

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD120420\
 Data File : VD067881.D
 Acq On : 04 Dec 2020 12:57
 Operator : VA/SY
 Sample : L4950-01
 Misc : 5.71G/5.00ml/MSVOA D/SOIL
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 PL-02-120420

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\82D120120S.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	2.144	33	38	41	rBV3	1680	2742	1.65%	0.180%
2	7.697	978	982	986	rBV2	1686	3053	1.83%	0.200%
3	7.920	1009	1020	1025	rBV2	26450	61274	36.81%	4.015%
4	7.985	1025	1031	1039	rVB2	50678	113873	68.40%	7.462%
5	8.332	1083	1090	1101	rVB2	28325	63767	38.30%	4.179%
6	8.867	1173	1181	1190	rBV	60275	118834	71.38%	7.787%
7	10.344	1423	1432	1442	rBV	86635	159763	95.97%	10.470%
8	11.414	1611	1614	1616	rVV2	2409	2883	1.73%	0.189%
9	11.455	1619	1621	1623	rVV3	3219	3381	2.03%	0.222%
10	11.650	1646	1654	1663	rBV	85738	146405	87.94%	9.594%
11	12.638	1815	1822	1832	rVB	70583	119161	71.58%	7.809%
12	12.897	1862	1866	1867	rVV2	1459	1928	1.16%	0.126%
13	12.914	1867	1869	1871	rVV2	2516	2652	1.59%	0.174%
14	12.955	1871	1876	1881	rVV6	7925	15404	9.25%	1.009%
15	13.126	1897	1905	1909	rBV5	3970	7215	4.33%	0.473%
