

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD081820\
 Data File : VD066303.D
 Acq On : 19 Aug 2020 04:56
 Operator : VA/SY
 Sample : L3689-04
 Misc : 5.47G/5.00ml/MSVOA D/SOIL
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 TP-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_D\METHOD\82D080520S.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	2.173	38	43	47	rVB3	4367	6394	1.06%	0.097%
2	4.955	509	516	526	rBV6	5209	18227	3.01%	0.276%
3	7.908	1009	1018	1024	rBV	102364	254621	42.07%	3.856%
4	7.973	1024	1029	1038	rVV2	204707	458915	75.83%	6.951%
5	8.067	1040	1045	1051	rVB6	6806	14896	2.46%	0.226%
6	8.179	1059	1064	1072	rBV5	3542	7560	1.25%	0.115%
7	8.326	1080	1089	1100	rBV	85888	190789	31.53%	2.890%
8	8.502	1112	1119	1129	rVB3	21831	52197	8.62%	0.791%
9	8.861	1169	1180	1192	rBV	245748	508898	84.09%	7.708%
10	9.355	1252	1264	1268	rBV4	15089	35783	5.91%	0.542%
11	9.408	1268	1273	1279	rVB3	6317	12351	2.04%	0.187%
12	9.773	1328	1335	1342	rVB3	19136	38723	6.40%	0.586%
13	9.908	1350	1358	1364	rBV2	21990	46282	7.65%	0.701%
14	9.973	1364	1369	1371	rBV3	3919	6471	1.07%	0.098%
15	10.114	1381	1393	1402	rVB4	9390	28017	4.63%	0.424%
16	10.261	1414	1418	1423	rVB2	8666	12294	2.03%	0.186%
17	10.337	1423	1431	1444	rBV	327029	605184	100.00%	9.166%
18	10.502	1455	1459	1467	rVB6	4930	12543	2.07%	0.190%
19	10.673	1483	1488	1492	rBV4	6440	10256	1.69%	0.155%
20	10.767	1499	1504	1505	rBV4	4829	6509	1.08%	0.099%
21	10.861	1515	1520	1524	rVB5	4607	8260	1.36%	0.125%
22	10.949	1529	1535	1541	rVB3	6553	10911	1.80%	0.165%
23	11.049	1544	1552	1556	rBV4	7876	16949	2.80%	0.257%
24	11.190	1568	1576	1580	rBV7	7221	13883	2.29%	0.210%
25	11.243	1580	1585	1590	rVB5	17671	30493	5.04%	0.462%
26	11.355	1598	1604	1608	rBV6	8229	16132	2.67%	0.244%
27	11.449	1615	1620	1627	rVB8	11836	25399	4.20%	0.385%
28	11.526	1627	1633	1636	rBV4	4708	8935	1.48%	0.135%
29	11.643	1644	1653	1663	rBV	244430	424949	70.22%	6.436%
30	11.778	1672	1676	1679	rVB5	5220	6244	1.03%	0.095%
31	11.837	1680	1686	1690	rBV6	9088	19460	3.22%	0.295%
32	11.884	1690	1694	1699	rBV4	16692	28606	4.73%	0.433%
33	11.984	1707	1711	1715	rBV5	3987	6908	1.14%	0.105%
34	12.067	1719	1725	1726	rBV5	7072	10241	1.69%	0.155%

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Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
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35	12.149	1732	1739	1748	rBV8	31615	86175	14.24%	1.305%
36	12.267	1751	1759	1764	rVB7	18272	40537	6.70%	0.614%
37	12.355	1764	1774	1776	rBV6	29313	68244	11.28%	1.034%
38	12.378	1776	1778	1785	rVB4	28372	50751	8.39%	0.769%
39	12.514	1796	1801	1808	rVB8	16120	42356	7.00%	0.642%
40	12.631	1814	1821	1828	rVB	167517	294550	48.67%	4.461%
41	12.720	1828	1836	1841	rBV8	19582	52928	8.75%	0.802%
42	12.796	1843	1849	1856	rVB6	48092	109514	18.10%	1.659%
43	12.884	1856	1864	1868	rBV7	31305	76928	12.71%	1.165%
44	12.961	1870	1877	1883	rBV7	58072	129831	21.45%	1.966%
45	13.102	1896	1901	1904	rVV4	24334	34091	5.63%	0.516%
46	13.143	1905	1908	1912	rVB5	11398	14478	2.39%	0.219%
47	13.202	1913	1918	1924	rVB8	16198	34841	5.76%	0.528%
48	13.267	1924	1929	1935	rBV3	45408	90257	14.91%	1.367%
49	13.320	1935	1938	1942	rBV6	12247	15360	2.54%	0.233%
50	13.402	1949	1952	1956	rBV5	13347	18954	3.13%	0.287%
51	13.461	1956	1962	1967	rVV7	24019	67093	11.09%	1.016%
52	13.502	1967	1969	1975	rVB7	19621	28488	4.71%	0.431%
53	13.578	1975	1982	1991	rBV2	128848	267009	44.12%	4.044%
54	13.720	2001	2006	2011	rBV8	19336	41373	6.84%	0.627%
55	13.902	2030	2037	2041	rBV4	114964	221431	36.59%	3.354%
56	14.008	2051	2055	2058	rBV6	28969	51482	8.51%	0.780%
57	14.072	2064	2066	2067	rBV2	11594	8595	1.42%	0.130%
58	14.120	2070	2074	2081	rVB2	74231	158877	26.25%	2.406%
59	14.184	2081	2085	2087	rBV5	11404	16185	2.67%	0.245%
60	14.231	2088	2093	2097	rVB5	45135	69377	11.46%	1.051%
61	14.296	2098	2104	2109	rVB9	36156	75481	12.47%	1.143%
62	14.355	2109	2114	2118	rBV7	25590	50593	8.36%	0.766%
63	14.408	2118	2123	2126	rBV3	72665	122112	20.18%	1.849%
64	14.443	2127	2129	2133	rVB	44786	53696	8.87%	0.813%
65	14.531	2139	2144	2145	rBV5	23566	32411	5.36%	0.491%
66	14.572	2148	2151	2157	rVB5	59134	98097	16.21%	1.486%
67	14.690	2166	2171	2174	rBV6	26197	46869	7.74%	0.710%
68	14.767	2180	2184	2188	rVB6	34586	58509	9.67%	0.886%
69	14.825	2188	2194	2201	rBV4	107571	258506	42.72%	3.915%
70	14.890	2201	2205	2209	rVB5	40178	65661	10.85%	0.994%
71	15.002	2220	2224	2226	rBV5	23765	31085	5.14%	0.471%

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Integration Parameters: RTEINT.P

Integrator: RTE
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Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
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72	15.096	2236	2240	2244	rBV7	51232	78752	13.01%	1.193%
73	15.214	2255	2260	2268	rVB7	59911	136303	22.52%	2.064%
74	15.449	2298	2300	2305	rVB6	26163	30381	5.02%	0.460%
75	15.549	2315	2317	2324	rVB8	51275	100157	16.55%	1.517%
76	16.225	2425	2432	2438	rBV10	37981	105699	17.47%	1.601%
77	16.702	2508	2513	2527	rVB8	71095	214201	35.39%	3.244%

