

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD090920\
 Data File : VD066603.D
 Acq On : 10 Sep 2020 02:11
 Operator : VA/SY
 Sample : L3870-01
 Misc : 6.62G/5.00ml/MSVOA D/SOIL
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 SU-04-090820

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\82D090320S.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	4.967	504	518	529	rBV2	75983	239196	10.33%	1.209%
2	7.914	1006	1019	1024	rBV2	342266	813789	35.14%	4.113%
3	7.979	1024	1030	1041	rVB	508824	1130260	48.81%	5.713%
4	8.326	1078	1089	1100	rBV	291015	659803	28.49%	3.335%
5	8.861	1171	1180	1197	rBV	838201	1682248	72.65%	8.503%
6	10.338	1422	1431	1439	rBV	1309993	2315592	100.00%	11.704%
7	11.644	1645	1653	1665	rBV	1182703	2015579	87.04%	10.188%
8	11.849	1680	1688	1693	rBV3	34467	67212	2.90%	0.340%
9	12.179	1732	1744	1750	rBV5	33771	83068	3.59%	0.420%
10	12.632	1814	1821	1831	rVB	898462	1471802	63.56%	7.439%
11	12.814	1845	1852	1857	rVB3	28179	55998	2.42%	0.283%
12	12.885	1857	1864	1867	rVV	66530	121569	5.25%	0.614%
13	12.955	1871	1876	1882	rVV4	116011	225966	9.76%	1.142%
14	13.008	1882	1885	1891	rVB6	15635	23662	1.02%	0.120%
15	13.120	1896	1904	1910	rBV	63278	129929	5.61%	0.657%
16	13.185	1911	1915	1922	rVB7	28535	52710	2.28%	0.266%
17	13.267	1924	1929	1935	rVV	91568	147355	6.36%	0.745%
18	13.402	1944	1952	1956	rVB3	33000	58482	2.53%	0.296%
19	13.461	1956	1962	1967	rBV5	59255	119130	5.14%	0.602%
20	13.514	1967	1971	1973	rVV5	34298	50266	2.17%	0.254%
21	13.579	1975	1982	1991	rVV2	1181789	2097177	90.57%	10.600%
22	13.738	2001	2009	2011	rBV7	29881	50174	2.17%	0.254%
23	13.773	2011	2015	2019	rVV2	118007	200274	8.65%	1.012%
24	13.814	2019	2022	2031	rVB4	88158	200362	8.65%	1.013%
25	13.896	2031	2036	2042	rBV3	229567	372988	16.11%	1.885%
26	13.973	2045	2049	2053	rVB	60434	86388	3.73%	0.437%
27	14.061	2061	2064	2068	rVB2	59033	81695	3.53%	0.413%
28	14.120	2069	2074	2078	rVB	105224	161277	6.96%	0.815%
29	14.226	2087	2092	2098	rVB2	141383	242134	10.46%	1.224%
30	14.296	2098	2104	2105	rBV4	26915	50700	2.19%	0.256%
31	14.367	2110	2116	2119	rBV5	59353	128939	5.57%	0.652%
32	14.408	2119	2123	2127	rVV4	90914	150418	6.50%	0.760%
33	14.449	2127	2130	2133	rVV4	50496	65311	2.82%	0.330%
34	14.479	2133	2135	2139	rVB	62187	78984	3.41%	0.399%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	14.526	2139	2143	2147	rBV2	77520	111491	4.81%	0.564%
36	14.573	2147	2151	2157	rVB5	65698	106122	4.58%	0.536%
37	14.632	2157	2161	2164	rBV5	35840	64298	2.78%	0.325%
38	14.667	2164	2167	2171	rVB3	50395	73594	3.18%	0.372%
39	14.720	2171	2176	2177	rBV4	65886	96103	4.15%	0.486%
40	14.755	2177	2182	2188	rVB4	198164	329753	14.24%	1.667%
41	14.832	2188	2195	2200	rBV4	239420	596310	25.75%	3.014%
42	14.985	2217	2221	2228	rVB2	167644	280729	12.12%	1.419%
43	15.096	2235	2240	2244	rBV4	82886	135144	5.84%	0.683%
44	15.208	2253	2259	2268	rBV6	95207	260041	11.23%	1.314%
45	15.285	2268	2272	2280	rVB9	41923	105077	4.54%	0.531%
46	15.367	2281	2286	2296	rBV6	78782	169151	7.30%	0.855%
47	15.461	2296	2302	2307	rBV2	106904	199508	8.62%	1.008%
48	15.520	2307	2312	2313	rBV4	60454	87051	3.76%	0.440%
49	15.596	2322	2325	2336	rVB4	99745	195179	8.43%	0.987%
50	15.708	2336	2344	2350	rBV6	69432	177796	7.68%	0.899%
51	15.779	2353	2356	2362	rVB6	34299	55003	2.38%	0.278%
52	15.873	2366	2372	2374	rVV5	53170	116990	5.05%	0.591%
53	15.914	2374	2379	2388	rVB4	144083	300705	12.99%	1.520%
54	16.002	2392	2394	2400	rVB7	20064	29802	1.29%	0.151%
55	16.079	2400	2407	2414	rBV6	44105	125297	5.41%	0.633%
56	16.237	2425	2434	2441	rBV3	129313	332008	14.34%	1.678%
57	16.355	2449	2454	2458	rBV7	19002	30511	1.32%	0.154%
58	16.437	2461	2468	2474	rVB3	106084	225777	9.75%	1.141%
59	16.502	2474	2479	2484	rBV5	32848	75107	3.24%	0.380%
60	16.708	2507	2514	2516	rBV7	34354	75278	3.25%	0.380%