16	13.173	1909	1913	1914	rBV2	1735	1685	1.01%	0.110%
17	13.279	1925	1931	1932	rBV	1945	2377	1.43%	0.156%
18	13.308	1934	1936	1939	rVB	1955	1874	1.13%	0.123%
19	13.402	1948	1952	1953	rVV3	2975	2907	1.75%	0.191%
20	13.444	1956	1959	1960	rVV3	3141	3187	1.91%	0.209%
21	13.467	1960	1963	1966	rVV5	4726	6198	3.72%	0.406%
22	13.514	1966	1971	1972	rVV4	2524	2830	1.70%	0.185%
23	13.579	1975	1982	1991	rVB	95988	166477	100.00%	10.910%
24	13.732	2004	2008	2009	rVV2	2849	2457	1.48%	0.161%
25	13.773	2009	2015	2018	rVV3	8640	15291	9.19%	1.002%
26	13.814	2018	2022	2031	rVB3	10365	20353	12.23%	1.334%
27	13.896	2031	2036	2040	rBV4	11562	17738	10.65%	1.162%
28	13.967	2045	2048	2054	rVV3	4919	7290	4.38%	0.478%
29	14.032	2054	2059	2060	rVV3	2383	2993	1.80%	0.196%
30	14.067	2062	2065	2068	rVV3	4386	6483	3.89%	0.425%
31	14.120	2070	2074	2079	rVV	8368	13483	8.10%	0.884%
32	14.226	2088	2092	2098	rVB5	10810	18026	10.83%	1.181%
33	14.355	2111	2114	2118	rBV5	3564	6122	3.68%	0.401%
34	14.414	2120	2124	2126	rVV4	4510	6152	3.70%	0.403%

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35	14.438	2126	2128	2133	rVV4	4788	7747	4.65%	0.508%
36	14.479	2133	2135	2140	rVB	7337	9078	5.45%	0.595%
