

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD102020\
 Data File : VD067317.D
 Acq On : 20 Oct 2020 12:06
 Operator : VA/SY
 Sample : L4222-01
 Misc : 6.99G/5.00ml/MSVOA D/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 EO-03-101920

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\82D101320S.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	7.920	1007	1020	1025	rBV	252891	589972	31.45%	2.242%
2	7.985	1025	1031	1042	rVB	416567	946985	50.48%	3.599%
3	8.338	1079	1091	1102	rBV	201323	471488	25.13%	1.792%
4	8.867	1170	1181	1192	rBV	594361	1200482	64.00%	4.562%
5	10.344	1422	1432	1441	rBV	805211	1454980	77.56%	5.529%
6	11.455	1615	1621	1629	rVB4	27945	66341	3.54%	0.252%
7	11.650	1646	1654	1665	rBV	800295	1389873	74.09%	5.282%
8	11.779	1672	1676	1681	rBV4	17933	26628	1.42%	0.101%
9	11.838	1681	1686	1689	rVV6	17210	34407	1.83%	0.131%
10	11.891	1690	1695	1708	rVB4	51289	153642	8.19%	0.584%
11	12.085	1721	1728	1732	rVB6	15279	32501	1.73%	0.124%
12	12.155	1732	1740	1747	rBV5	87277	225860	12.04%	0.858%
13	12.273	1757	1760	1765	rVB3	60545	82790	4.41%	0.315%
14	12.385	1765	1779	1786	rBV5	134944	422788	22.54%	1.607%
15	12.479	1790	1795	1798	rBV4	49936	97539	5.20%	0.371%
16	12.638	1816	1822	1832	rBV	735805	1275073	67.97%	4.846%
17	12.726	1833	1837	1840	rBV3	51850	76069	4.06%	0.289%
18	12.802	1846	1850	1856	rVB3	151225	284280	15.15%	1.080%
19	12.885	1856	1864	1868	rBV4	106339	260748	13.90%	0.991%
20	12.961	1868	1877	1883	rVV4	305192	857900	45.73%	3.260%
21	13.014	1883	1886	1891	rVB4	125369	204869	10.92%	0.779%
22	13.102	1896	1901	1905	rBV3	146471	202329	10.79%	0.769%
23	13.155	1905	1910	1912	rBV2	144908	249052	13.28%	0.946%
24	13.191	1912	1916	1922	rVB	392217	586453	31.26%	2.229%
25	13.249	1923	1926	1930	rVB	92044	130334	6.95%	0.495%
26	13.320	1933	1938	1942	rBV2	152715	255165	13.60%	0.970%
27	13.367	1943	1946	1950	rVV4	85812	116848	6.23%	0.444%
28	13.467	1955	1963	1967	rVV7	316207	758250	40.42%	2.882%
29	13.508	1967	1970	1974	rVV2	301371	471044	25.11%	1.790%
30	13.579	1974	1982	1991	rVB	972347	1875885	100.00%	7.129%
31	13.649	1991	1994	2000	rBV5	195396	296365	15.80%	1.126%
32	13.720	2001	2006	2010	rVV5	132269	237864	12.68%	0.904%
33	13.902	2032	2037	2042	rBV2	864971	1483069	79.06%	5.636%
34	13.949	2042	2045	2050	rVB3	194753	280117	14.93%	1.065%

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Sampling : 1
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Stop Thrs : 0
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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

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Title : SW846 8260

35	14.008	2051	2055	2059	rBV4	184392	277804	14.81%	1.056%
36	14.055	2059	2063	2067	rBV4	162318	317560	16.93%	1.207%
37	14.161	2078	2081	2087	rVB2	306266	483707	25.79%	1.838%
38	14.226	2087	2092	2097	rBV4	268977	498802	26.59%	1.896%
39	14.291	2097	2103	2109	rVV5	314723	583401	31.10%	2.217%
40	14.349	2109	2113	2116	rBV3	172275	269253	14.35%	1.023%
41	14.414	2118	2124	2129	rVB3	812843	1480814	78.94%	5.628%
42	14.532	2139	2144	2147	rBV4	196362	325687	17.36%	1.238%
43	14.573	2147	2151	2157	rVV4	571269	1089134	58.06%	4.139%
44	14.632	2158	2161	2168	rVB5	244139	473711	25.25%	1.800%
45	14.767	2180	2184	2190	rBV7	185327	373348	19.90%	1.419%
46	14.832	2191	2195	2199	rVV5	276269	531892	28.35%	2.021%
47	14.885	2200	2204	2211	rVB5	432206	834599	44.49%	3.172%
48	15.008	2221	2225	2228	rBV2	141713	209785	11.18%	0.797%
49	15.102	2236	2241	2245	rBV4	247043	393983	21.00%	1.497%
50	15.208	2256	2259	2261	rBV2	99945	144454	7.70%	0.549%
51	15.373	2283	2287	2292	rBV3	266046	439626	23.44%	1.671%
52	15.685	2337	2340	2343	rBV4	73556	99178	5.29%	0.377%
53	16.226	2427	2432	2438	rVB5	112797	232570	12.40%	0.884%
54	16.355	2451	2454	2460	rVB6	89957	156178	8.33%	0.594%

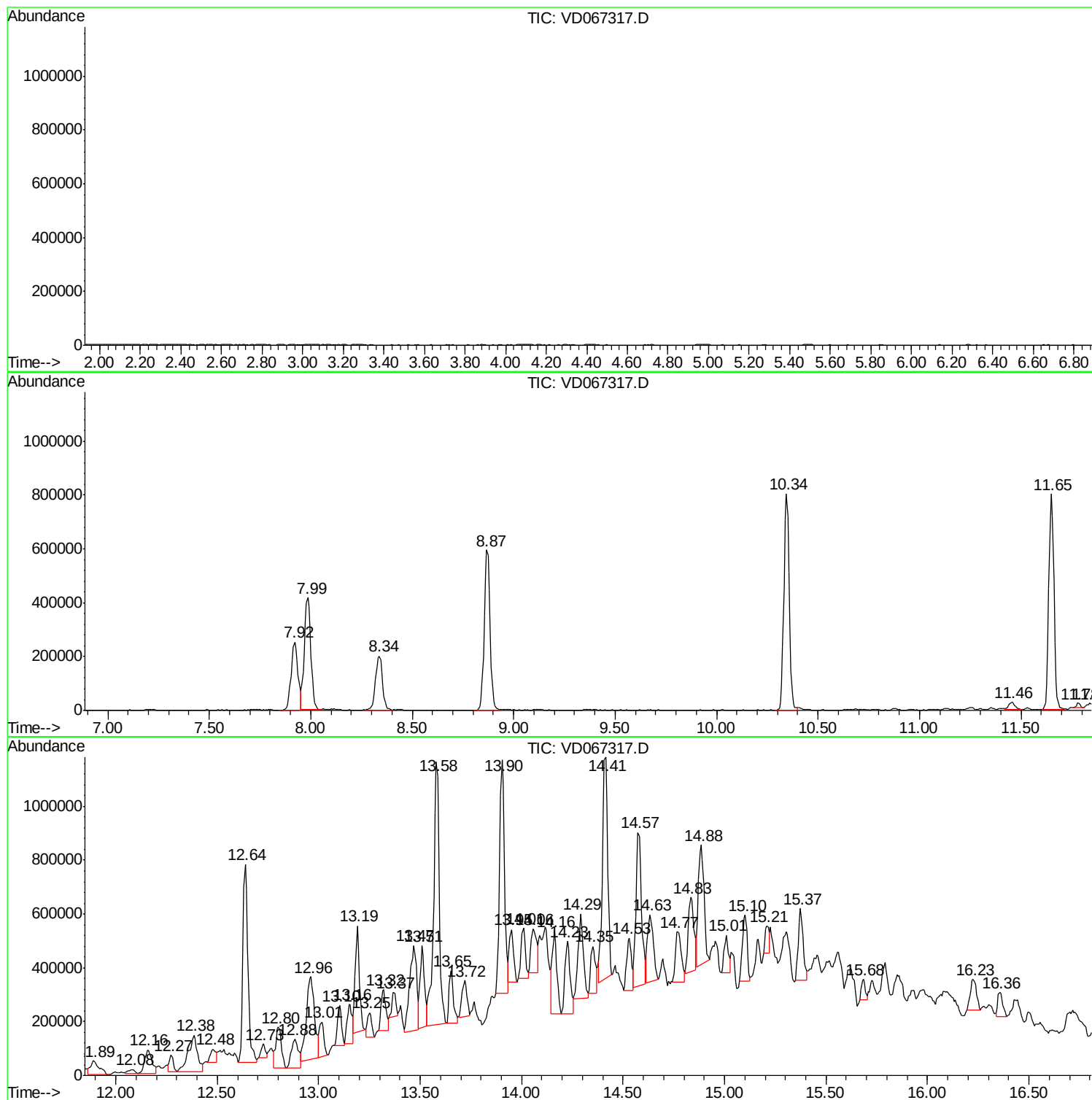
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D101320S.M
Quant Title : SW846 8260