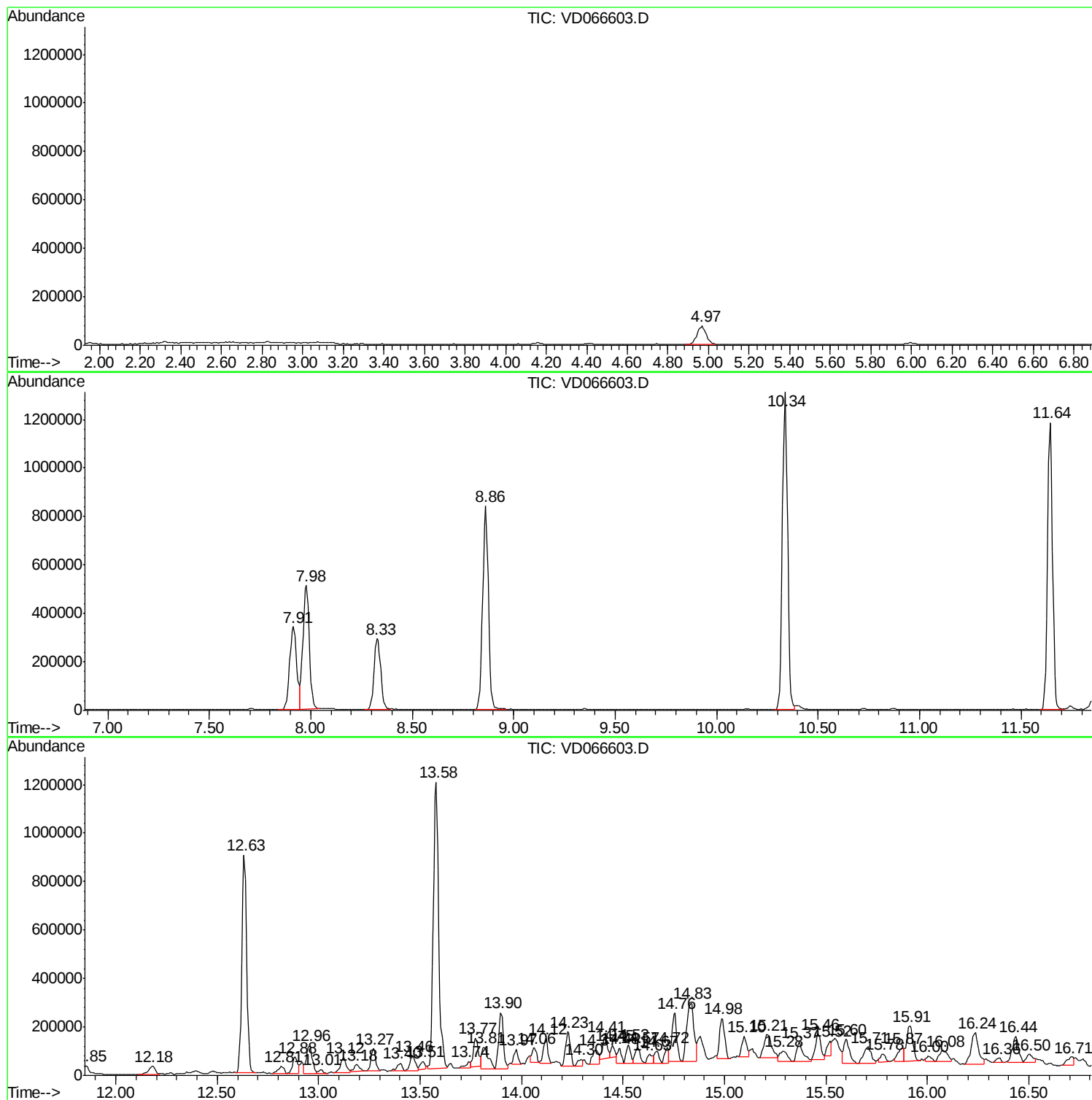
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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ClientSampleId :
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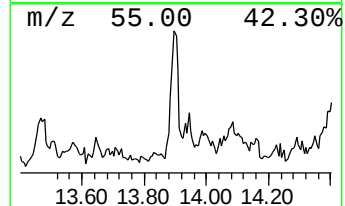
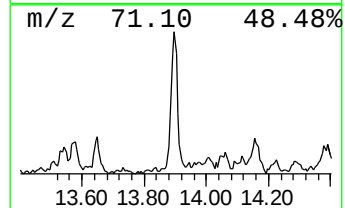
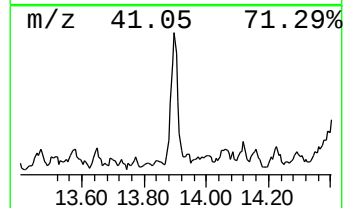
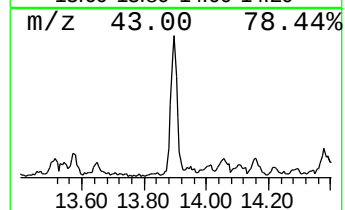
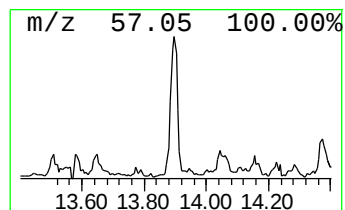
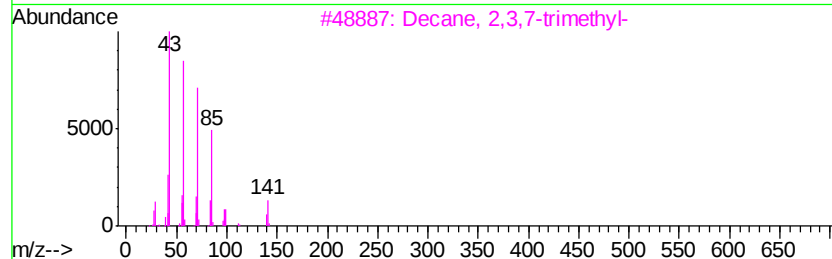
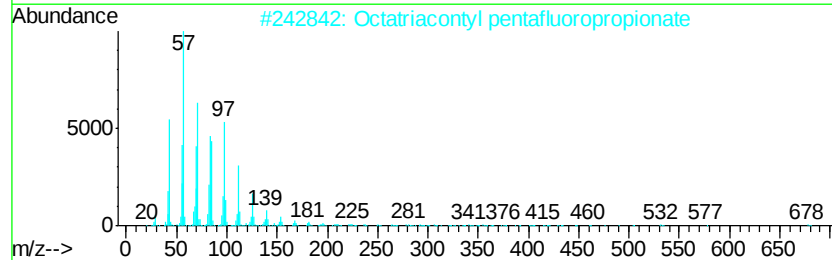
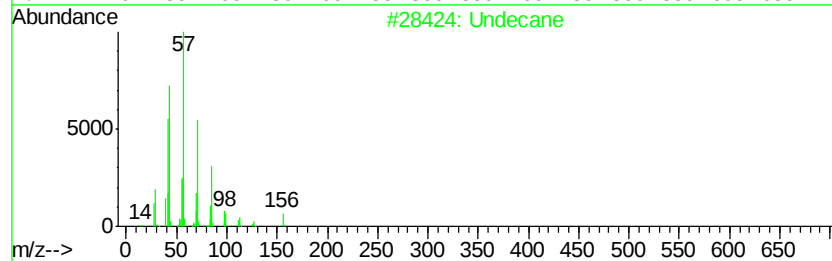
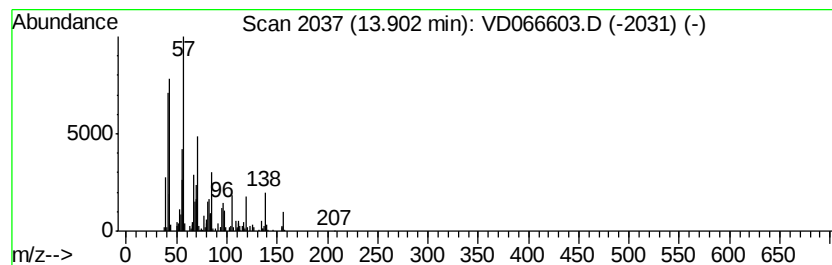
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 1 Undecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	8.89 ug/l	372988	1,4-Dichlorobenzene-d4	13.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane		156	C11H24	001120-21-4	60
2	Octatriacontyl pentafluoropropio...		697	C41H77F5O2	1000351-89-1	47
3	Decane, 2,3,7-trimethyl-		184	C13H28	062238-13-5	43
4	Tridecane		184	C13H28	000629-50-5	43
5	Dodecane		170	C12H26	000112-40-3	43