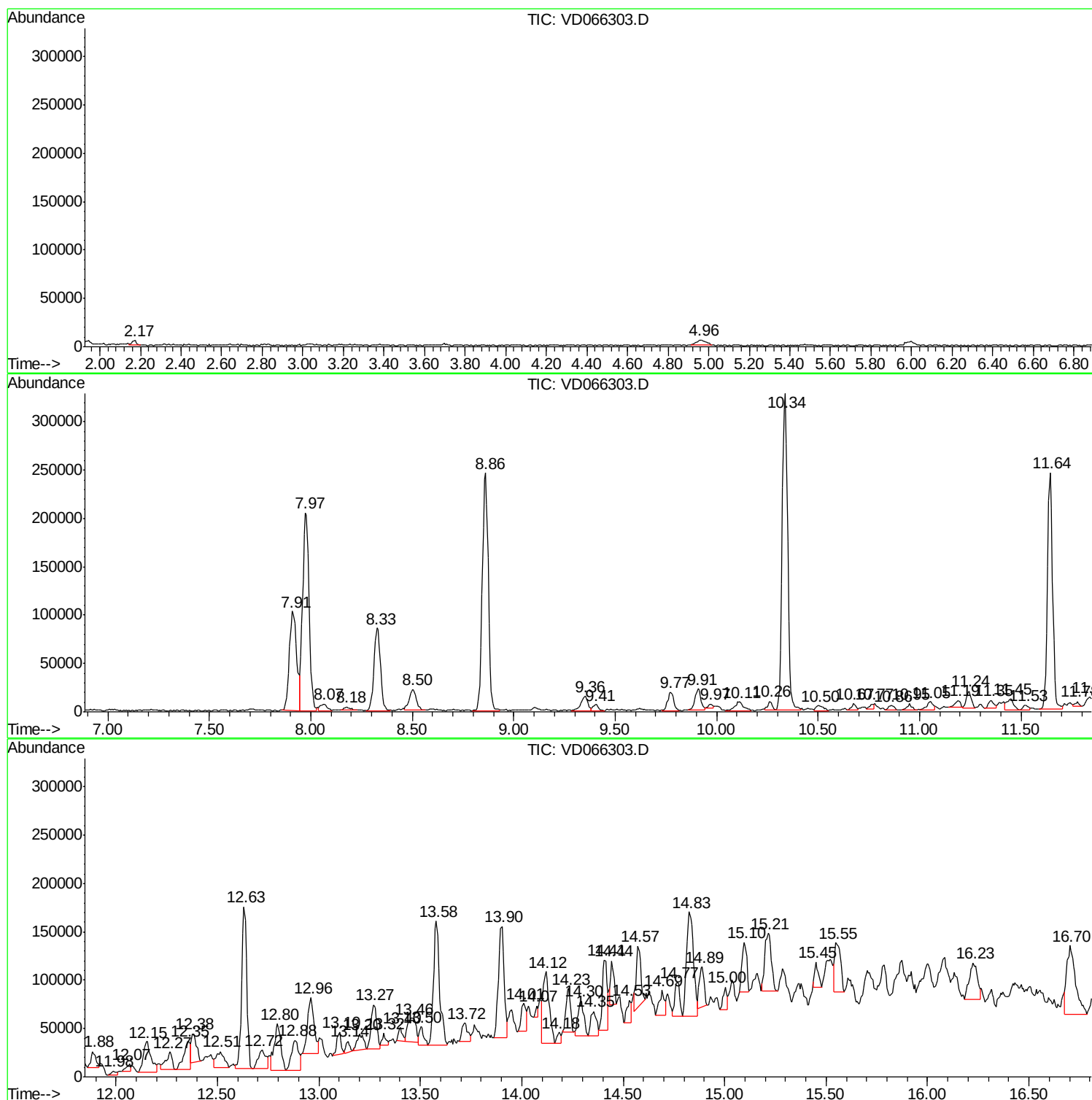
Sum of corrected areas: 6602498

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Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080520S.M
Quant Title : SW846 8260

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



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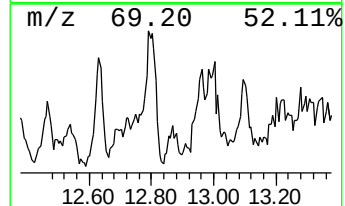
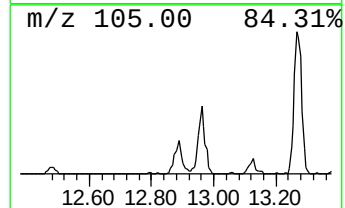
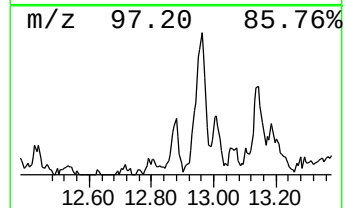
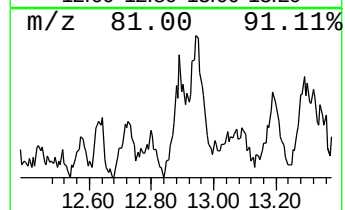
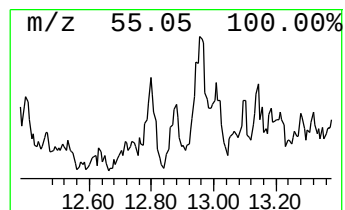
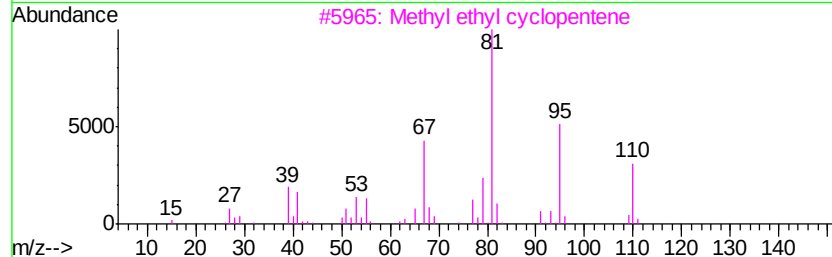
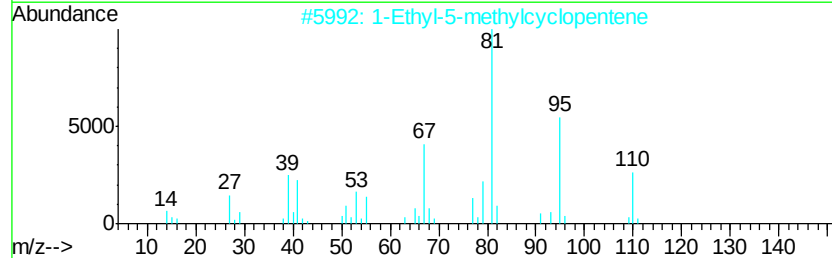
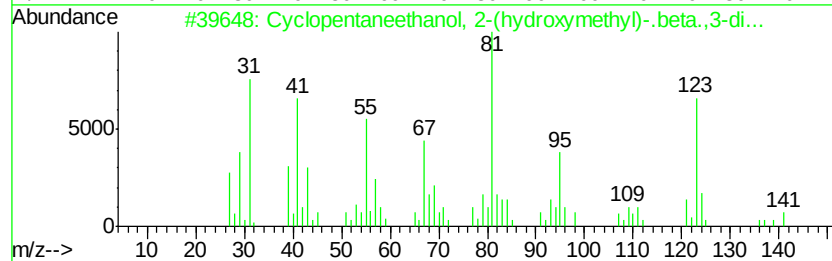
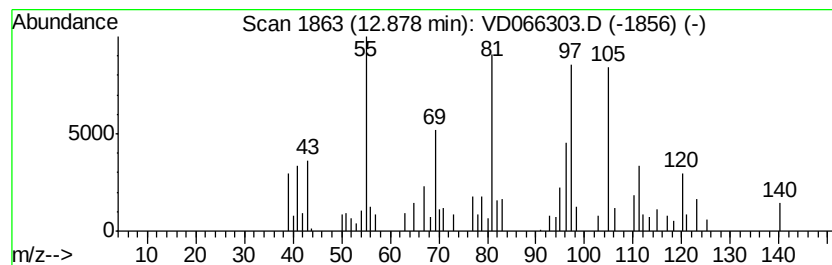
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080520S.M
Quant Title : SW846 8260