37	14.526	2140	2143	2147	rBV3	6027	8106	4.87%	0.531%
38	14.567	2147	2150	2159	rVB4	6699	10832	6.51%	0.710%
39	14.626	2159	2160	2162	rBV2	2091	1922	1.15%	0.126%
40	14.661	2162	2166	2171	rVB8	6076	10342	6.21%	0.678%
41	14.714	2171	2175	2178	rBV3	6065	10560	6.34%	0.692%
42	14.755	2178	2182	2186	rVB5	12172	20673	12.42%	1.355%
43	14.820	2186	2193	2200	rBV4	21499	56091	33.69%	3.676%
44	14.985	2217	2221	2227	rBV5	13192	20490	12.31%	1.343%
45	15.061	2231	2234	2236	rVB4	2124	1928	1.16%	0.126%
46	15.096	2236	2240	2243	rBV4	9517	13707	8.23%	0.898%
47	15.132	2245	2246	2249	rVV3	5693	3982	2.39%	0.261%
48	15.220	2253	2261	2266	rVB7	10659	24044	14.44%	1.576%
49	15.461	2297	2302	2307	rBV6	7413	14677	8.82%	0.962%
50	15.520	2307	2312	2314	rVV5	7494	13500	8.11%	0.885%
51	15.543	2314	2316	2318	rVV3	6576	7595	4.56%	0.498%
52	15.602	2322	2326	2330	rVB6	8593	13473	8.09%	0.883%
53	15.702	2335	2343	2344	rBV7	7584	12707	7.63%	0.833%
54	15.773	2350	2355	2356	rBV4	2109	2511	1.51%	0.165%
55	15.838	2363	2366	2368	rBV5	2859	3104	1.86%	0.203%
56	15.861	2368	2370	2371	rVV2	4736	4170	2.50%	0.273%
57	15.914	2371	2379	2383	rVB5	14979	29560	17.76%	1.937%
58	16.002	2388	2394	2398	rBV5	4291	8040	4.83%	0.527%