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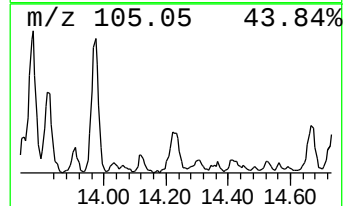
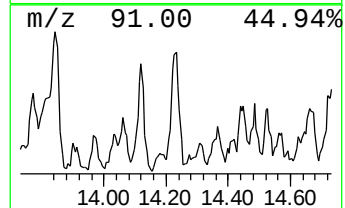
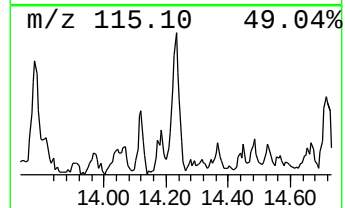
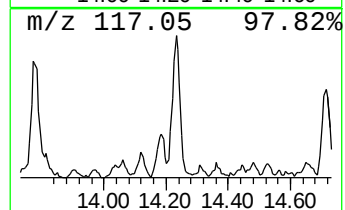
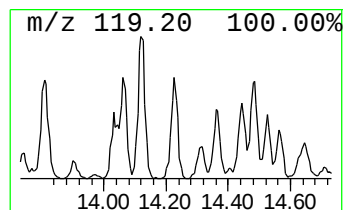
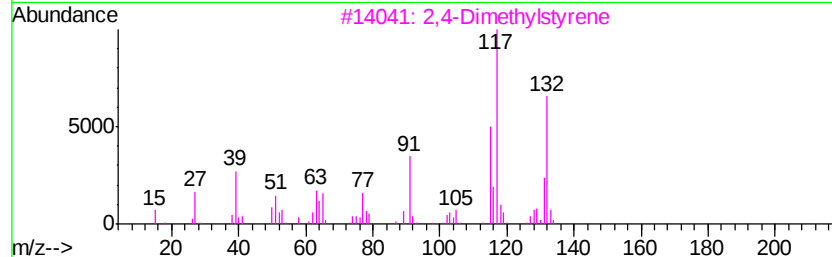
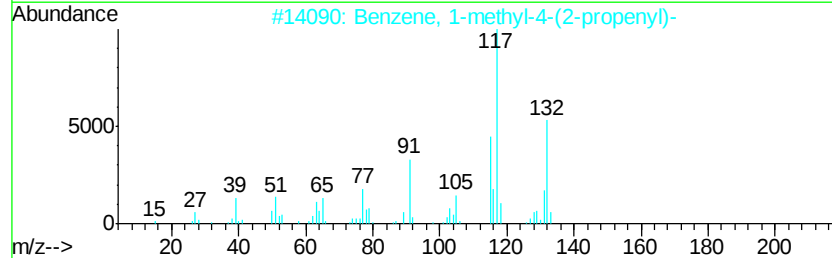
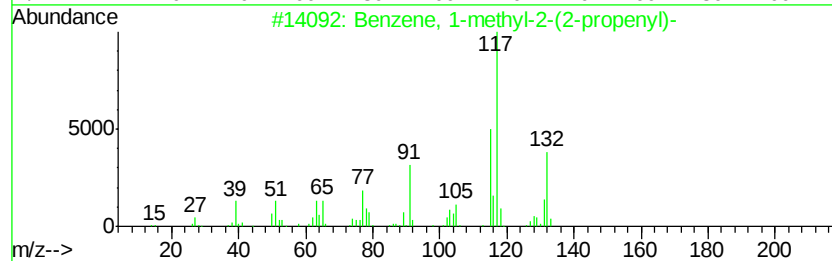
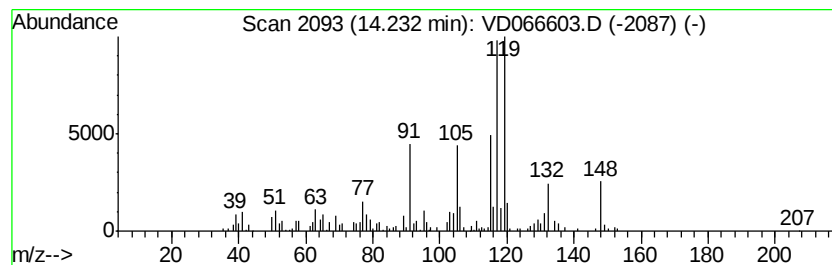
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 2 unknown14.23 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.23	5.77 ug/l	242134	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	46
2		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	46
3		2,4-Dimethylstyrene	132	C10H12	002234-20-0	42
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	38
5		Propiophenone, 2'-methyl-	148	C10H12O	002040-14-4	38