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



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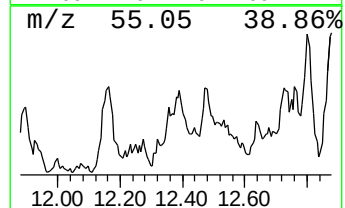
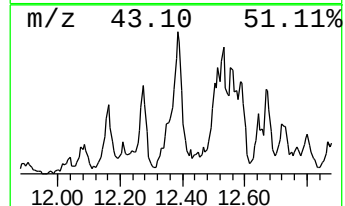
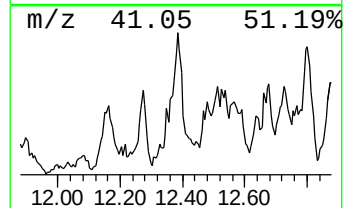
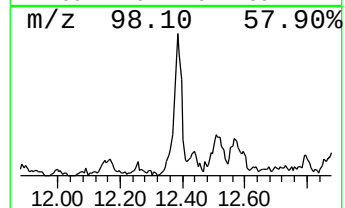
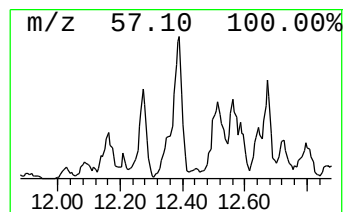
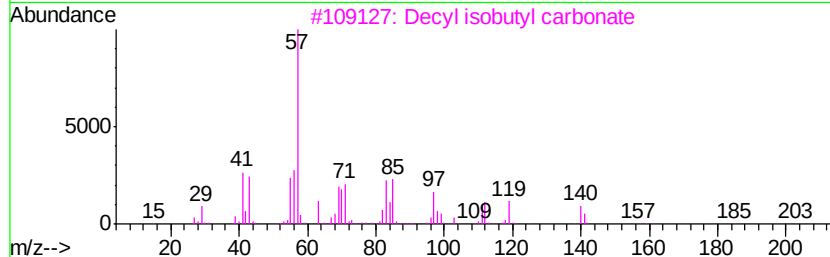
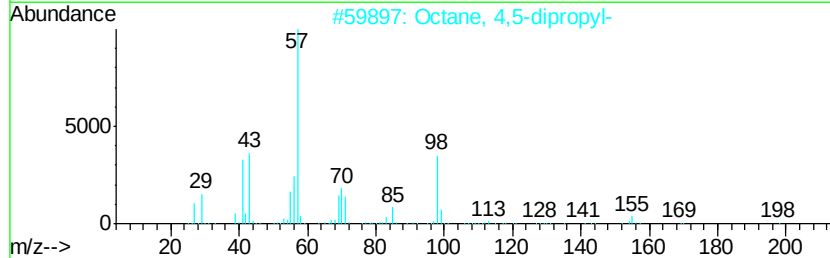
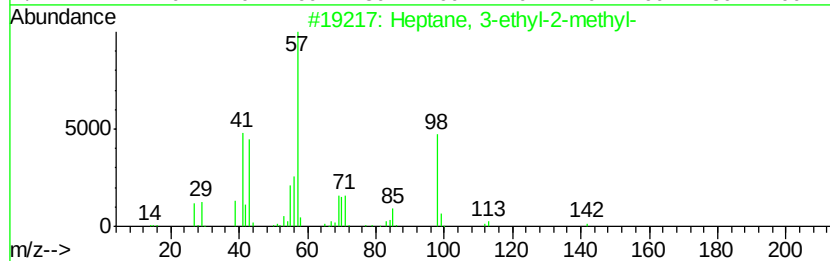
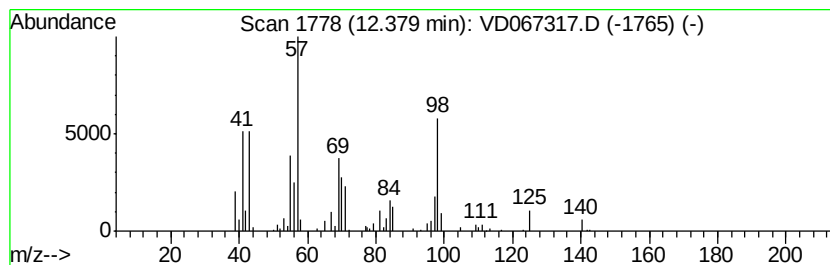
Instrument :
MSVOA_D
ClientSampleId :
EO-03-101920

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D101320S.M
Quant Title : SW846 8260

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

Peak Number 1 Heptane, 3-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.38	15.21 ug/l	422788	Chlorobenzene-d5	11.65		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	74
2		Octane, 4,5-dipropyl-	198	C14H30	020905-05-9	72
3		Decyl isobutyl carbonate	258	C15H30O3	959226-07-4	53
4		Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	52
5		1-Butanamine, 2-methyl-N-(2-meth...	155	C10H21N	054518-97-7	47



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Misc : 6.99G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