59	16.079	2401	2407	2411	rBV7	4331	8585	5.16%	0.563%
60	16.185	2424	2425	2427	rBV	2045	1699	1.02%	0.111%
61	16.226	2427	2432	2439	rVB4	9804	20537	12.34%	1.346%
62	16.349	2449	2453	2456	rBV3	2629	3307	1.99%	0.217%
63	16.438	2461	2468	2473	rVV7	10201	23553	14.15%	1.543%
64	16.502	2473	2479	2480	rVV5	5064	9136	5.49%	0.599%
65	16.543	2483	2486	2487	rVV3	3522	3701	2.22%	0.243%
66	16.561	2487	2489	2492	rVV3	3914	4758	2.86%	0.312%
67	16.696	2508	2512	2513	rBV2	7986	7216	4.33%	0.473%
68	16.708	2513	2514	2519	rVB4	5901	8386	5.04%	0.550%

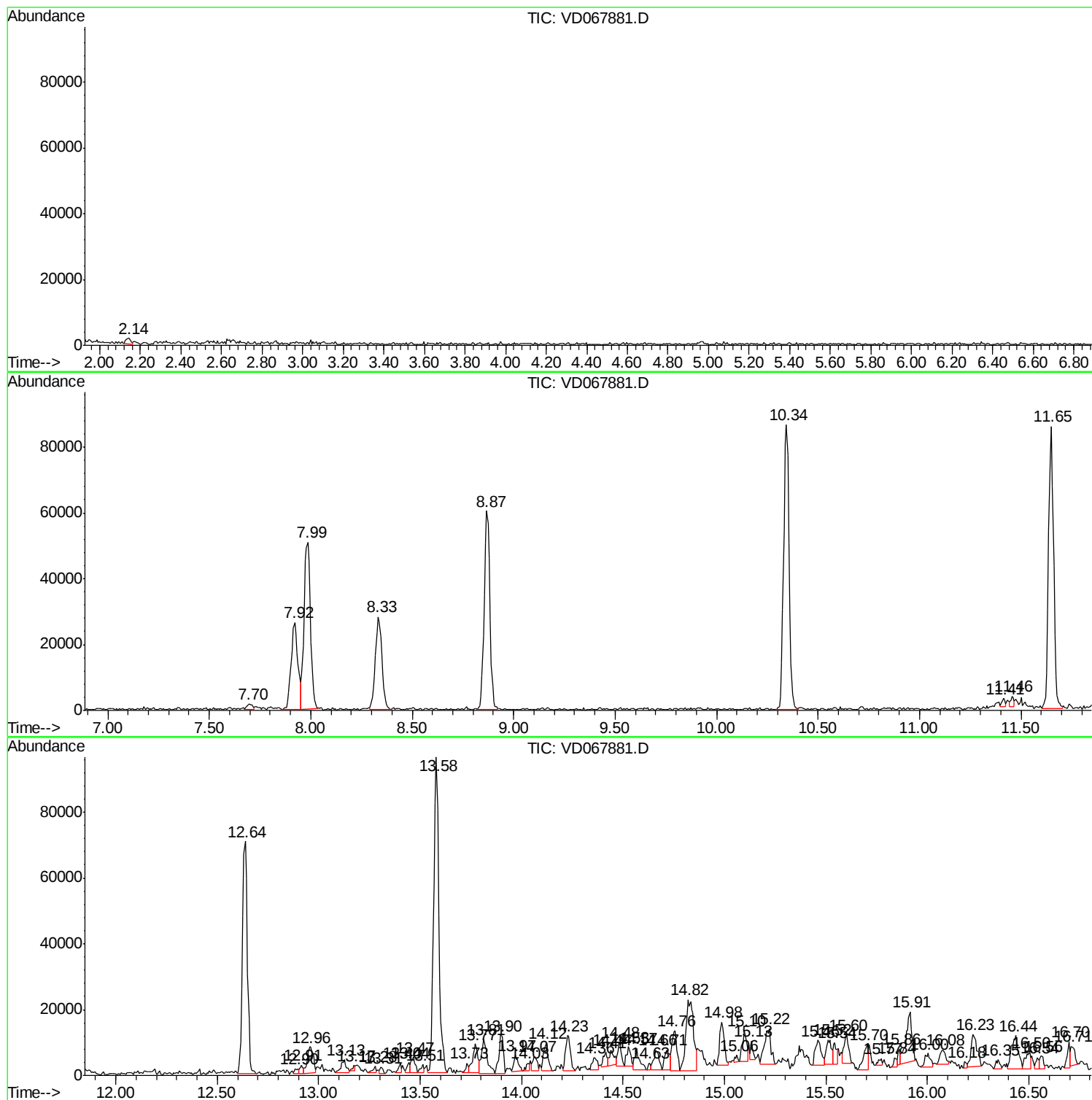
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Quant Title : SW846 8260

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



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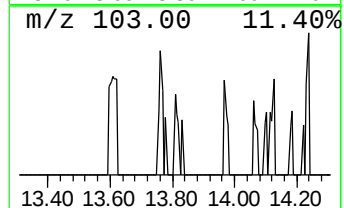
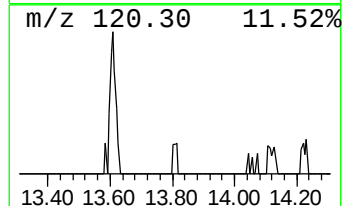
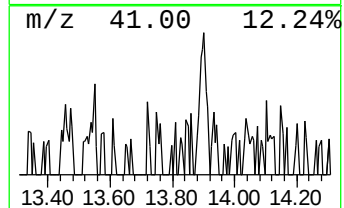
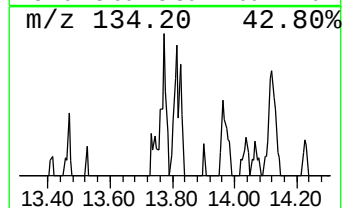
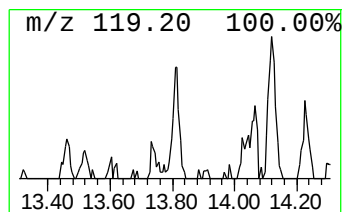
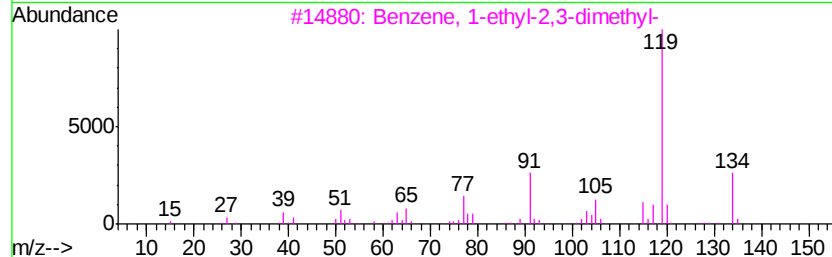
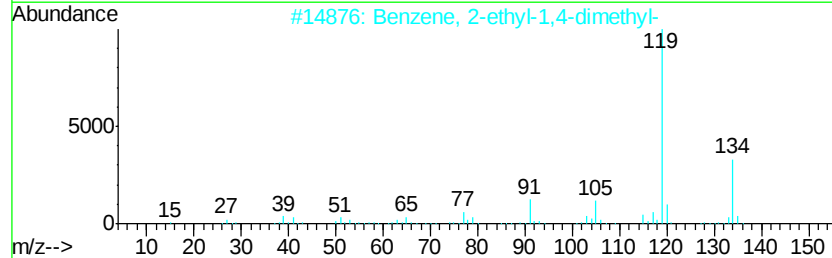
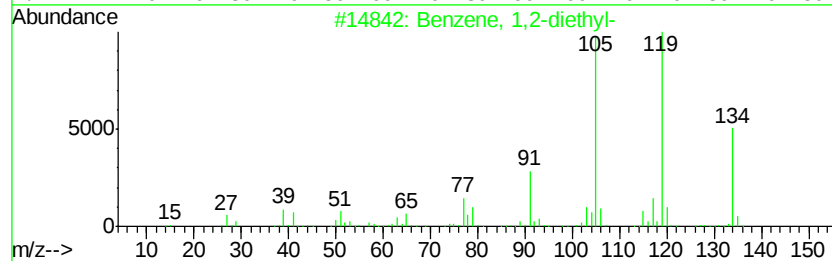
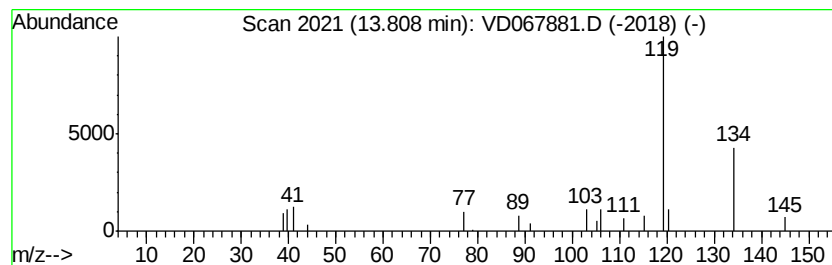