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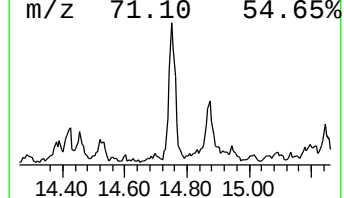
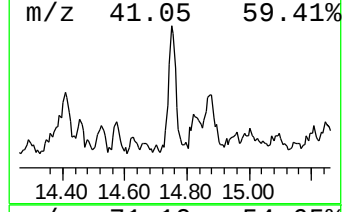
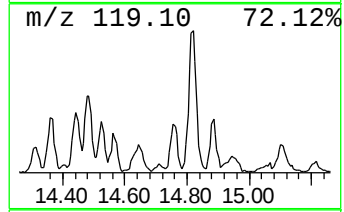
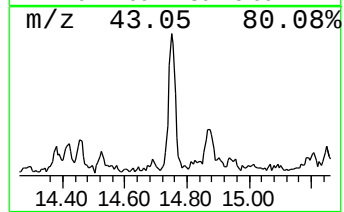
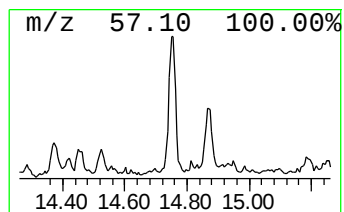
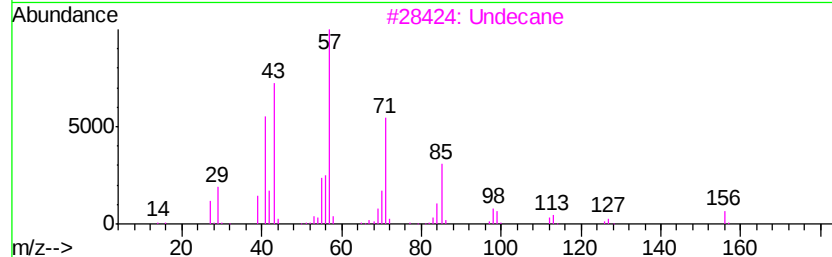
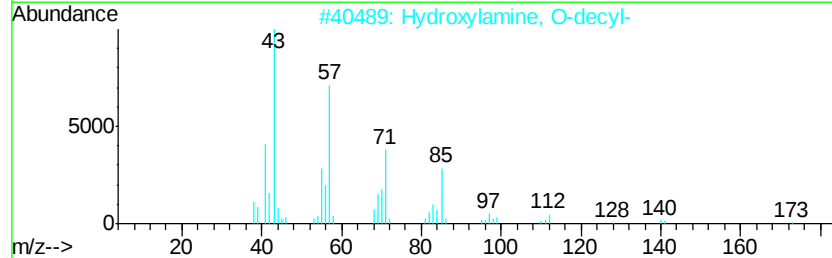
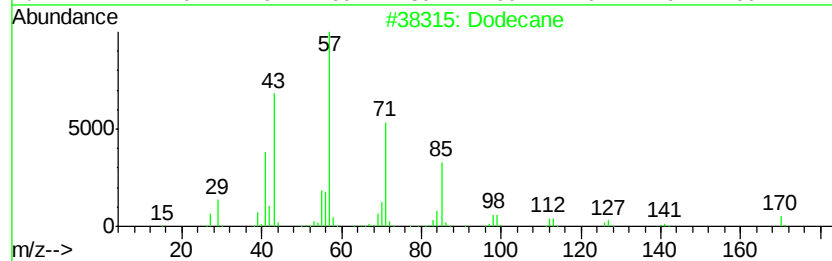
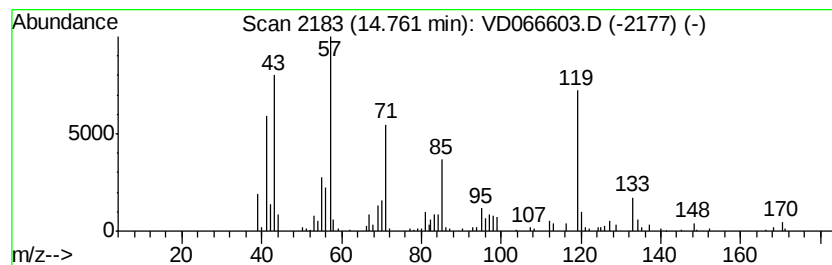
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Dodecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	7.86 ug/l	329753	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane	170	C12H26	000112-40-3	60
2		Hydroxylamine, O-decyl-	173	C10H23NO	029812-79-1	52
3		Undecane	156	C11H24	001120-21-4	50
4		Octane	114	C8H18	000111-65-9	49
5		Decane	142	C10H22	000124-18-5	47



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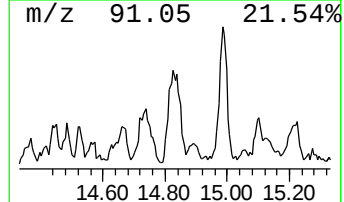
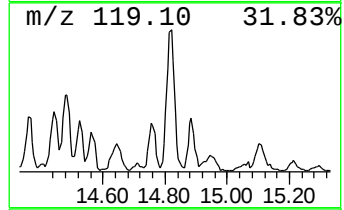
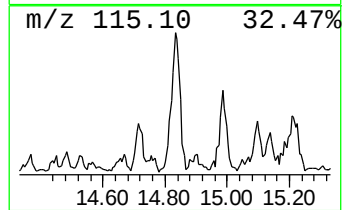
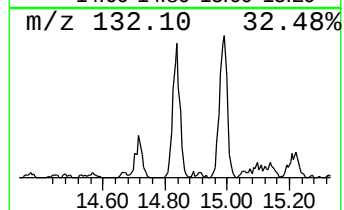
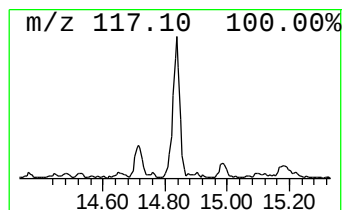
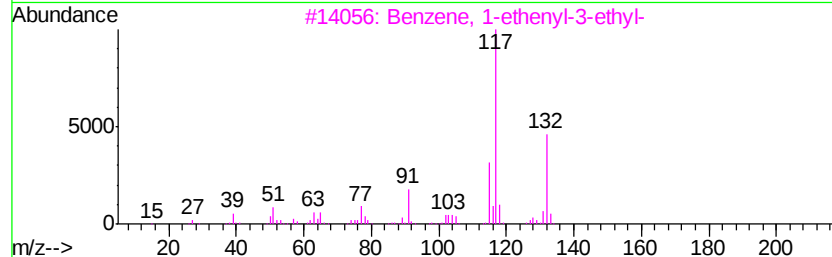
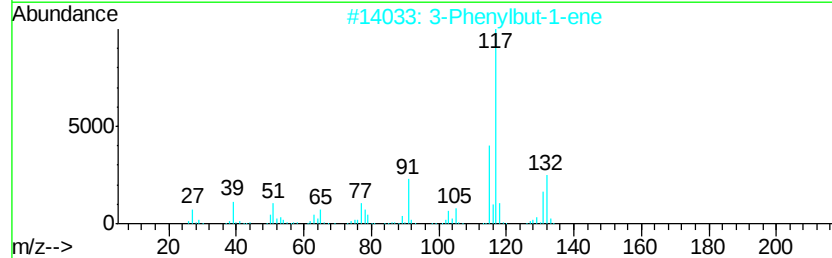
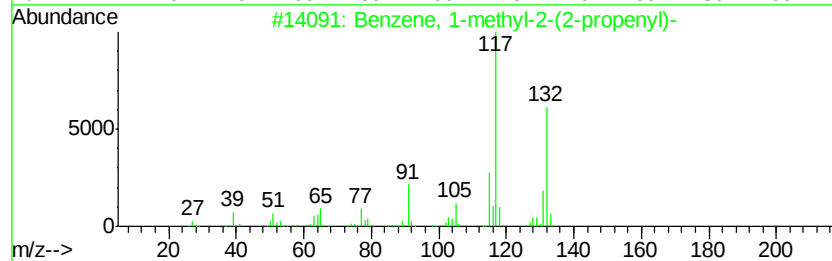
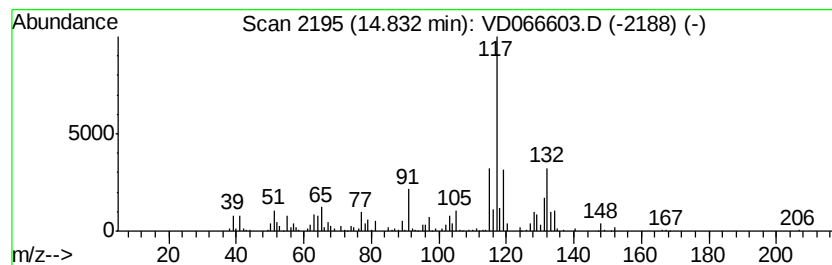
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Benzene, 1-methyl-2-(2-prop... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	14.22 ug/l	596310	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-methyl-2-(2-propenyl)-	132 C10H12	001587-04-8 76
2		3-Phenylbut-1-ene	132 C10H12	000934-10-1 76
3		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 76
4		1-Phenyl-1-butene	132 C10H12	000824-90-8 76
5		Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000767-99-7 76