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

Peak Number 1 unknown12.88 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.88	14.41 ug/l	76928	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentaneethanol, 2-(hydroxym...	172	C10H20O2	000485-42-7	43
2		1-Ethyl-5-methylcyclopentene	110	C8H14	097797-57-4	43
3		Methyl ethyl cyclopentene	110	C8H14	019780-56-4	43
4		Cyclohexene, 4-methyl-	96	C7H12	000591-47-9	43
5		Cyclohexanepropanal	140	C9H16O	004361-28-8	38



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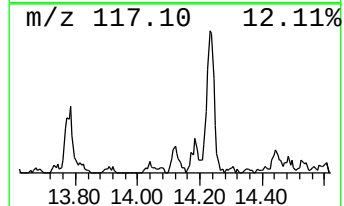
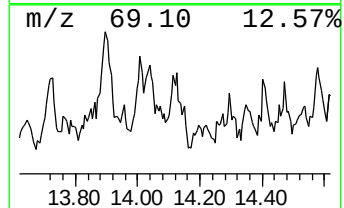
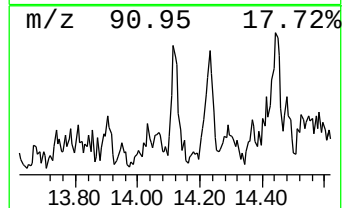
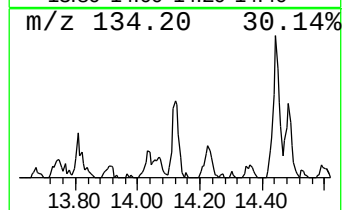
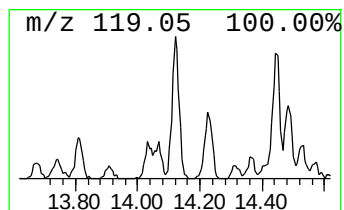
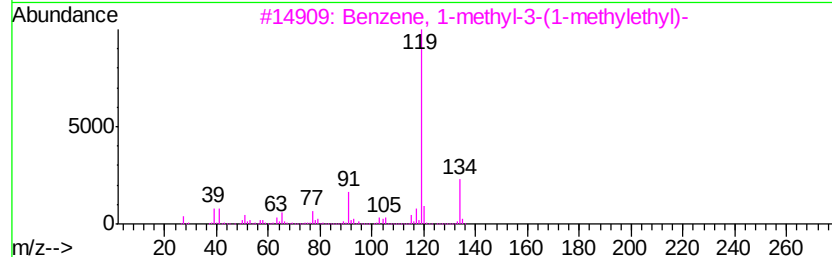
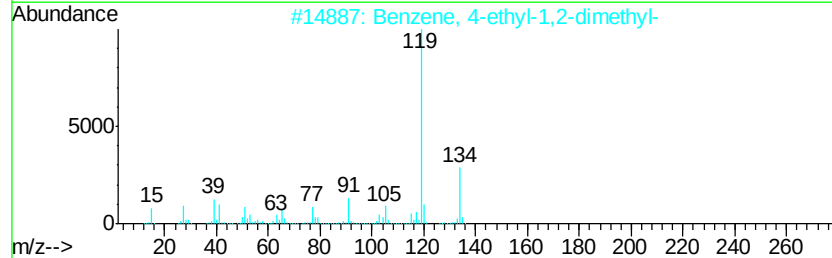
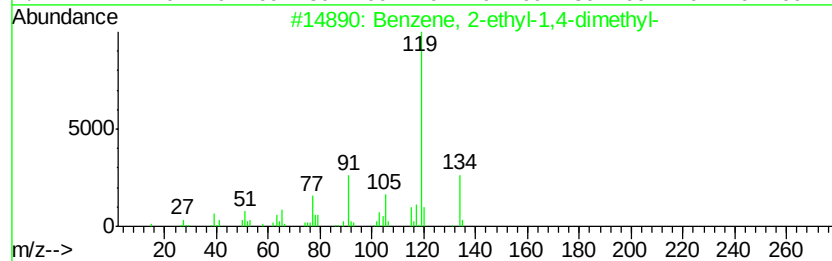
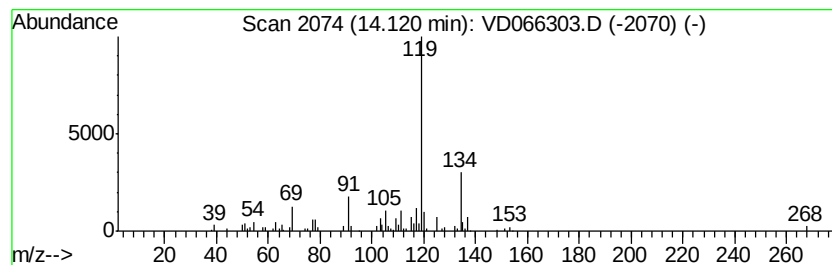
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.12	29.75 ug/l	158877	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93
3		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	93
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5		p-Cymene	134	C10H14	000099-87-6	90