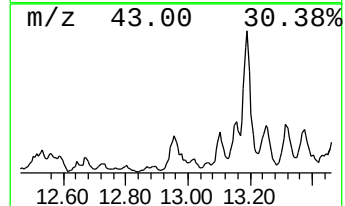
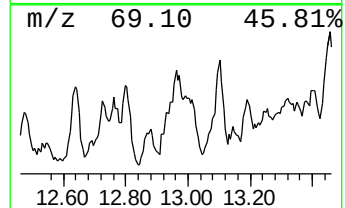
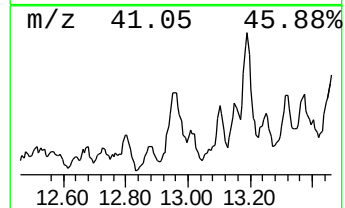
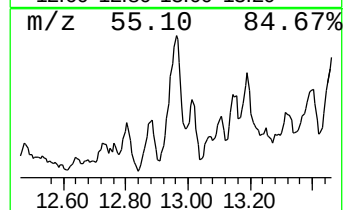
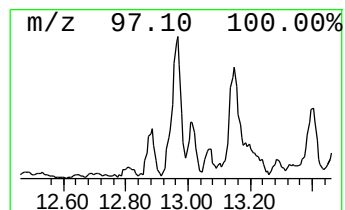
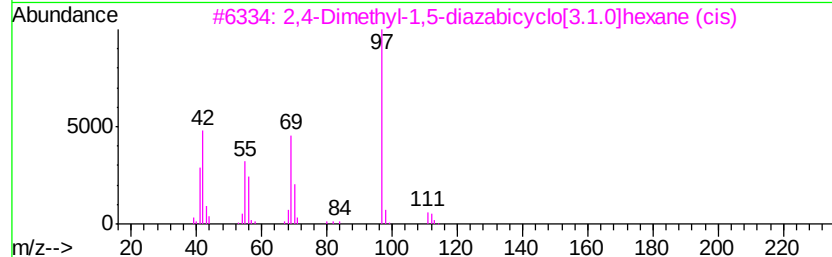
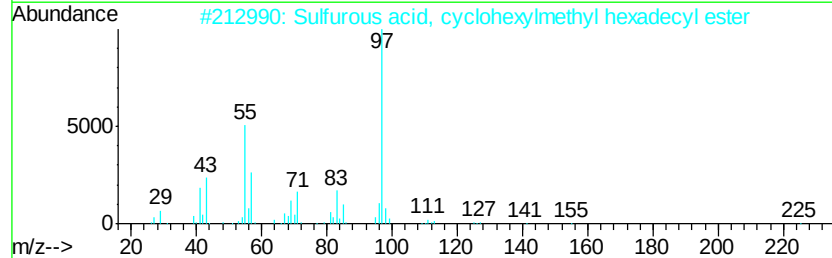
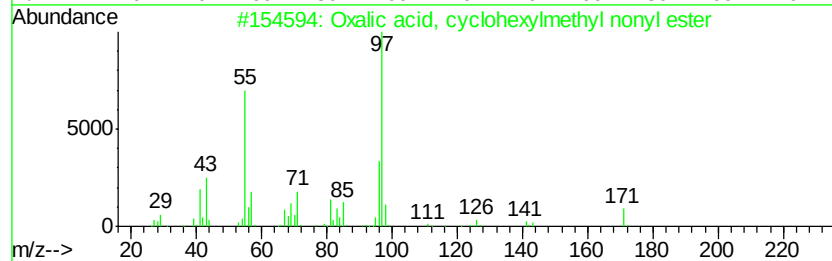
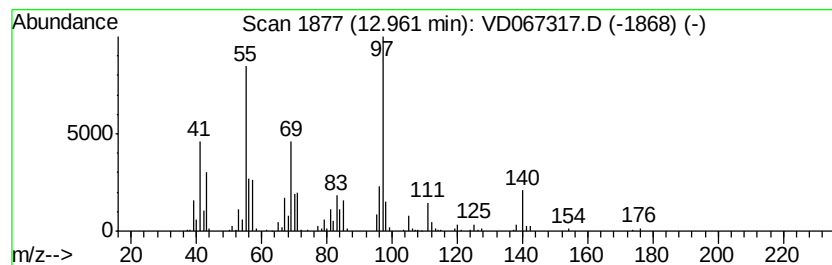
Instrument :
MSVOA_D
ClientSampleId :
EO-03-101920

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D101320S.M
Quant Title : SW846 8260

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

Peak Number 2 Oxalic acid, cyclohexylmeth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	22.87 ug/l	857900	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Oxalic acid, cyclohexylmethvl no...	312 C18H32O4	1000309-68-6	53
2	Sulfurous acid, cyclohexylmethvl...	402 C23H46O3S	1000309-22-4	50
3	2,4-Dimethyl-1,5-diazabicyclo[3.1.0]hexane (cis)	112 C6H12N2	100463-01-2	49
4	Cyclohexane, 1-ethyl-2-methyl-	126 C9H18	003728-54-9	47
5	3-Heptene, 4-propyl-	140 C10H20	004485-13-6	46



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Misc : 6.99G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

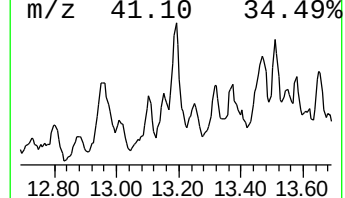
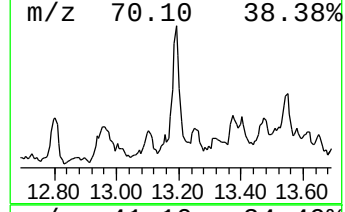
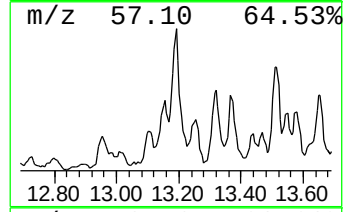
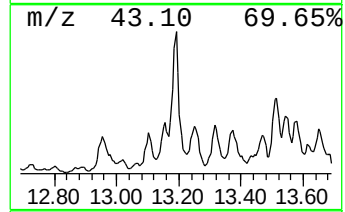
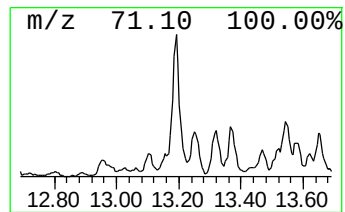
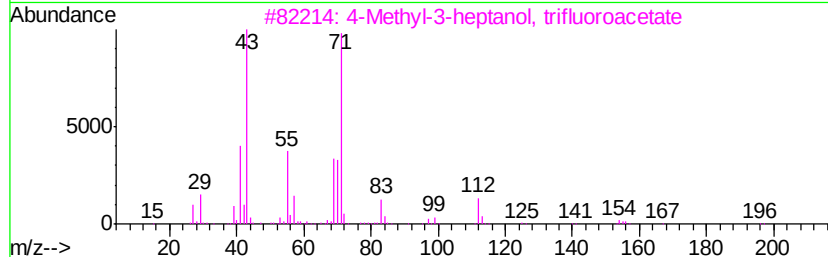
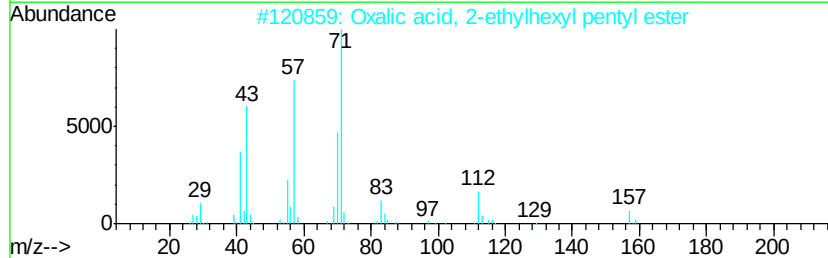
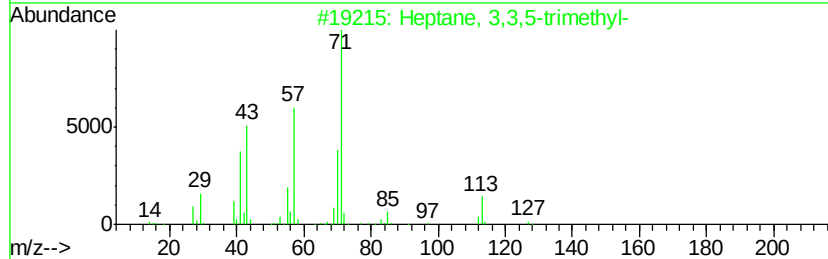
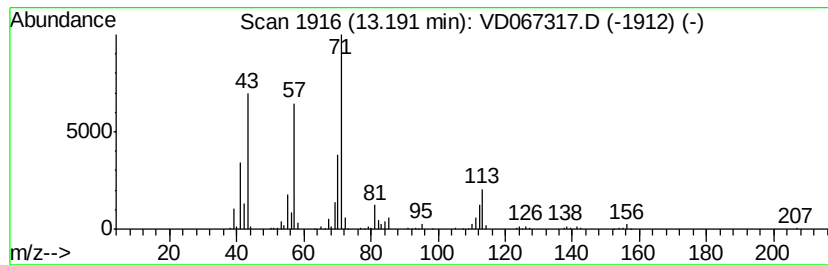
Instrument :
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ClientSampled :
EO-03-101920