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SU-04-090820

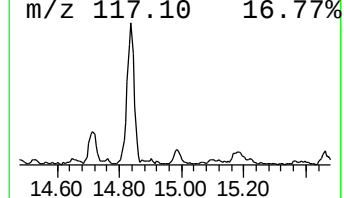
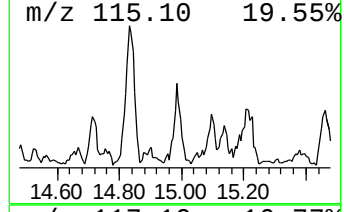
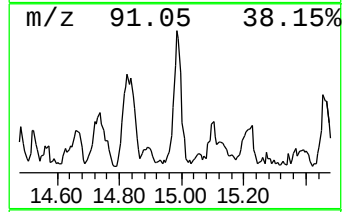
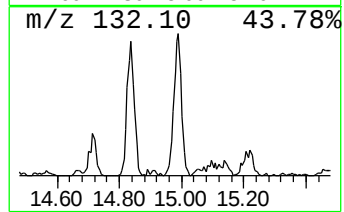
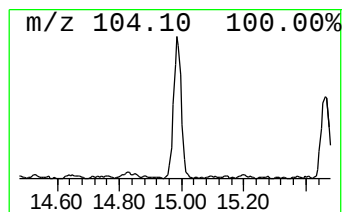
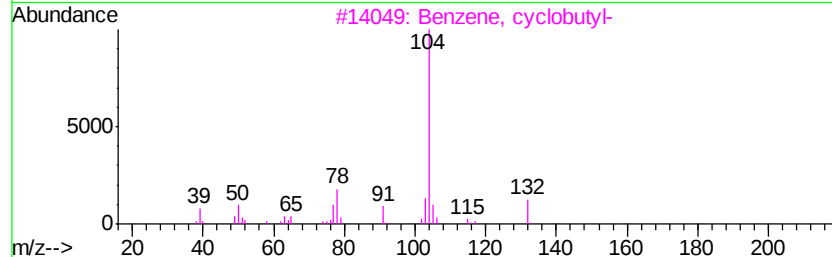
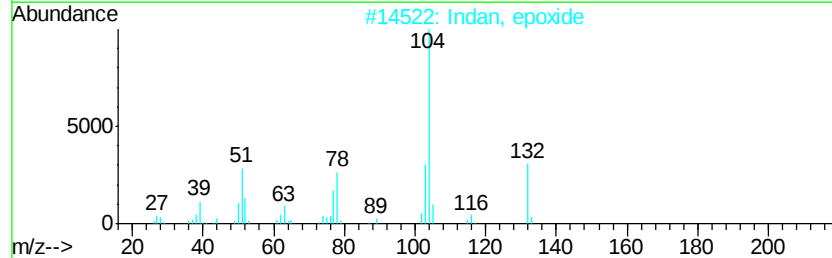
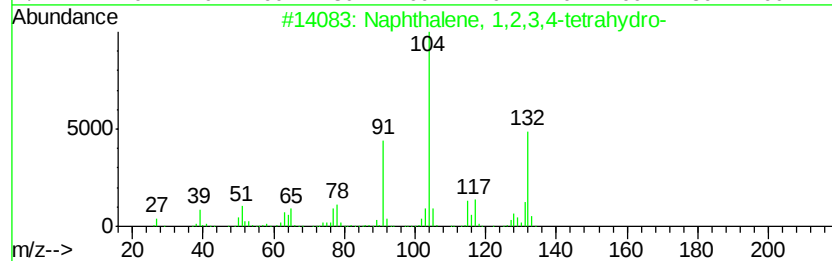
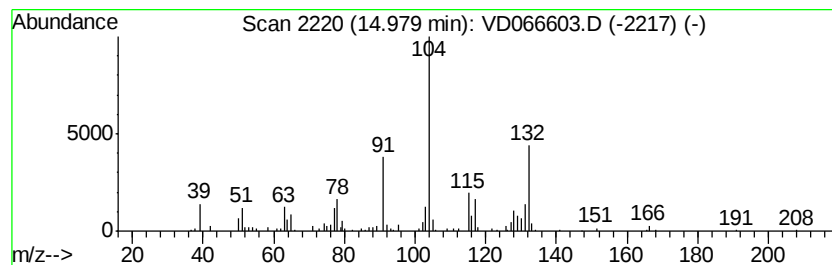
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Quant Title : SW846 8260

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

Peak Number 5 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.98	6.69 ug/l	280729	1,4-Dichlorobenzene-d4	13.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2		Indan, epoxide	132	C9H8O	000768-22-9	55
3		Benzene, cyclobutyl-	132	C10H12	004392-30-7	50
4		2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	49
5		2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD090920\
Data File : VD066603.D
Acq On : 10 Sep 2020 02:11
Operator : VA/SY
Sample : L3870-01
Misc : 6.62G/5.00ml/MSVOA D/SOIL
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-04-090820