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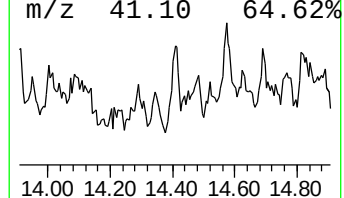
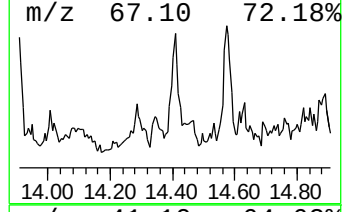
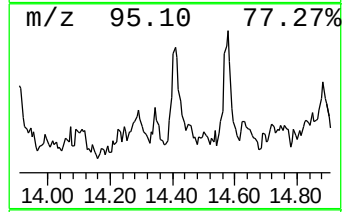
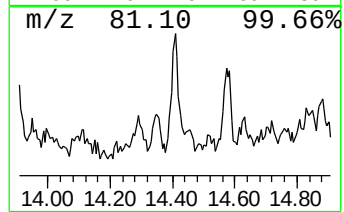
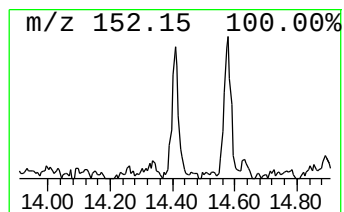
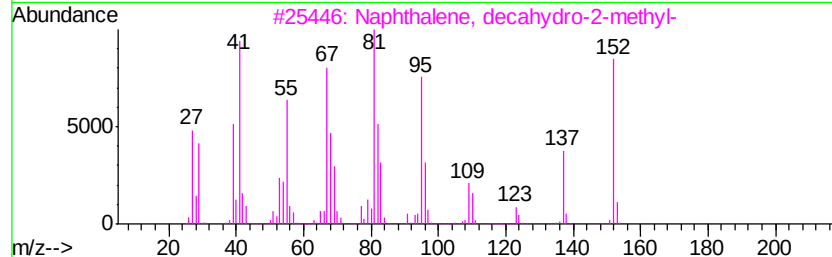
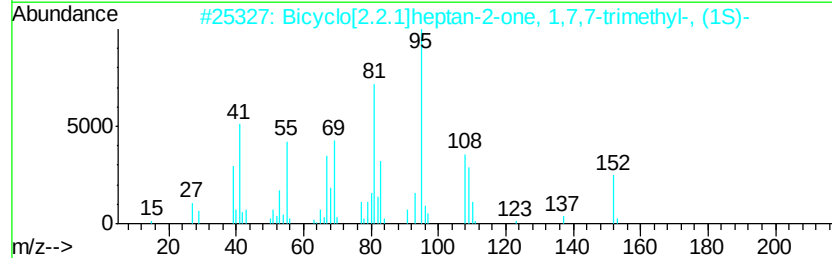
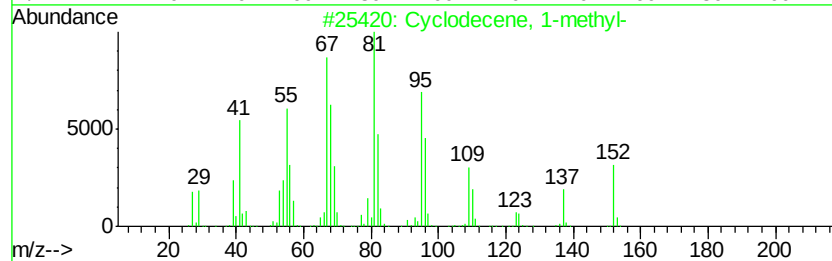
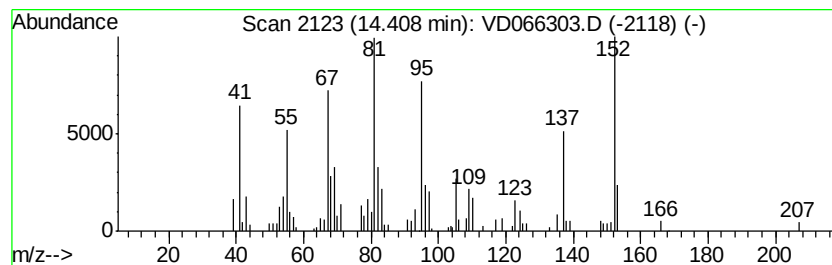
Instrument :
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ClientSampled :
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Quant Title : SW846 8260

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

Peak Number 3 Cyclodecene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.41	22.87 ug/l	122112	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclodecene, 1-methyl-	152 C11H20	066633-38-3 81
2		Bicyclo[2.2.1]heptan-2-one, 1,7,...	152 C10H16O	000464-48-2 64
3		Naphthalene, decahydro-2-methyl-	152 C11H20	002958-76-1 64
4		Decalin, anti-1-methyl-, cis-	152 C11H20	1000158-89-0 62
5		Decalin, syn-1-methyl-, cis-	152 C11H20	1000158-89-1 62



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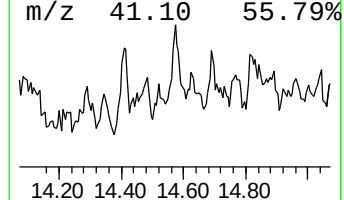
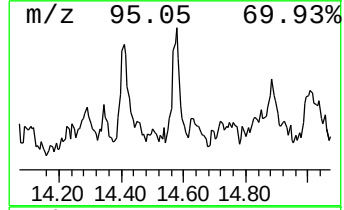
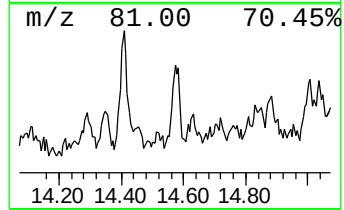
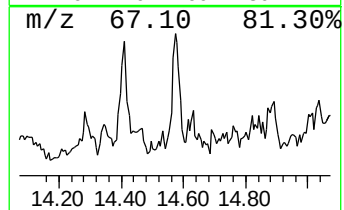
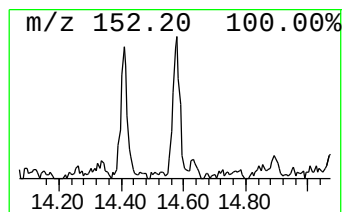
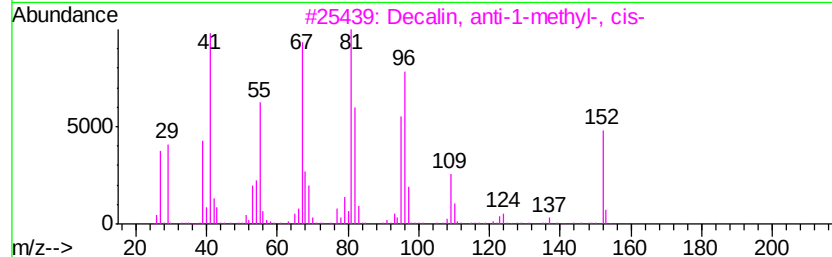
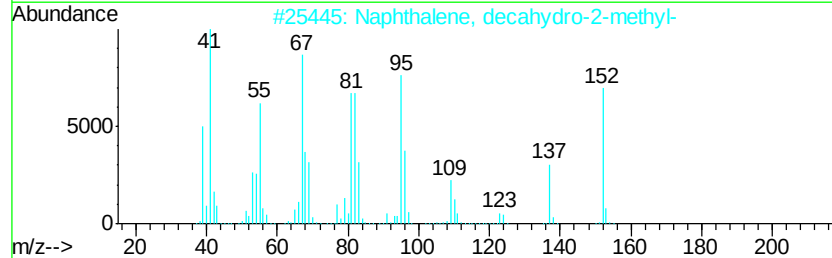
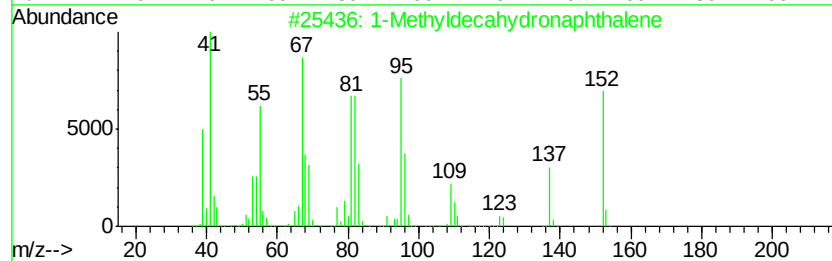
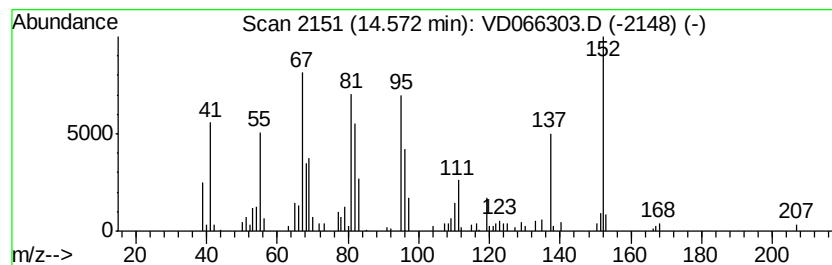
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Quant Title : SW846 8260