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Quant Title : SW846 8260

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

Peak Number 3 Heptane, 3,3,5-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.19	15.63 ug/l	586453	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptane, 3,3,5-trimethyl-	142 C10H22	007154-80-5	80
2	Oxalic acid, 2-ethylhexyl pentyl...	272 C15H28O4	1000309-38-7	59
3	4-Methyl-3-heptanol, trifluoroac...	226 C10H17F3O2	1000365-51-8	59
4	Octane, 3-ethyl-	142 C10H22	005881-17-4	59
5	Octane, 3,3-dimethyl-	142 C10H22	004110-44-5	53



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ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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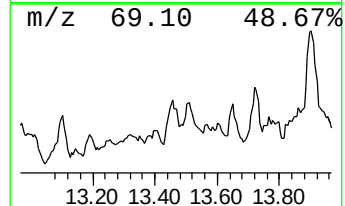
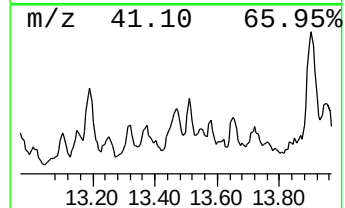
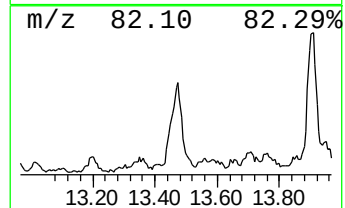
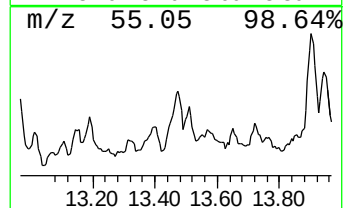
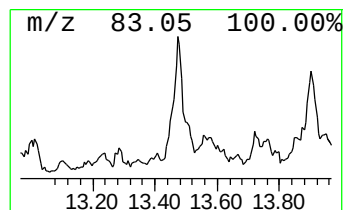
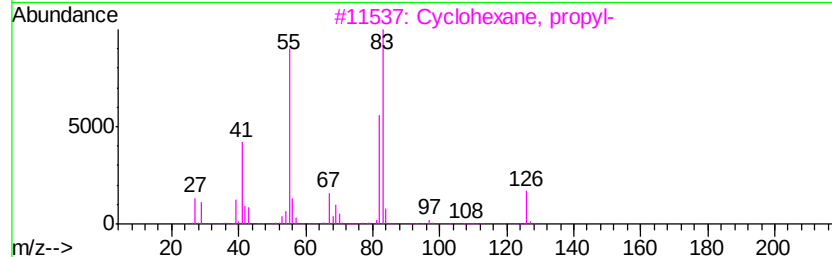
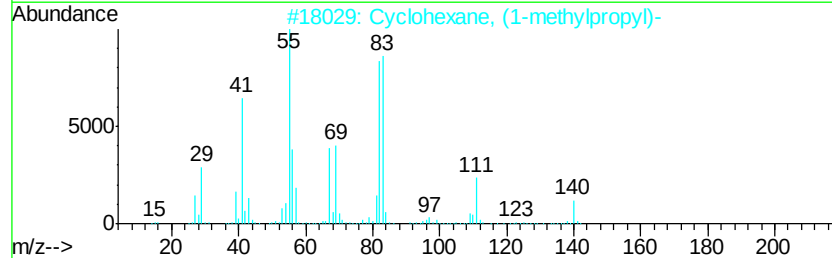
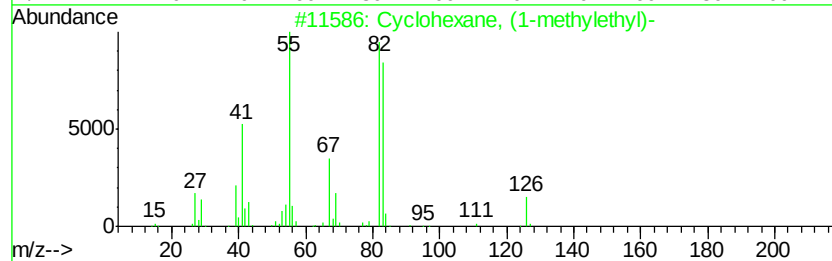
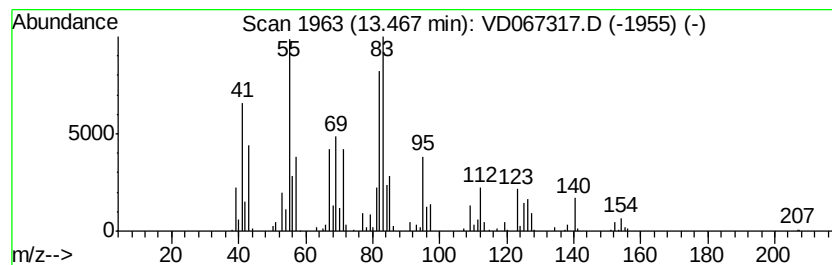
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D101320S.M
Quant Title : SW846 8260

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

Peak Number 4 unknown13.47 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	20.21 ug/l	758250	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	46
2		Cyclohexane, (1-methylpropyl)-	140	C10H20	007058-01-7	46
3		Cyclohexane, propyl-	126	C9H18	001678-92-8	38
4		Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	38
5		Cyclohexane, pentyl-	154	C11H22	004292-92-6	38