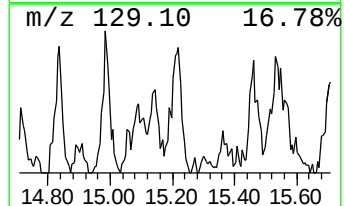
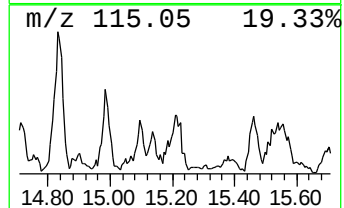
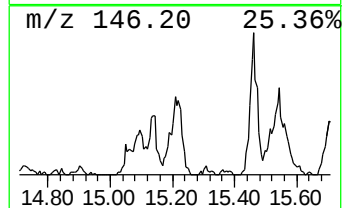
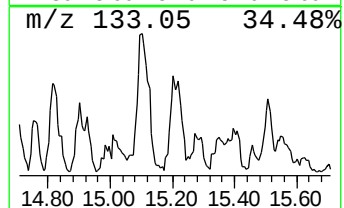
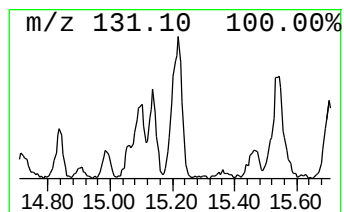
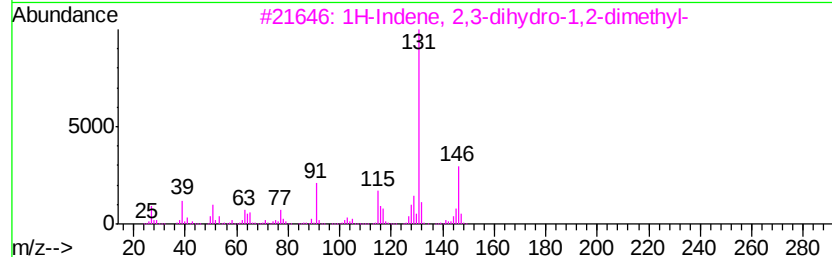
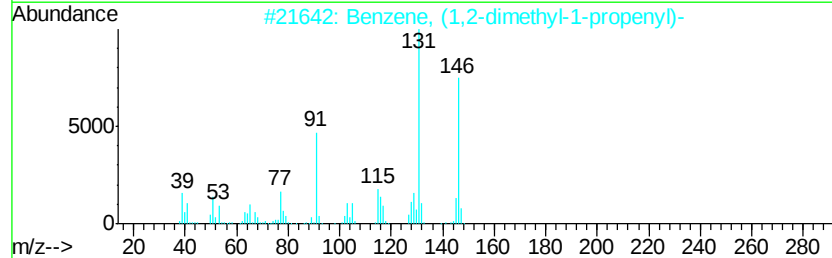
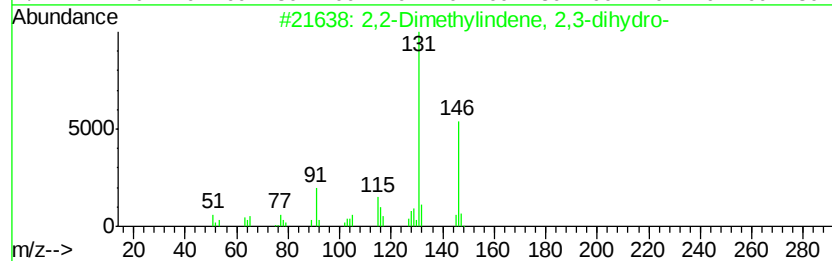
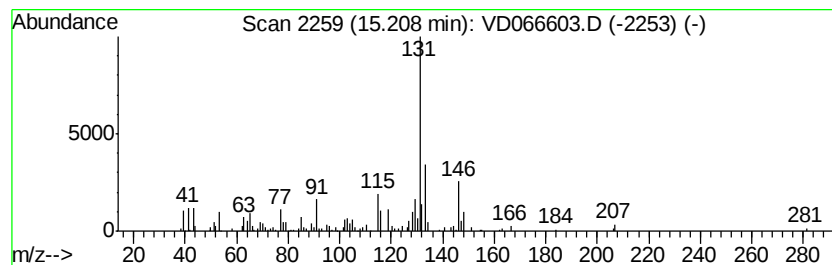
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Quant Title : SW846 8260

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

Peak Number 6 2,2-Dimethylindene, 2,3-dih... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.21	6.20 ug/l	260041	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	81
2		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	81
3		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	81
4		Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	76
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD090920\
Data File : VD066603.D
Acq On : 10 Sep 2020 02:11
Operator : VA/SY
Sample : L3870-01
Misc : 6.62G/5.00ml/MSVOA D/SOIL
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-04-090820