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1,2-diethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.81	6.11 ug/l	20353	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	80
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	72
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	72
4		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	72
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	72



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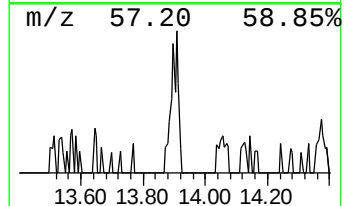
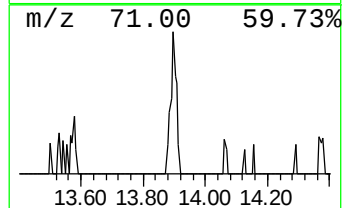
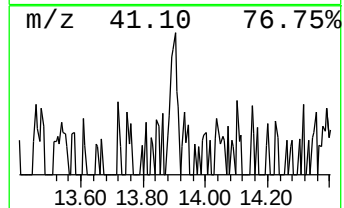
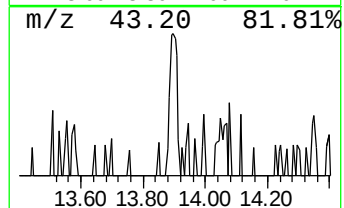
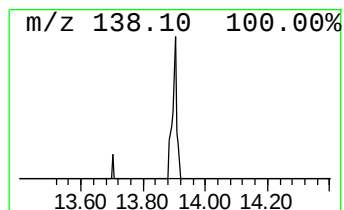
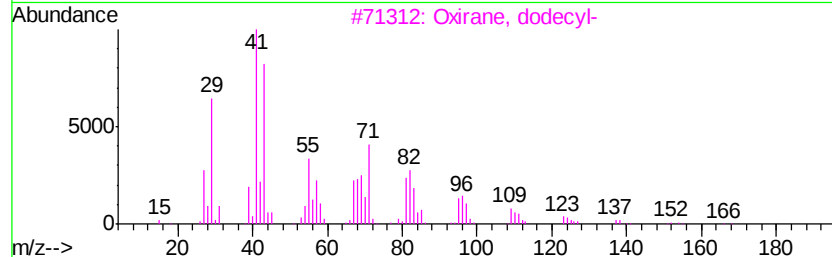
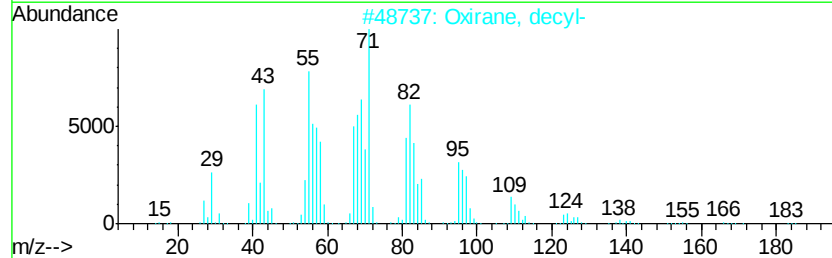
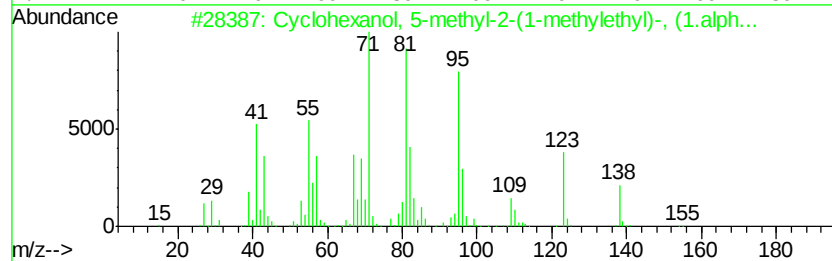
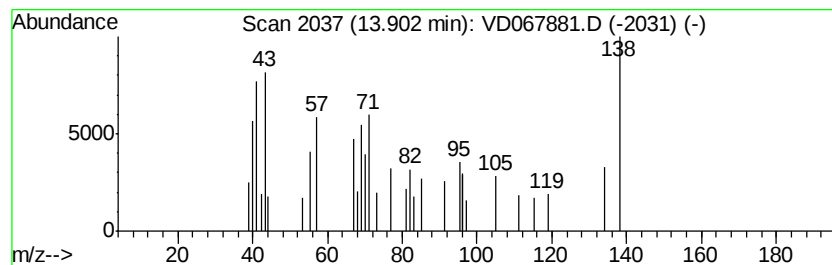
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown13.90 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	5.33 ug/l	17738	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanol, 5-methyl-2-(1-meth...	156	C10H20O	015356-70-4	42
2		Oxirane, decyl-	184	C12H24O	002855-19-8	40
3		Oxirane, dodecyl-	212	C14H28O	003234-28-4	38
4		Hydroxylamine, O-decyl-	173	C10H23NO	029812-79-1	27