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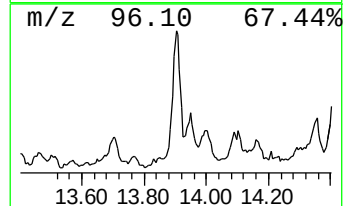
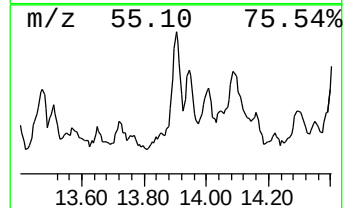
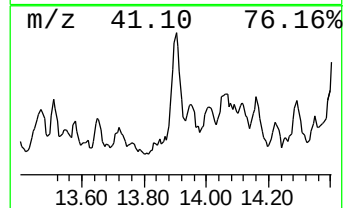
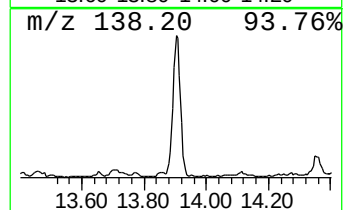
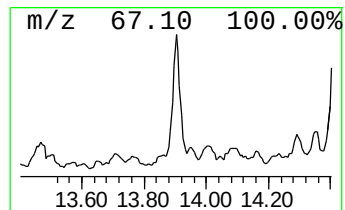
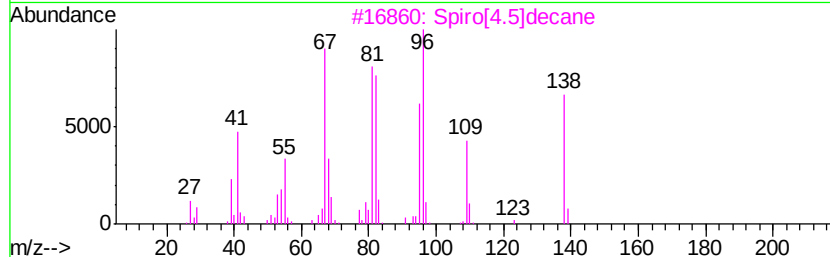
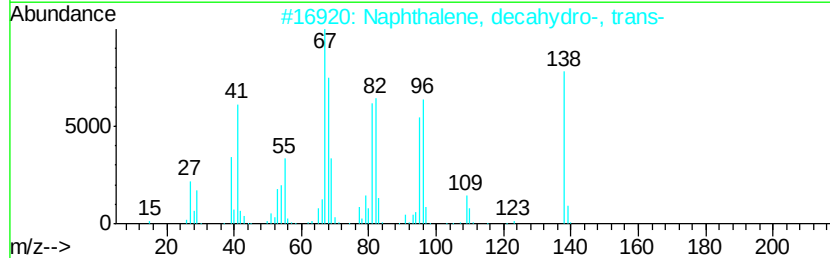
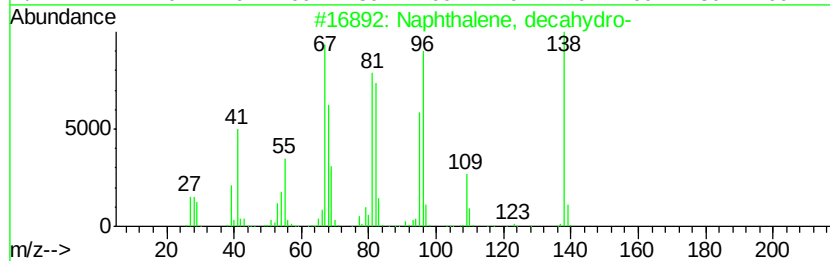
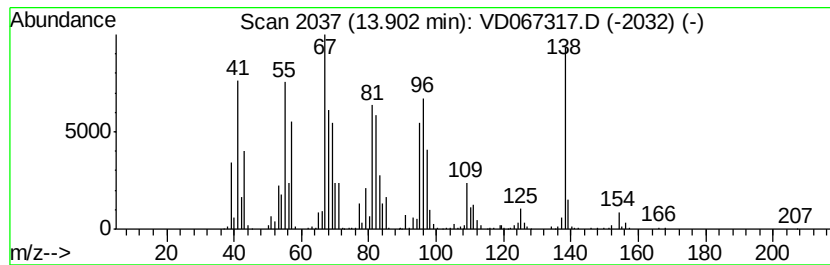
Instrument :
MSVOA_D
ClientSampleId :
EO-03-101920

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D101320S.M
Quant Title : SW846 8260

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

Peak Number 5 Naphthalene, decahydro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	39.53 ug/l	1483070	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, decahydro-	138 C10H18	000091-17-8 97
2		Naphthalene, decahydro-, trans-	138 C10H18	000493-02-7 87
3		Spiro[4.5]decane	138 C10H18	000176-63-6 76
4		1,1'-Bicyclopentyl	138 C10H18	001636-39-1 76
5		Naphthalene, decahydro-, cis-	138 C10H18	000493-01-6 76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD102020\
Data File : VD067317.D
Acq On : 20 Oct 2020 12:06
Operator : VA/SY
Sample : L4222-01
Misc : 6.99G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

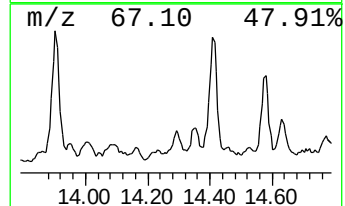
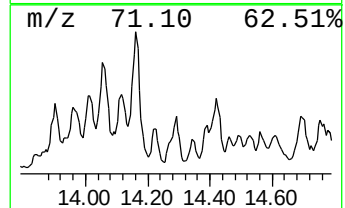
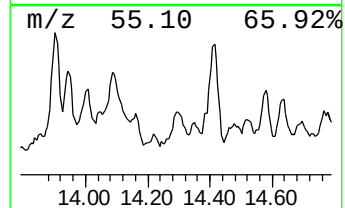
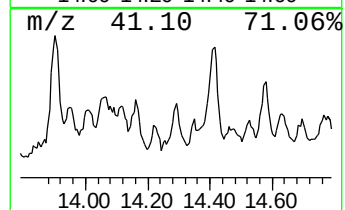
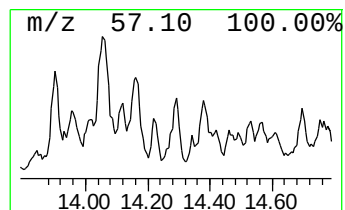
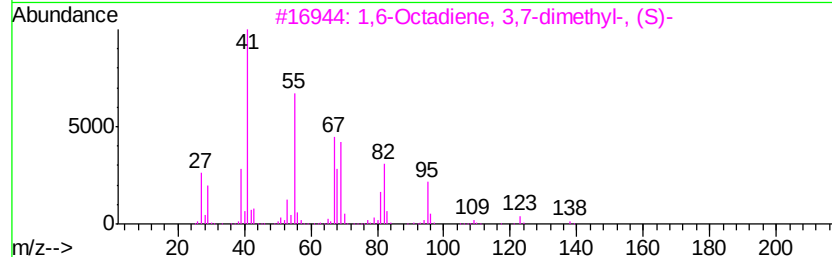
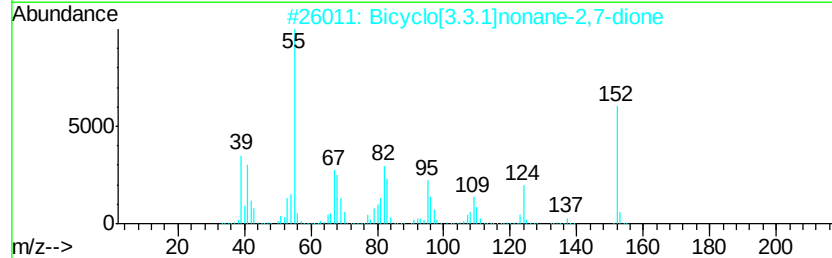
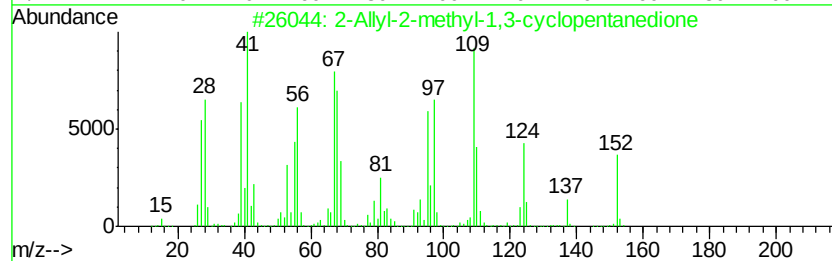
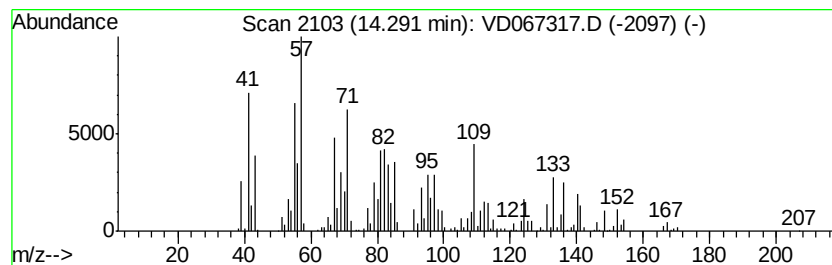
Instrument :
MSVOA_D
ClientSampleID :
EO-03-101920