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

Peak Number 4 1-Methyldecahydronaphthalene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	18.37 ug/l	98097	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Methyldecahydronaphthalene	152	C11H20	002958-75-0	93
2		Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	93
3		Decalin, anti-1-methyl-, cis-	152	C11H20	1000158-89-0	64
4		Decalin, syn-1-methyl-, cis-	152	C11H20	1000158-89-1	55
5		Bicyclo[5.4.0]undecane (trans)	152	C11H20	021394-30-9	52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD081820\
Data File : VD066303.D
Acq On : 19 Aug 2020 04:56
Operator : VA/SY
Sample : L3689-04
Misc : 5.47G/5.00ml/MSVOA D/SOIL
ALS Vial : 39 Sample Multiplier: 1

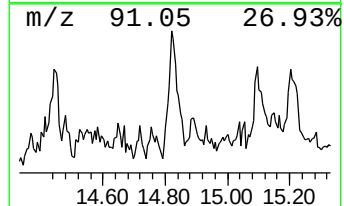
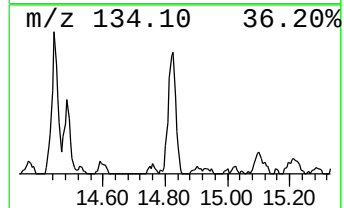
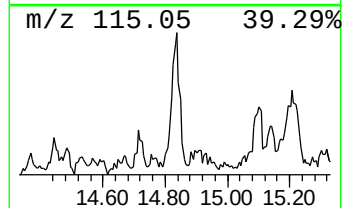
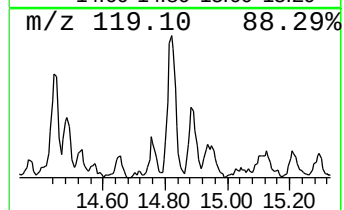
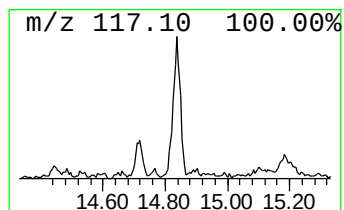
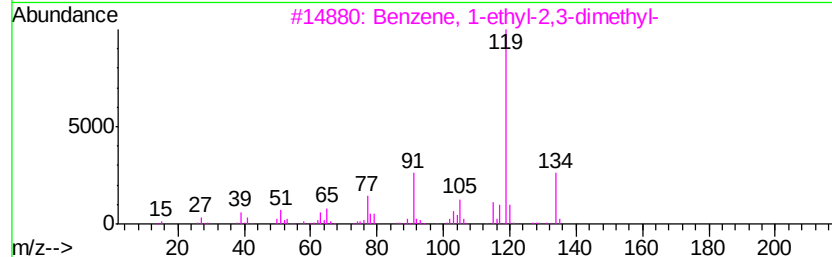
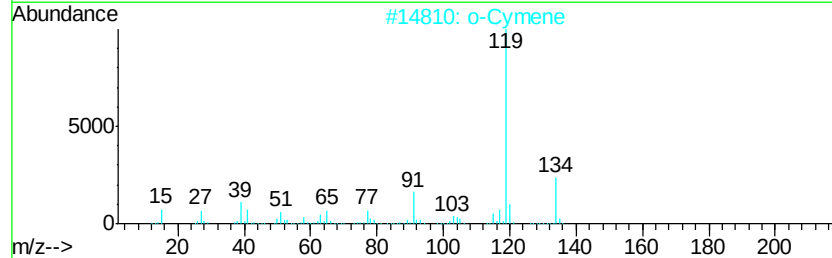
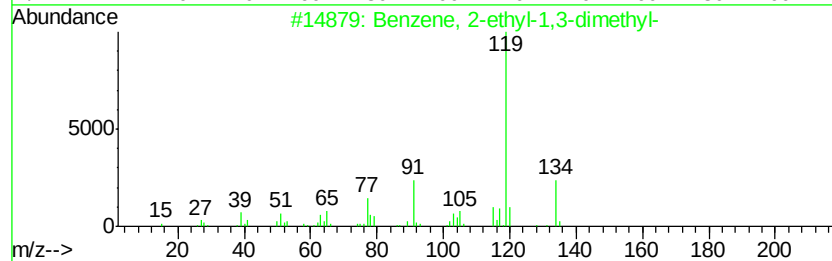
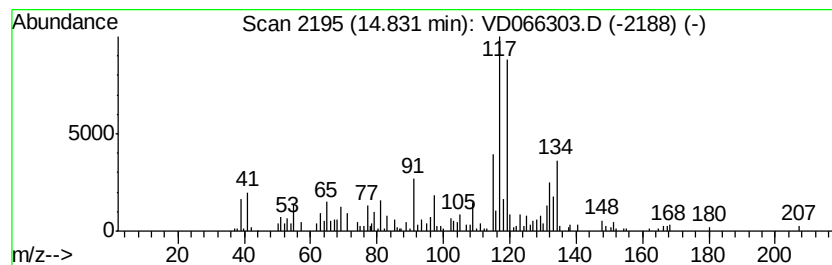
Instrument :
MSVOA_D
ClientSampleId :
TP-5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080520S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	48.41 ug/l	258506	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4 55
2		o-Cymene	134 C10H14	000527-84-4 55
3		Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2 55
4		Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3 55
5		1,3-Cyclopentadiene, 1,2,3,4-tet...	134 C10H14	076089-59-3 49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD081820\
Data File : VD066303.D
Acq On : 19 Aug 2020 04:56
Operator : VA/SY
Sample : L3689-04
Misc : 5.47G/5.00ml/MSVOA D/SOIL
ALS Vial : 39 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampled :
TP-5