5		R(-)3,7-Dimethyl-1,6-octadiene	138	C10H18	010281-56-8	27



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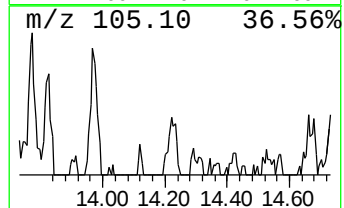
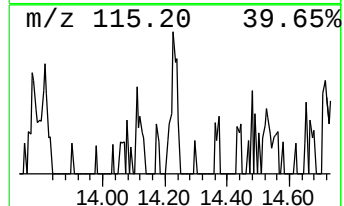
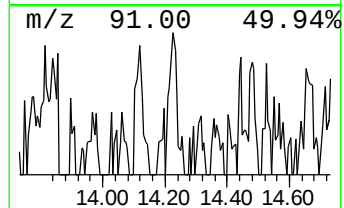
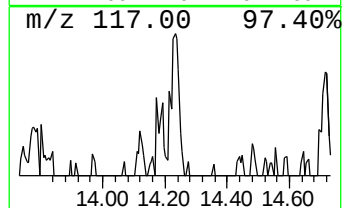
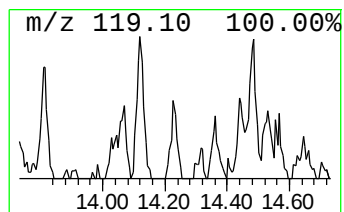
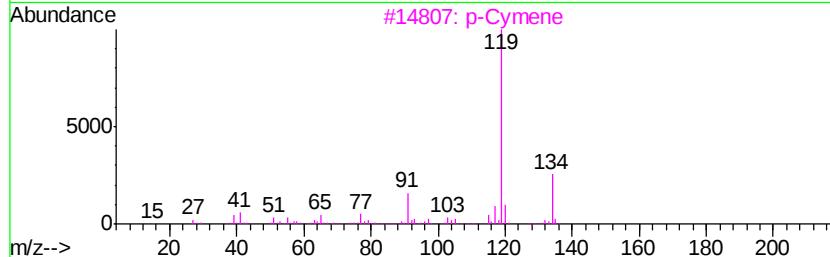
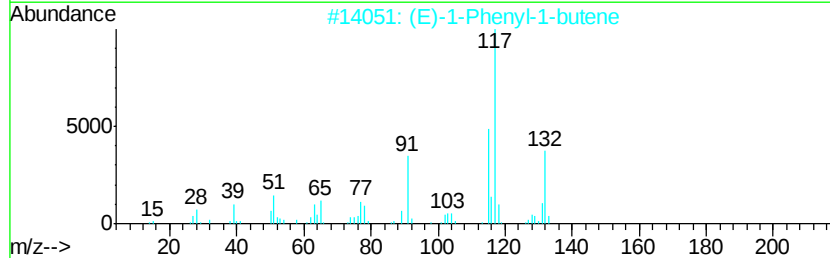
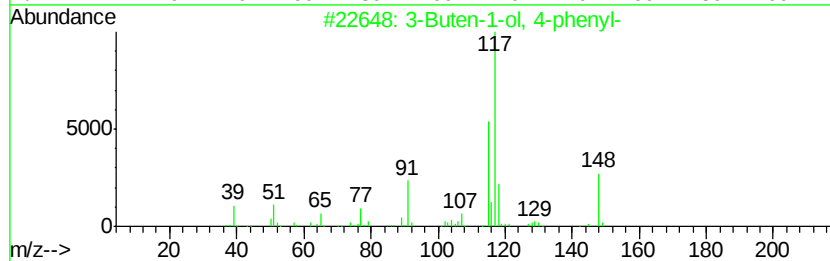
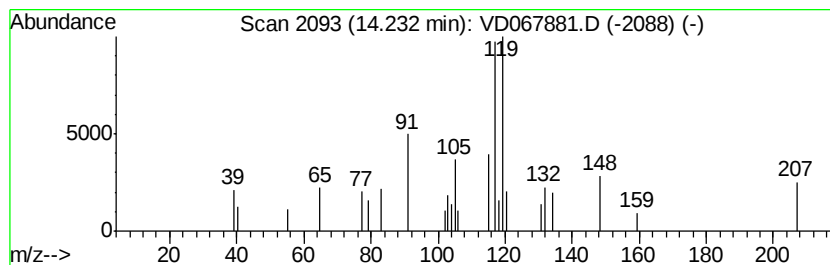
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown14.23 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.23	5.41 ug/l	18026	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Buten-1-ol, 4-phenyl-	148	C10H12O	000937-58-6	43
2		(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	32
3		p-Cymene	134	C10H14	000099-87-6	27
4		2-(3H-Benzo[c]isoxazol-1-yl)methy...	199	C11H9N3O	1000193-59-7	27
5		5-Methyl-6-phenyltetrahydro-1,3-...	207	C11H13NO5	086071-95-6	25



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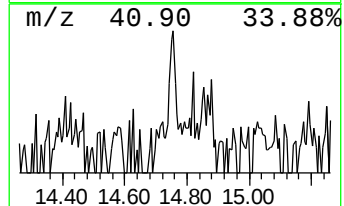
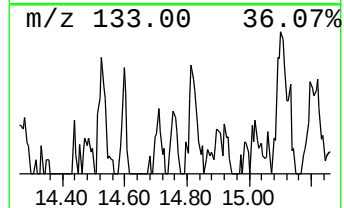
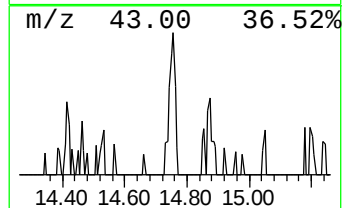
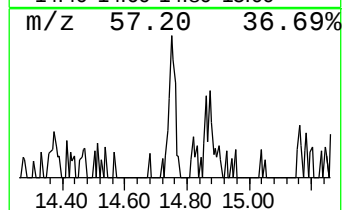
