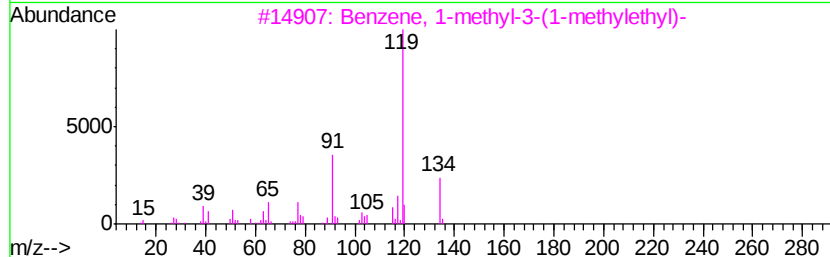
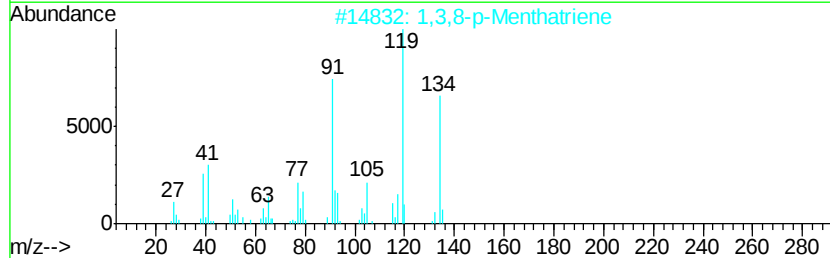
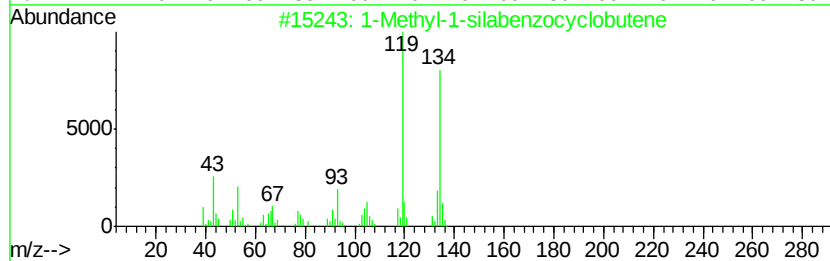
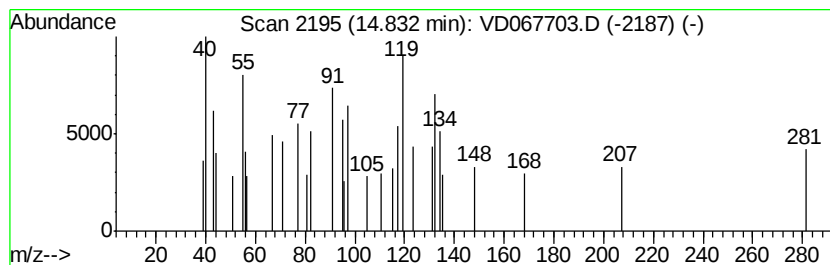
Instrument :
MSVOA_D
ClientSampleId :
VNJ-223

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.M
Quant Title : SW846 8260

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

Peak Number 9 unknown14.83 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.83	8.94 ug/l	16441	1,4-Dichlorobenzene-d4	13.58	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Methyl-1-silabenzocyclobutene	134	C8H10Si	160841-06-5	25
2	1,3,8-p-Menthatriene	134	C10H14	018368-95-1	22
3	Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	22
4	3-Methyl-2,3-dihydro-benzofuran	134	C9H10O	013524-73-7	18
5	4,4-Dimethyl-2-cyclopenten-1-one	110	C7H10O	022748-16-9	11



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD111920\
Data File : VD067703.D
Acq On : 19 Nov 2020 12:54
Operator : VA/SY
Sample : L4778-06
Misc : 4.07G/5.00ml/MSVOA_D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
VNJ-223

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.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
.beta.-Ocimene	12.50	6.5	ug/l	14433	3	11.65	110906	50.0
unknown13.90	13.90	12.4	ug/l	22715	4	13.58	91997	50.0
unknown14.11	14.11	5.6	ug/l	10333	4	13.58	91997	50.0
unknown14.39	14.39	5.5	ug/l	10164	4	13.58	91997	50.0
unknown14.42	14.42	7.0	ug/l	12852	4	13.58	91997	50.0
Chloroacetic acid...	14.73	8.3	ug/l	15266	4	13.58	91997	50.0
unknown14.75	14.75	8.5	ug/l	15639	4	13.58	91997	50.0
unknown14.83	14.83	8.9	ug/l	16441	4	13.58	91997	50.0