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D101320S.M
Quant Title : SW846 8260

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

Peak Number 6 unknown14.29 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	15.55 ug/l	583401	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-Allyl-2-methyl-1,3-cyclopentan...	152 C9H12O2	026828-48-8	38
2	Bicyclo[3.3.1]nonane-2,7-dione	152 C9H12O2	1000210-30-3	30
3	1,6-Octadiene, 3,7-dimethyl-, (S)-	138 C10H18	010281-55-7	25
4	Oxalic acid, allyl hexadecyl ester	354 C21H38O4	1000309-24-4	25
5	5-Ethyldecane	170 C12H26	017302-36-2	25



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD102020\
Data File : VD067317.D
Acq On : 20 Oct 2020 12:06
Operator : VA/SY
Sample : L4222-01
Misc : 6.99G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

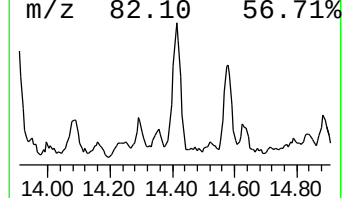
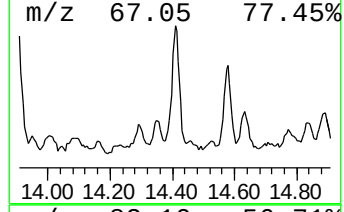
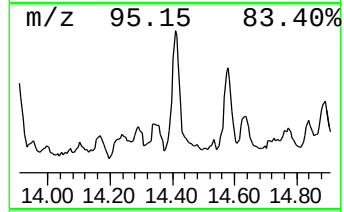
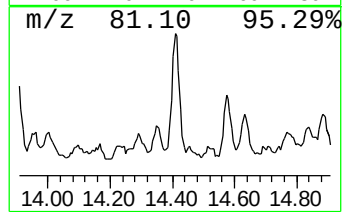
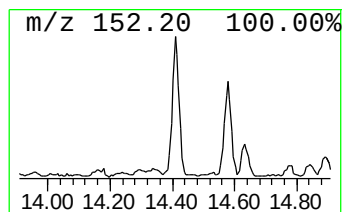
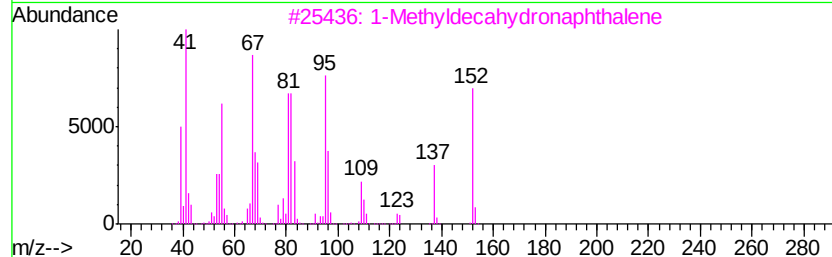
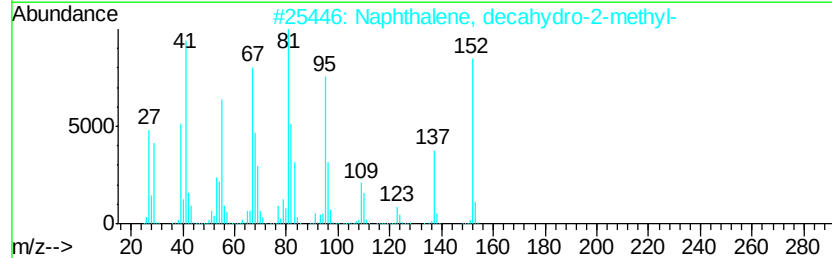
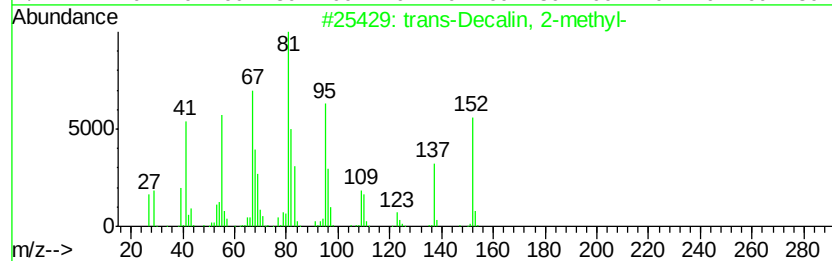
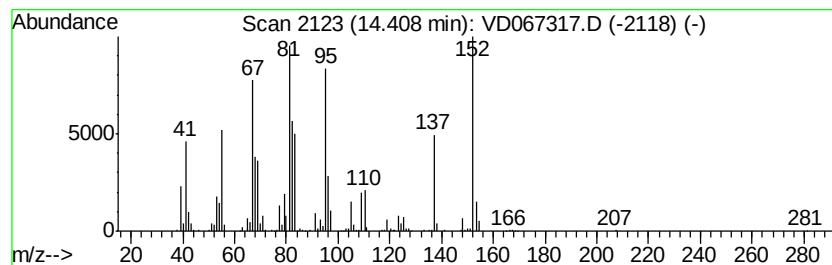
Instrument :
MSVOA_D
ClientSampleId :
EO-03-101920

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D101320S.M
Quant Title : SW846 8260

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

Peak Number 7 trans-Decalin, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.41	39.47 ug/l	1480810	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	trans-Decalin, 2-methyl-	152 C11H20	1000152-47-3	94
2	Naphthalene, decahydro-2-methyl-	152 C11H20	002958-76-1	94
3	1-Methyldecahydronaphthalene	152 C11H20	002958-75-0	86
4	cis-Decalin, 2-syn-methyl-	152 C11H20	1000155-85-6	72
5	Cyclohexanone, 5-methyl-2-(1-met...	152 C10H16O	015932-80-6	62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD102020\
Data File : VD067317.D
Acq On : 20 Oct 2020 12:06
Operator : VA/SY
Sample : L4222-01
Misc : 6.99G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