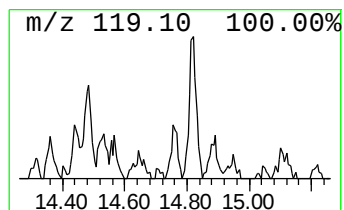
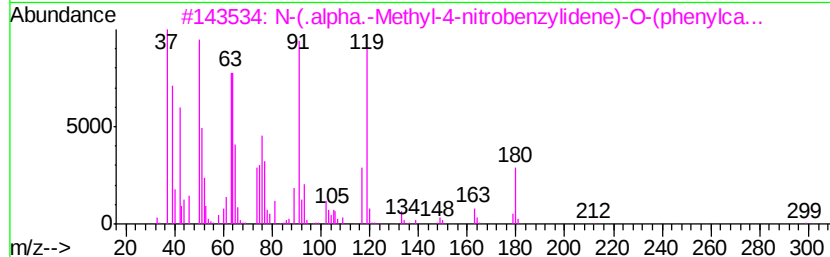
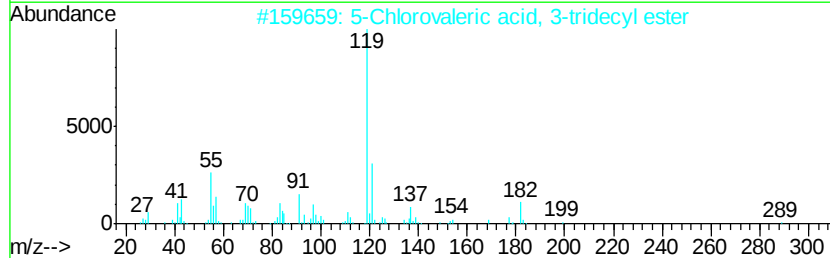
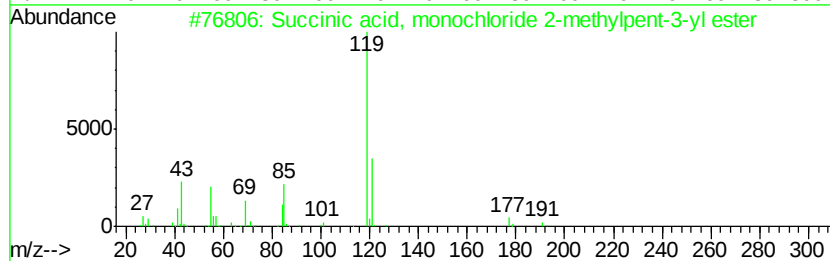
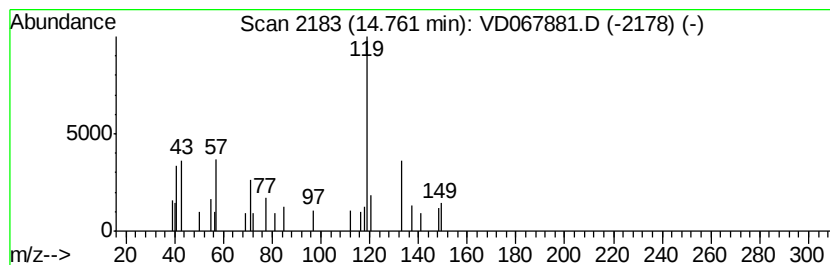
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Peak Number 4 unknown14.76 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	6.21 ug/l	20673	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Succinic acid, monochloride 2-me...	220	C10H17ClO3	1000349-40-1	32
2		5-Chlorovaleric acid, 3-tridecyl...	318	C18H35ClO2	1000299-91-3	23
3		N-(.alpha.-Methyl-4-nitrobenzyl)...	299	C15H13N3O4	1000223-36-2	10
4		4'-Methylpropiofenone	148	C10H12O	005337-93-9	9
5		Propiofenone, 2'-methyl-	148	C10H12O	002040-14-4	9



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD120420\
Data File : VD067881.D
Acq On : 04 Dec 2020 12:57
Operator : VA/SY
Sample : L4950-01
Misc : 5.71G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
PL-02-120420

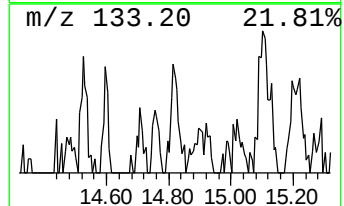
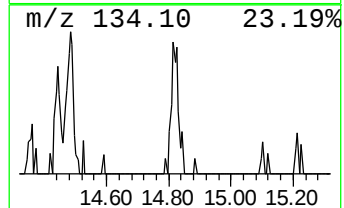
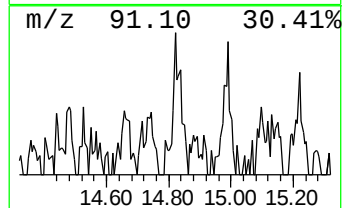
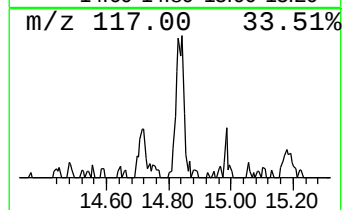
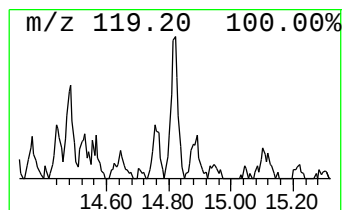