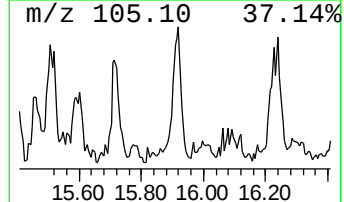
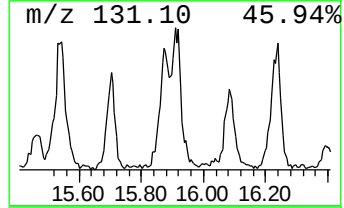
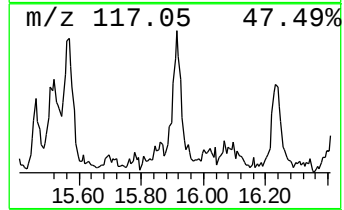
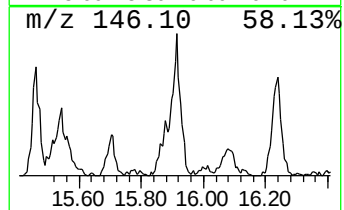
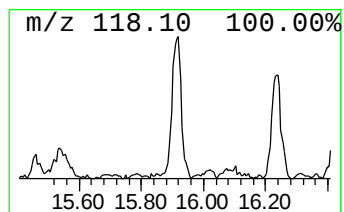
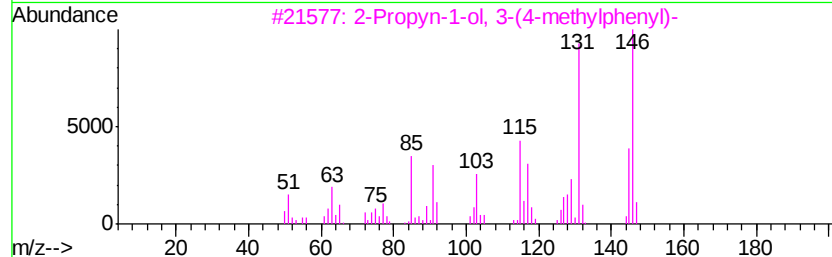
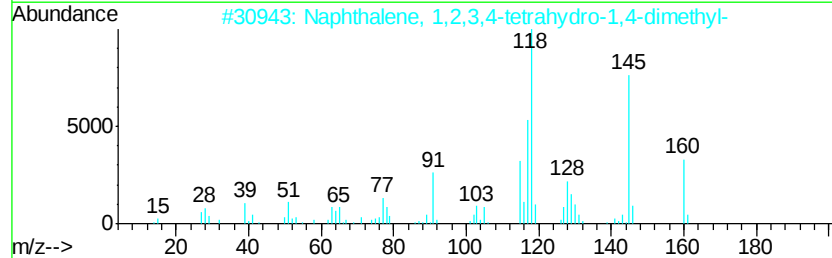
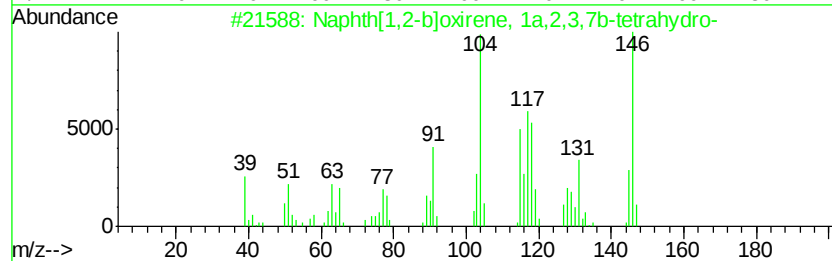
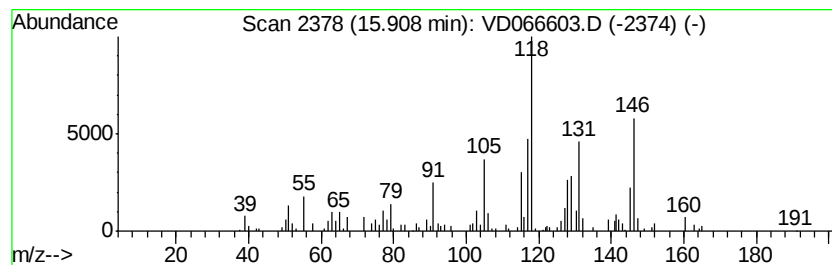
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Quant Title : SW846 8260

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

Peak Number 7 Naphth[1,2-b]oxirene, 1a,2,... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	7.17 ug/l	300705	1,4-Dichlorobenzene-d4	13.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	64
2		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	52
3		2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	49
4		2-Ethyl-2,3-dihydro-1H-indene	146	C11H14	056147-63-8	46
5		Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD090920\
Data File : VD066603.D
Acq On : 10 Sep 2020 02:11
Operator : VA/SY
Sample : L3870-01
Misc : 6.62G/5.00ml/MSVOA D/SOIL
ALS Vial : 35 Sample Multiplier: 1

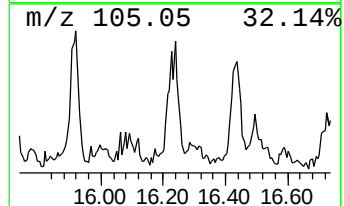
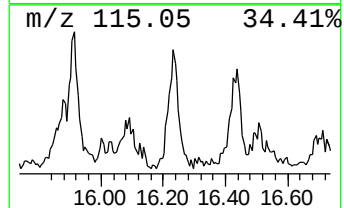
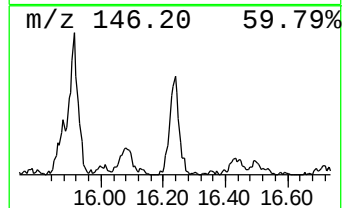
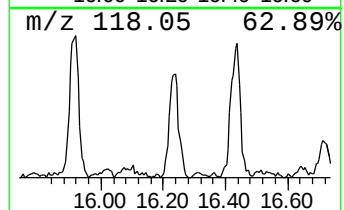
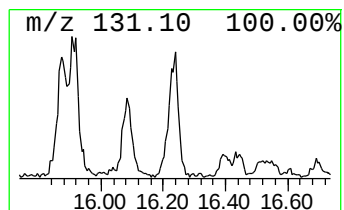
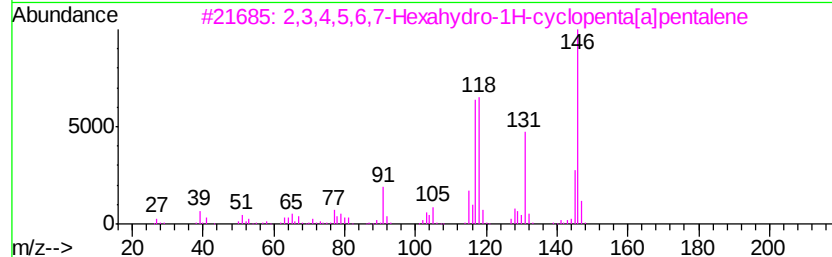
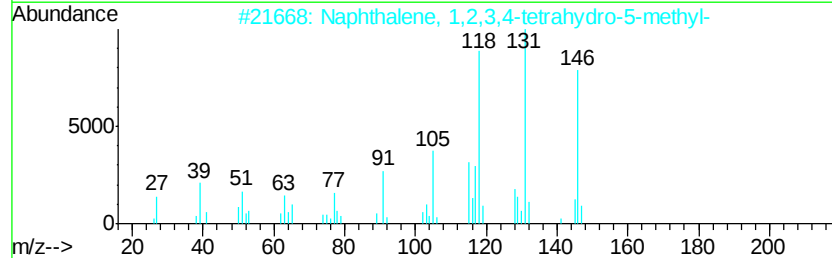
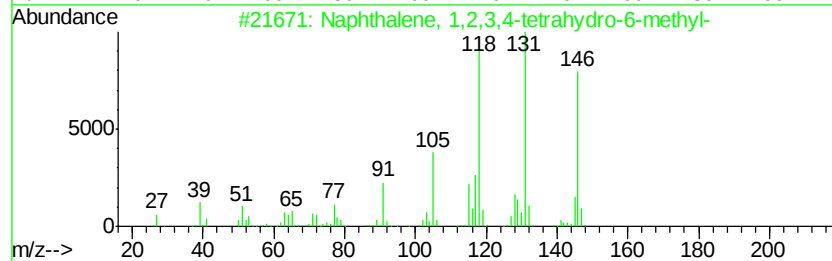
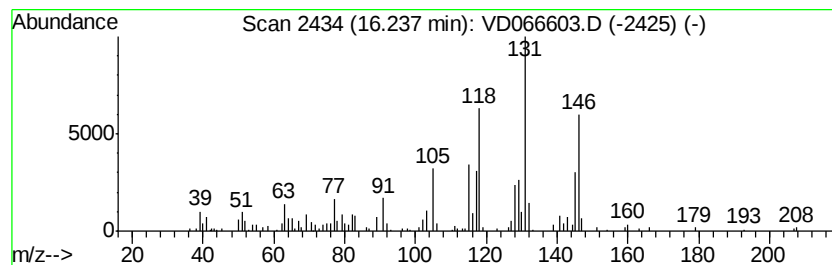
Instrument :
MSVOA_D
ClientSampled :
SU-04-090820