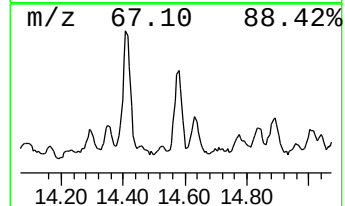
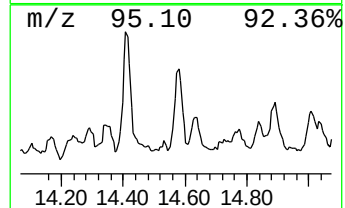
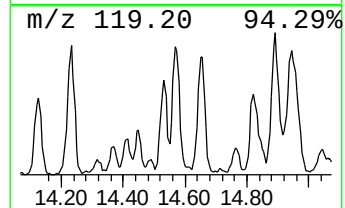
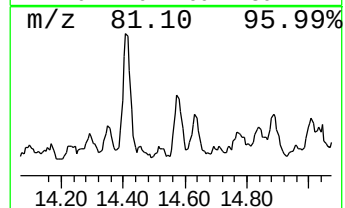
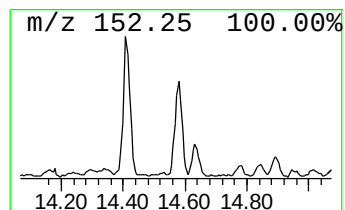
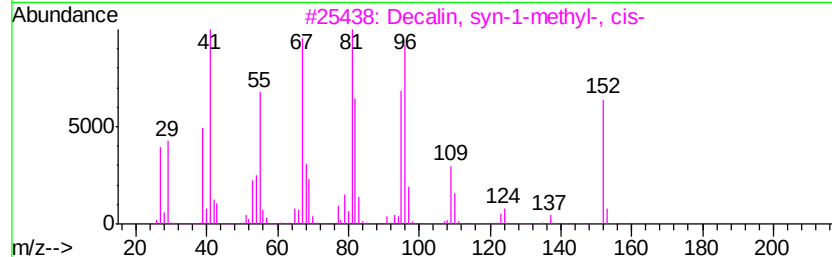
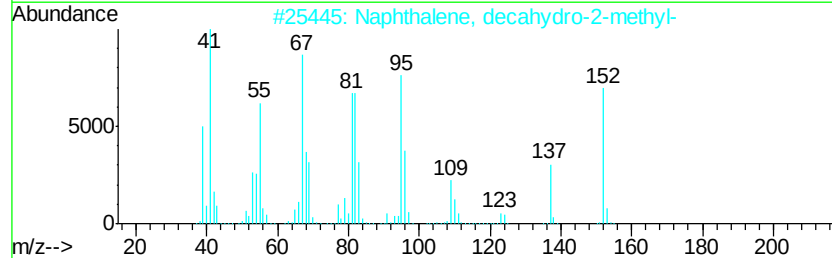
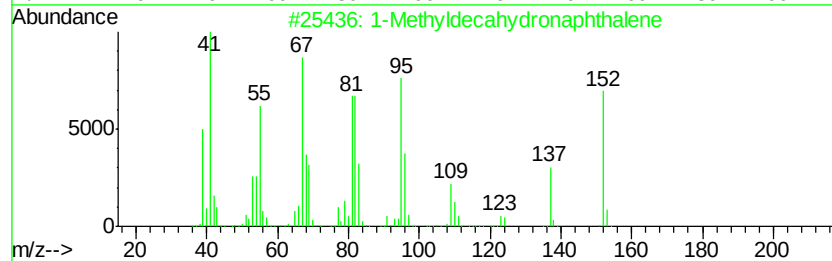
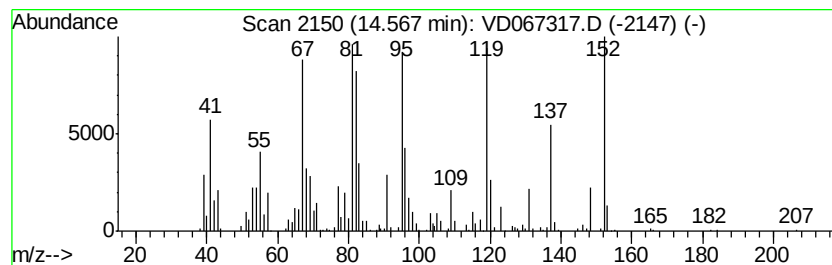
Instrument :
MSVOA_D
ClientSampleId :
EO-03-101920

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D101320S.M
Quant Title : SW846 8260

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

Peak Number 8 1-Methyldecahydronaphthalene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	29.03 ug/l	1089130	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1-Methyldecahydronaphthalene	152 C11H20	002958-75-0 86
2		Naphthalene, decahydro-2-methyl-	152 C11H20	002958-76-1 86
3		Decalin, syn-1-methyl-, cis-	152 C11H20	1000158-89-1 64
4		cis-Decalin, 2-syn-methyl-	152 C11H20	1000155-85-6 62
5		trans-Decalin, 2-methyl-	152 C11H20	1000152-47-3 52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD102020\
Data File : VD067317.D
Acq On : 20 Oct 2020 12:06
Operator : VA/SY
Sample : L4222-01
Misc : 6.99G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

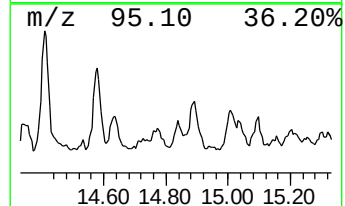
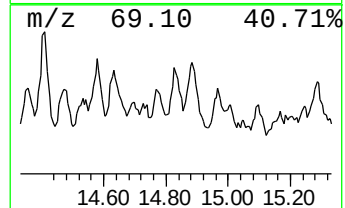
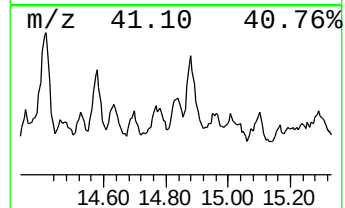
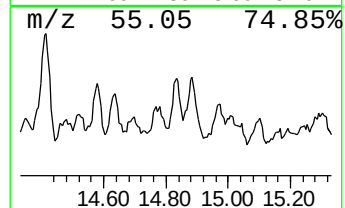
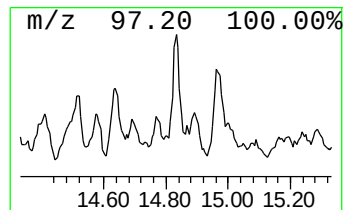
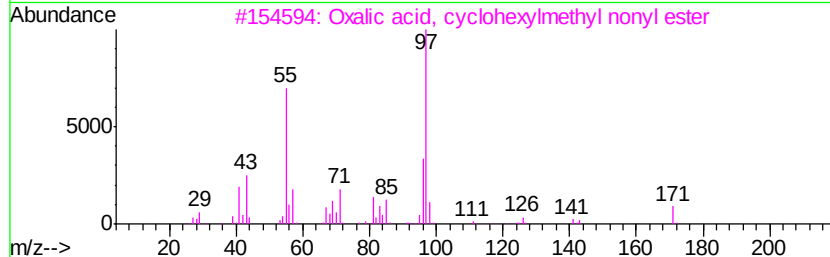
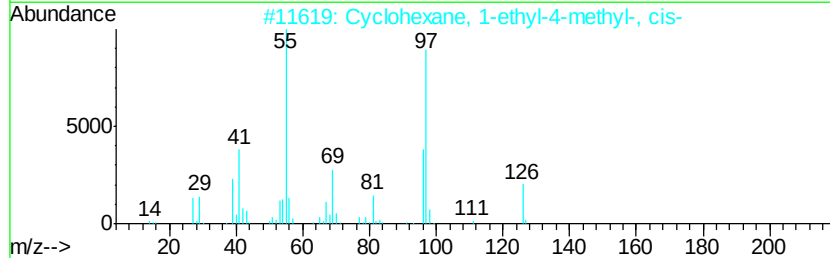
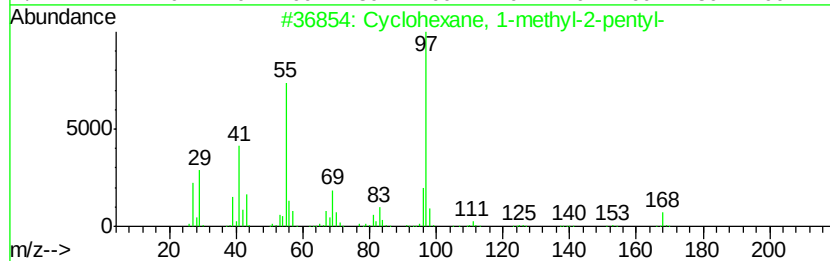
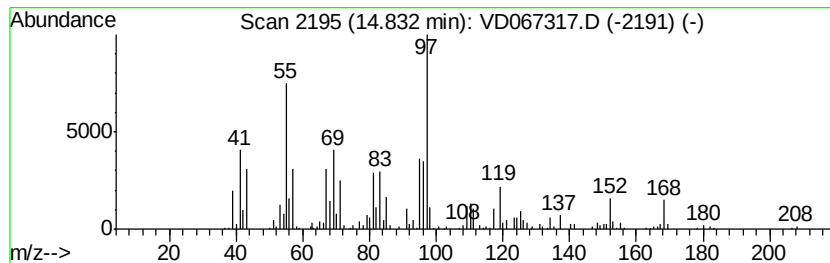
Instrument :
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EO-03-101920