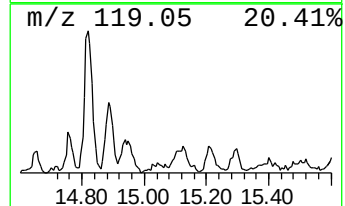
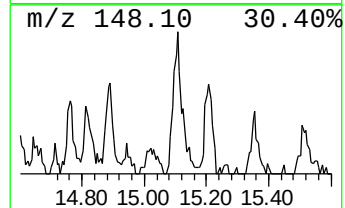
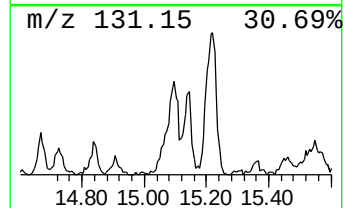
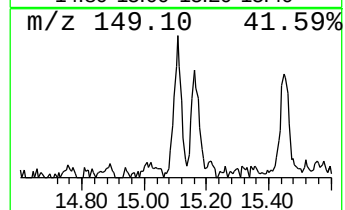
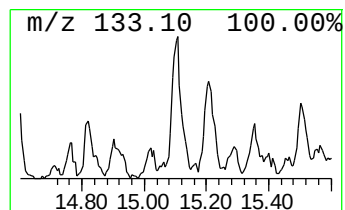
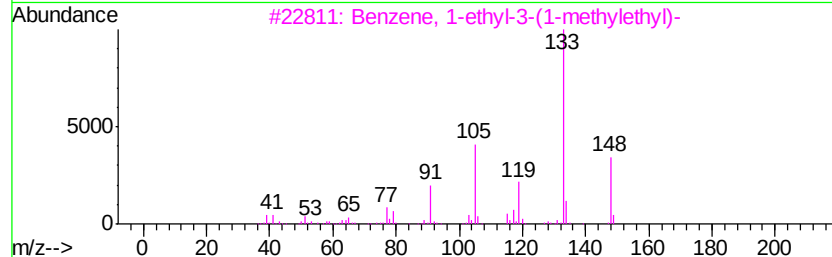
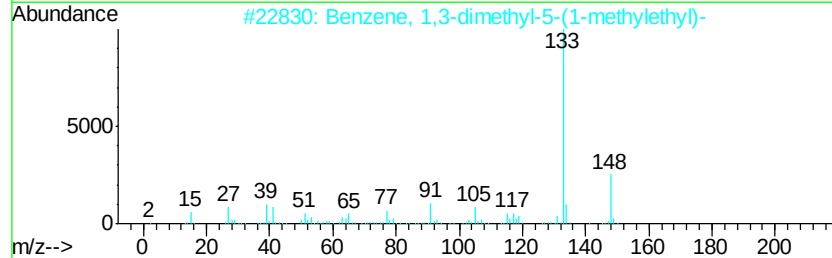
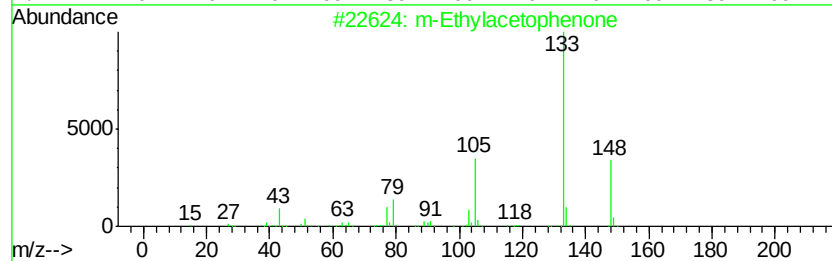
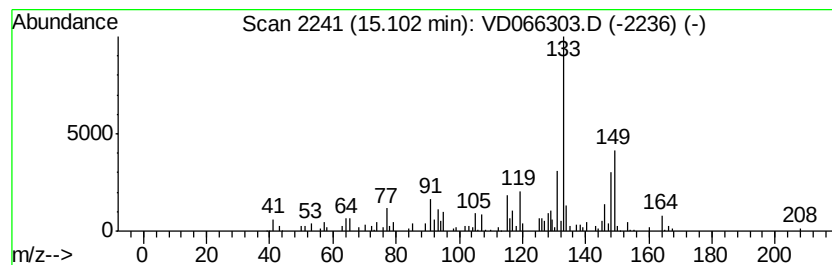
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Quant Title : SW846 8260

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

Peak Number 6 unknown15.10 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.10	14.75 ug/l	78752	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	m-Ethylacetophenone	148	C10H12O	022699-70-3	43
2		Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	43
3		Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	43
4		5-Aminoindazole	133	C7H7N3	019335-11-6	38
5		2-tert-Butyltoluene	148	C11H16	001074-92-6	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD081820\
Data File : VD066303.D
Acq On : 19 Aug 2020 04:56
Operator : VA/SY
Sample : L3689-04
Misc : 5.47G/5.00ml/MSVOA D/SOIL
ALS Vial : 39 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
TP-5