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.24	7.92 ug/l	332008	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9	87
2	Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5	87
3	2,3,4,5,6,7-Hexahydro-1H-cyclope...	146 C11H14	1000189-31-0	80
4	1H-Inden-1-one, 2,3-dihydro-2-me...	146 C10H10O	017496-14-9	64
5	Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5	62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD090920\
Data File : VD066603.D
Acq On : 10 Sep 2020 02:11
Operator : VA/SY
Sample : L3870-01
Misc : 6.62G/5.00ml/MSVOA D/SOIL
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-04-090820

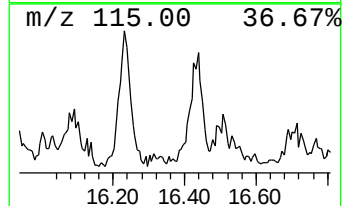
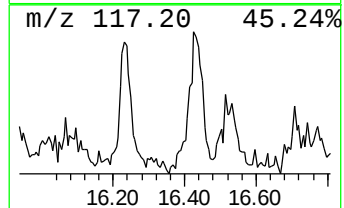
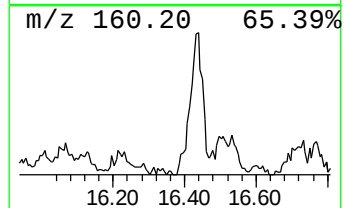
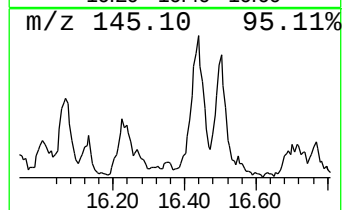
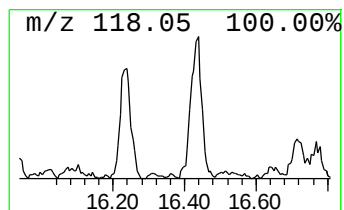
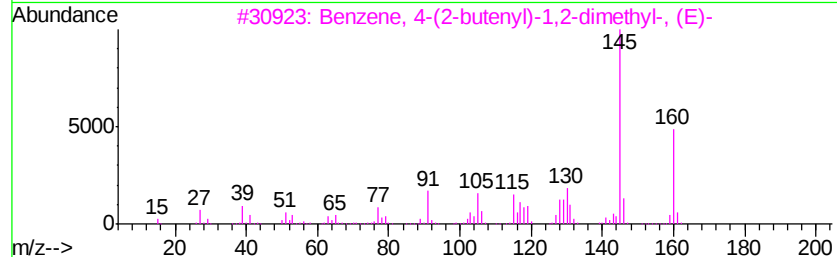
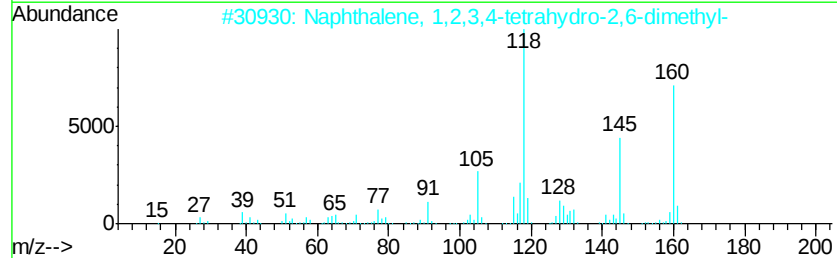
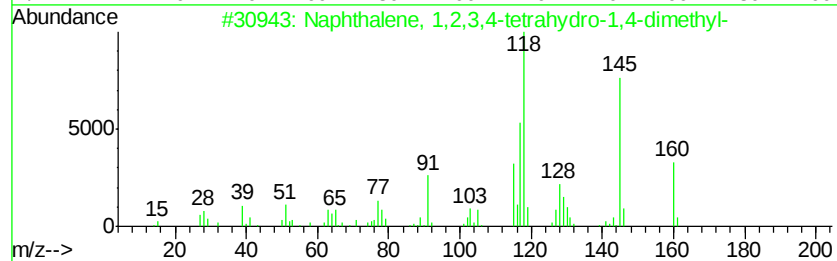
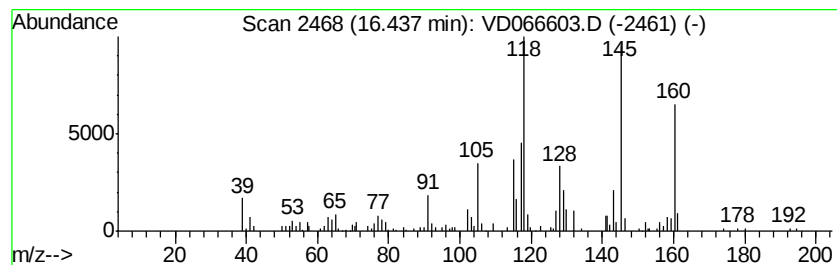
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.44	5.38 ug/l	225777	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	87
2		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	52
3		Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	52
4		Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	49
5		Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	49



Data Path : Z:\V0ASRV\HPCHEM1\MSV0A_D\DATA\VD090920\
Data File : VD066603.D
Acq On : 10 Sep 2020 02:11
Operator : VA/SY
Sample : L3870-01
Misc : 6.62G/5.00ml/MSV0A_D/SOIL
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSV0A_D
ClientSampleId :
SU-04-090820

Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_D\METHOD\82D090320S.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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unknown14.23	14.23	5.8	ug/l	242134	4	13.58	2097180	50.0
Dodecane	14.76	7.9	ug/l	329753	4	13.58	2097180	50.0
Benzene, 1-methyl...	14.83	14.2	ug/l	596310	4	13.58	2097180	50.0
Naphthalene, 1,2,...	14.98	6.7	ug/l	280729	4	13.58	2097180	50.0
2,2-Dimethylinden...	15.21	6.2	ug/l	260041	4	13.58	2097180	50.0
Naphth[1,2-b]oxir...	15.91	7.2	ug/l	300705	4	13.58	2097180	50.0
Naphthalene, 1,2,...	16.24	7.9	ug/l	332008	4	13.58	2097180	50.0
Naphthalene, 1,2,...	16.44	5.4	ug/l	225777	4	13.58	2097180	50.0