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D101320S.M
Quant Title : SW846 8260

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

Peak Number 9 unknown14.83 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	14.18 ug/l	531892	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexane, 1-methyl-2-pentyl-	168 C12H24	054411-01-7 46
2		Cyclohexane, 1-ethyl-4-methyl-, ...	126 C9H18	004926-78-7 45
3		Oxalic acid, cyclohexylmethyl no...	312 C18H32O4	1000309-68-6 43
4		Cyclohexanecarboxylic acid, 4-pe...	368 C22H28N2O3	133856-87-8 43
5		2-Butyl-3,4,5,6-tetrahydropyridine	139 C9H17N	001462-94-8 43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD102020\
Data File : VD067317.D
Acq On : 20 Oct 2020 12:06
Operator : VA/SY
Sample : L4222-01
Misc : 6.99G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
EO-03-101920

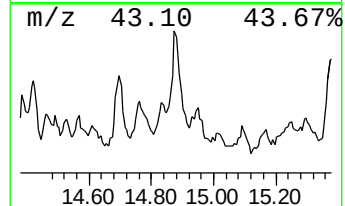
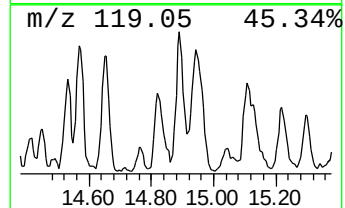
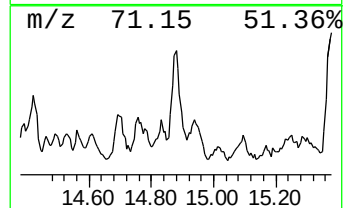
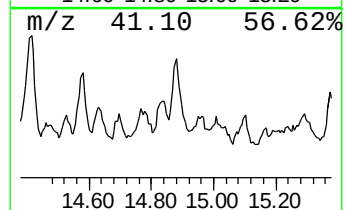
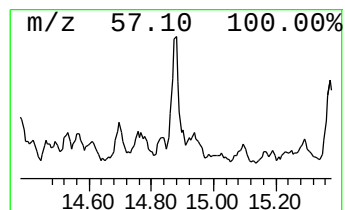
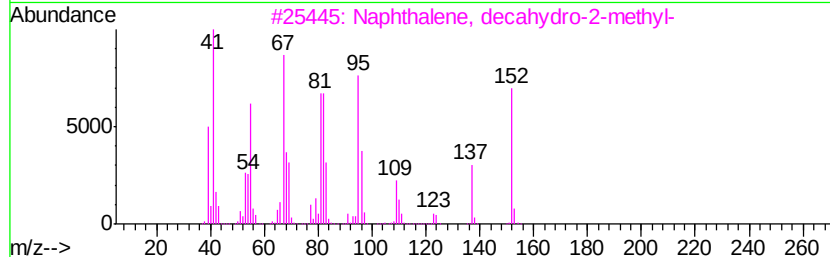
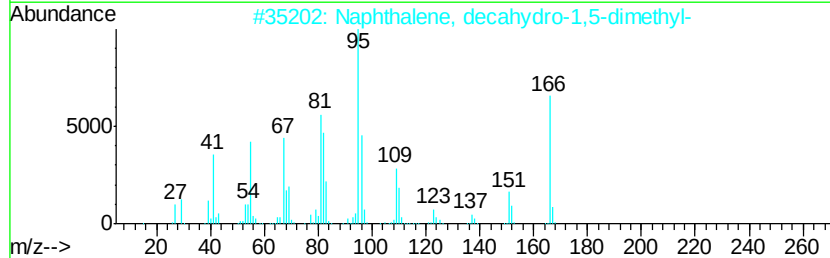
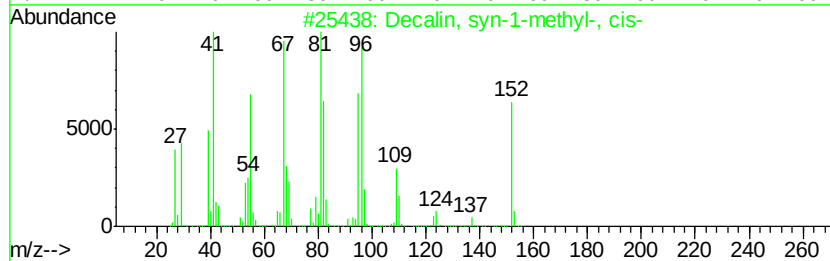
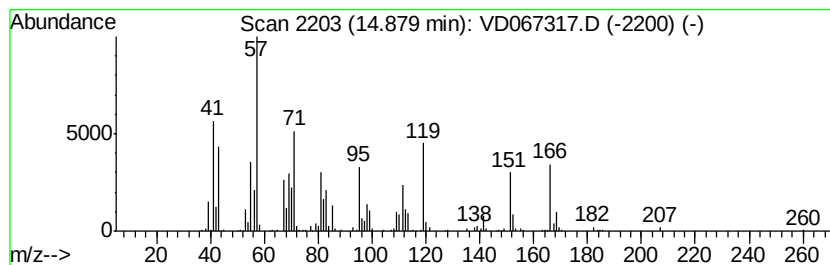
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D101320S.M
Quant Title : SW846 8260

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

Peak Number 10 unknown14.88 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.88	22.25 ug/l	834599	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decalin, syn-1-methyl-, cis-	152	C11H20	1000158-89-1	47
2		Naphthalene, decahydro-1,5-dimet...	166	C12H22	066552-62-3	38
3		Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	38
4		1-Methyldecahydronaphthalene	152	C11H20	002958-75-0	38
5		2-(1-Hydroxybut-2-enylidene)cycl...	166	C10H14O2	1000186-18-9	35



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD102020\
Data File : VD067317.D
Acq On : 20 Oct 2020 12:06
Operator : VA/SY
Sample : L4222-01
Misc : 6.99G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
EO-03-101920

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D101320S.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
Heptane, 3-ethyl-...	12.38	15.2	ug/l	422788	3	11.65	1389870	50.0
Oxalic acid, cycl...	12.96	22.9	ug/l	857900	4	13.58	1875890	50.0
Heptane, 3,3,5-tr...	13.19	15.6	ug/l	586453	4	13.58	1875890	50.0
unknown13.47	13.47	20.2	ug/l	758250	4	13.58	1875890	50.0
Naphthalene, deca...	13.90	39.5	ug/l	1483070	4	13.58	1875890	50.0
unknown14.29	14.29	15.6	ug/l	583401	4	13.58	1875890	50.0
trans-Decalin, 2-...	14.41	39.5	ug/l	1480810	4	13.58	1875890	50.0
1-Methyldecahydro...	14.57	29.0	ug/l	1089130	4	13.58	1875890	50.0
unknown14.83	14.83	14.2	ug/l	531892	4	13.58	1875890	50.0
unknown14.88	14.88	22.3	ug/l	834599	4	13.58	1875890	50.0