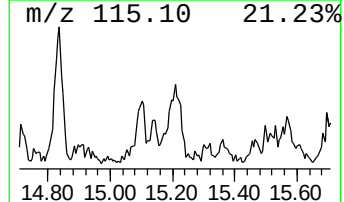
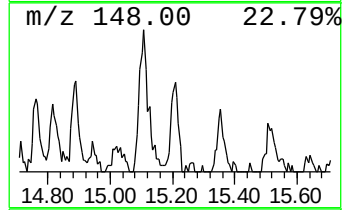
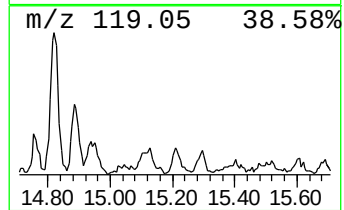
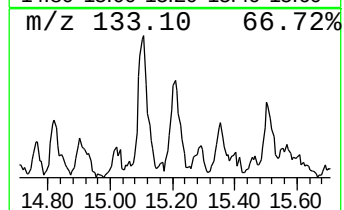
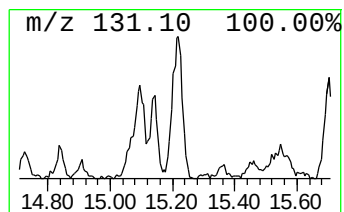
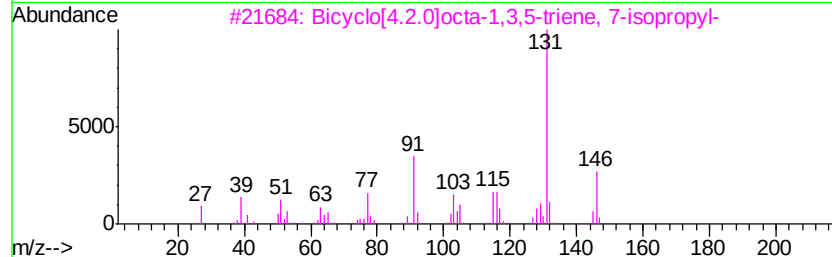
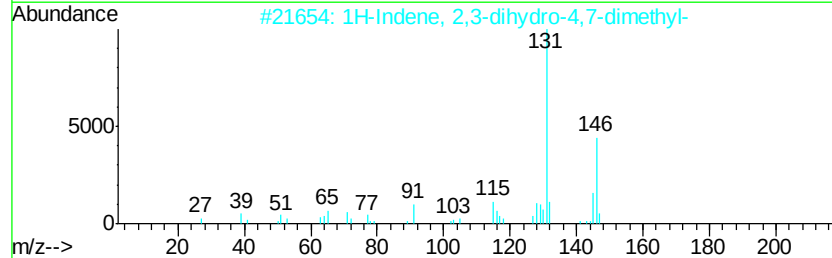
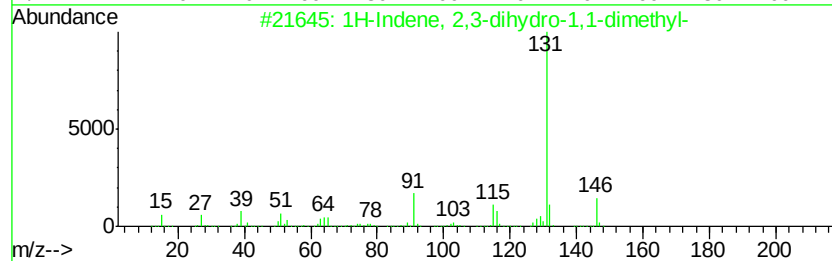
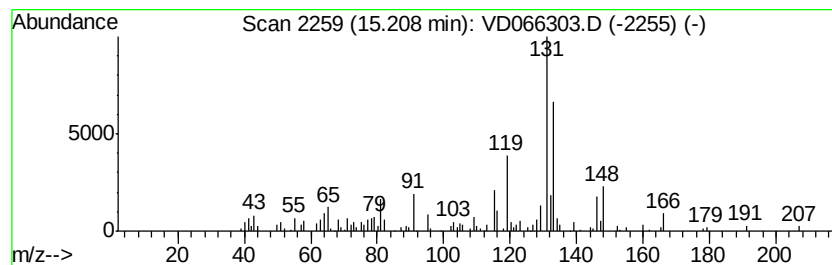
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Quant Title : SW846 8260

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

Peak Number 7 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.21	25.52 ug/l	136303	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	60
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	49
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	49
4		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	46
5		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD081820\
Data File : VD066303.D
Acq On : 19 Aug 2020 04:56
Operator : VA/SY
Sample : L3689-04
Misc : 5.47G/5.00ml/MSVOA D/SOIL
ALS Vial : 39 Sample Multiplier: 1

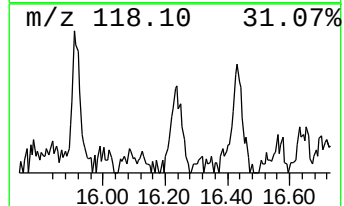
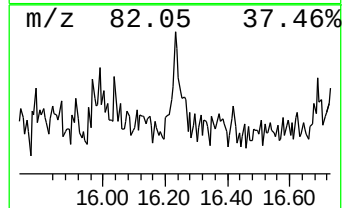
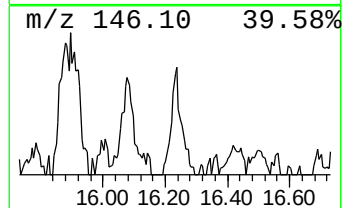
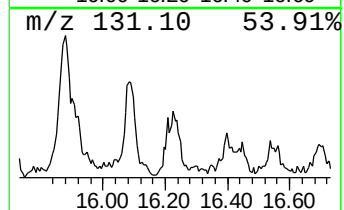
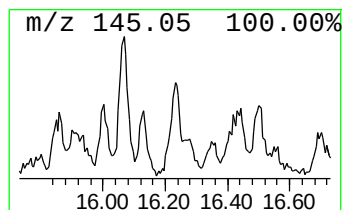
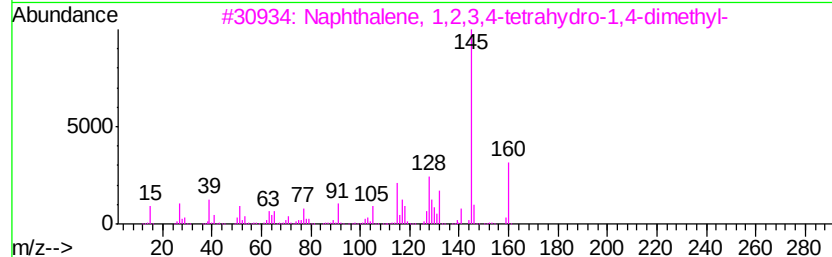
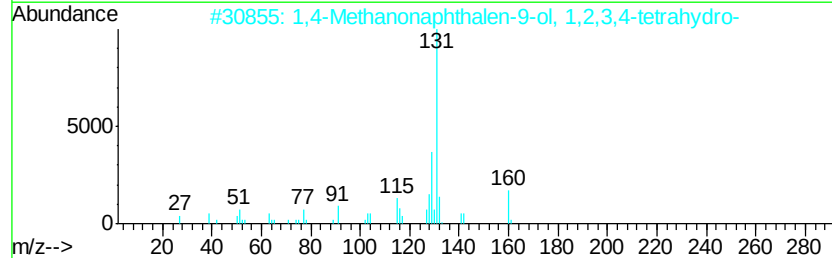
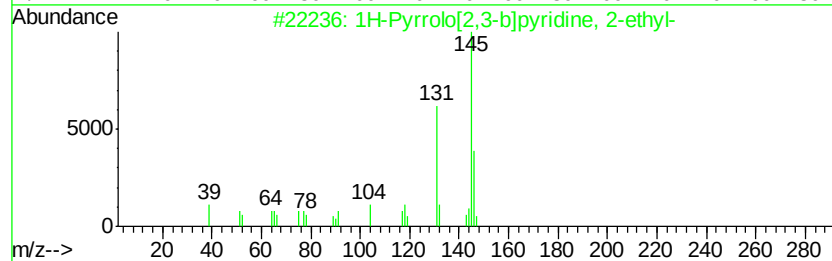
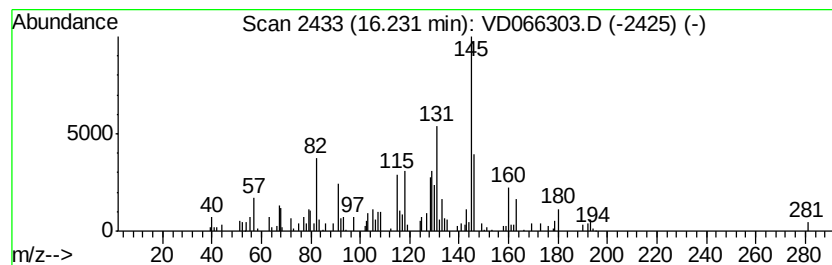
Instrument :
MSVOA_D
ClientSampleId :
TP-5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080520S.M
Quant Title : SW846 8260

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

Peak Number 9 1H-Pyrrolo[2,3-b]pyridine, ... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.23	19.79 ug/l	105699	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Pyrrolo[2,3-b]pyridine, 2-ethyl-	146 C9H10N2	023612-49-9 53
2		1,4-Methanonaphthalen-9-ol, 1,2,...	160 C11H12O	055255-94-2 46
3		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	004175-54-6 41
4		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 38
5		Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3 38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD081820\
Data File : VD066303.D
Acq On : 19 Aug 2020 04:56
Operator : VA/SY
Sample : L3689-04
Misc : 5.47G/5.00ml/MSVOA D/SOIL
ALS Vial : 39 Sample Multiplier: 1

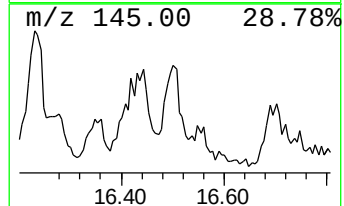
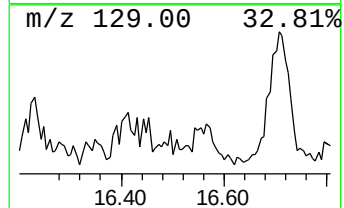
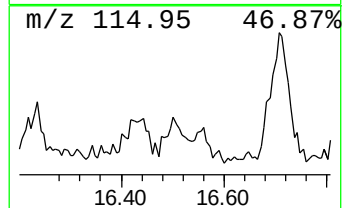
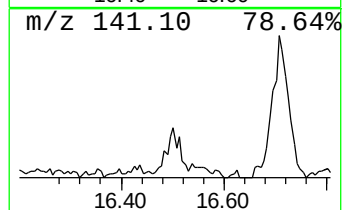
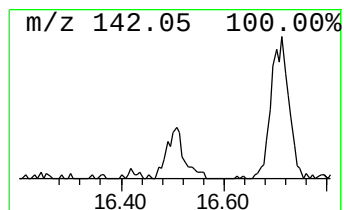
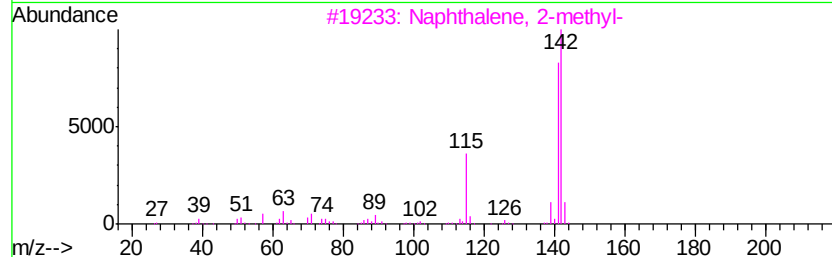
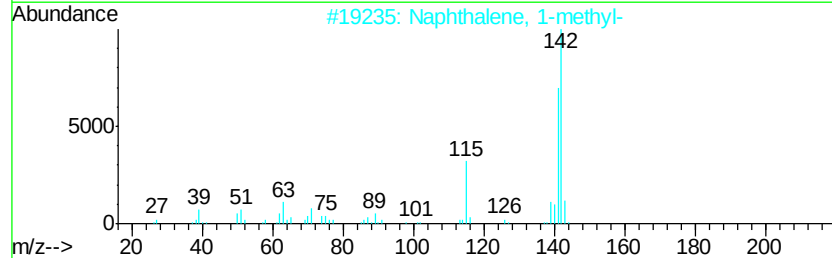
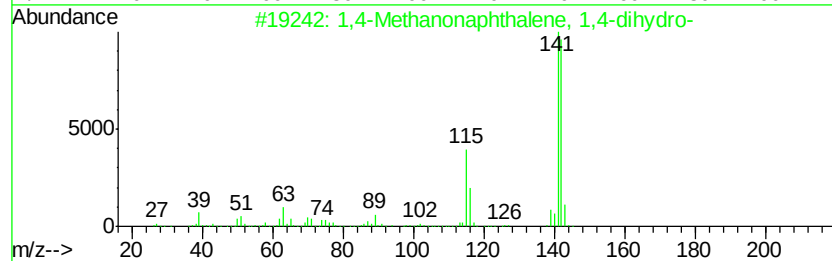
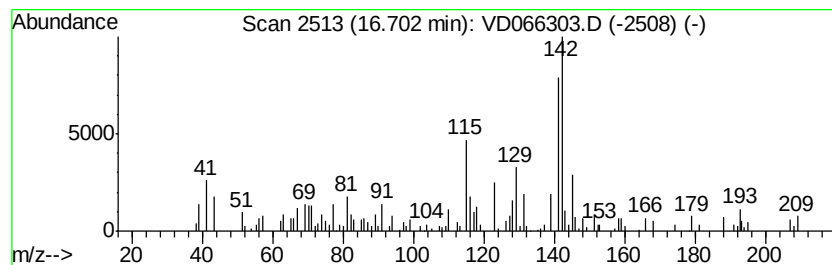
Instrument :
MSVOA_D
ClientSampleId :
TP-5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080520S.M
Quant Title : SW846 8260

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

Peak Number 10 1,4-Methanonaphthalene, 1,4... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.70	40.11 ug/l	214201	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1,4-Methanonaphthalene, 1,4-dihy...	142 C11H10	004453-90-1	64
2	Naphthalene, 1-methyl-	142 C11H10	000090-12-0	64
3	Naphthalene, 2-methyl-	142 C11H10	000091-57-6	58
4	Benzocycloheptatriene	142 C11H10	000264-09-5	52
5	Spiro(tricyclo[6.2.1.0(2,7)]unde...	186 C13H14O	1000210-02-3	47



Data Path : Z:\V0ASRV\HPCHEM1\MSV0A_D\DATA\VD081820\
Data File : VD066303.D
Acq On : 19 Aug 2020 04:56
Operator : VA/SY
Sample : L3689-04
Misc : 5.47G/5.00ml/MSV0A_D/SOIL
ALS Vial : 39 Sample Multiplier: 1

Instrument :
MSV0A_D
ClientSampleId :
TP-5

Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_D\METHOD\82D080520S.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Benzene, 2-ethyl-...	14.12	29.8	ug/l	158877	4	13.58	267009	50.0
Cyclodecene, 1-me...	14.41	22.9	ug/l	122112	4	13.58	267009	50.0
1-Methyldecahydro...	14.57	18.4	ug/l	98097	4	13.58	267009	50.0
Benzene, 2-ethyl-...	14.83	48.4	ug/l	258506	4	13.58	267009	50.0
unknown15.10	15.10	14.8	ug/l	78752	4	13.58	267009	50.0
1H-Indene, 2,3-di...	15.21	25.5	ug/l	136303	4	13.58	267009	50.0
1H-Pyrrolo[2,3-b]...	16.23	19.8	ug/l	105699	4	13.58	267009	50.0
1,4-Methanonaphth...	16.70	40.1	ug/l	214201	4	13.58	267009	50.0