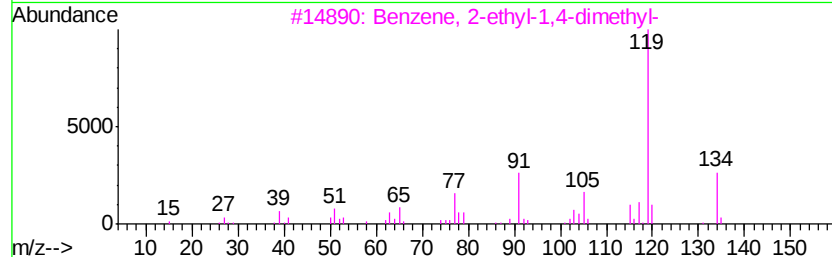
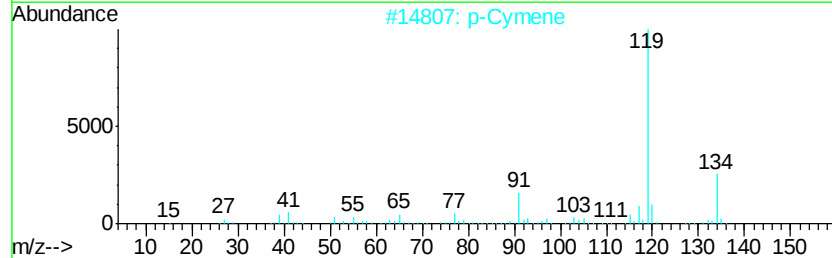
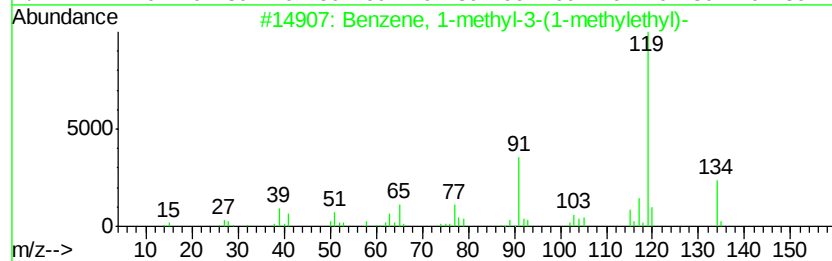
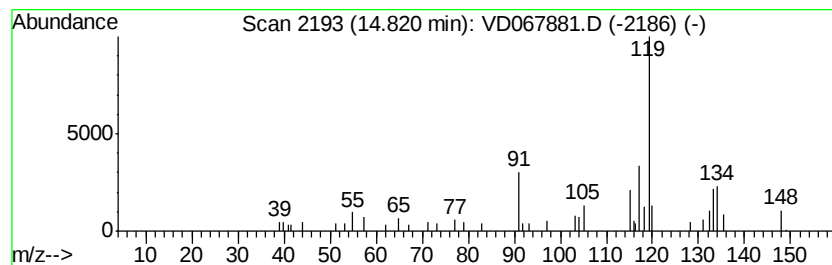
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-methyl-3-(1-meth... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	16.85 ug/l	56091	1,4-Dichlorobenzene-d4	13.58

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene,	1-methyl-3-(1-methyleth...	134	C10H14		000535-77-3	52
2	p-Cymene		134	C10H14		000099-87-6	50
3	Benzene,	2-ethyl-1,4-dimethyl-	134	C10H14		001758-88-9	49
4	Benzene,	1-ethyl-3,5-dimethyl-	134	C10H14		000934-74-7	49
5	Benzene,	2-ethyl-1,3-dimethyl-	134	C10H14		002870-04-4	49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD120420\
Data File : VD067881.D
Acq On : 04 Dec 2020 12:57
Operator : VA/SY
Sample : L4950-01
Misc : 5.71G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
PL-02-120420

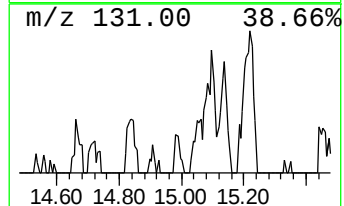
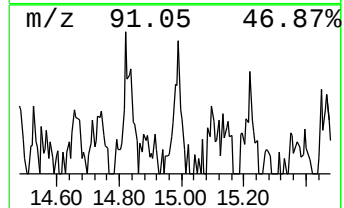
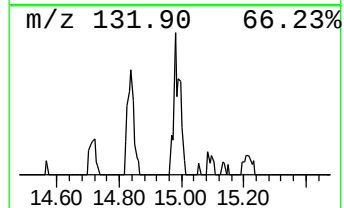
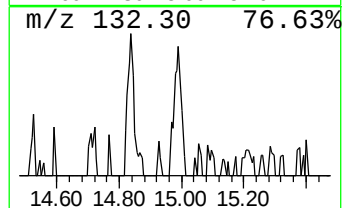
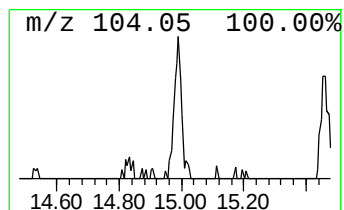
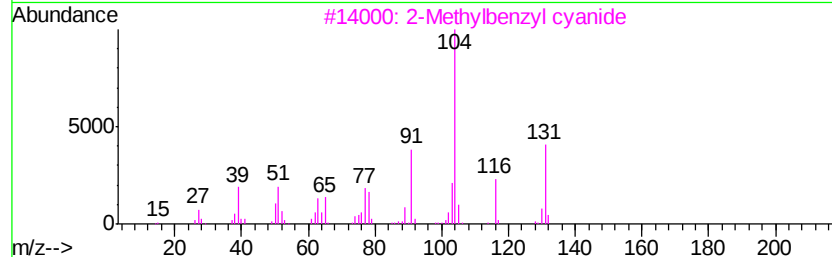
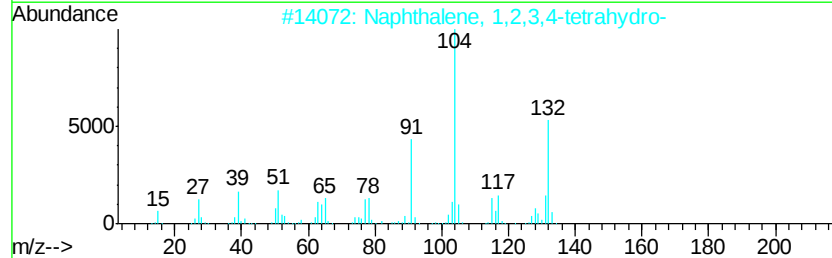
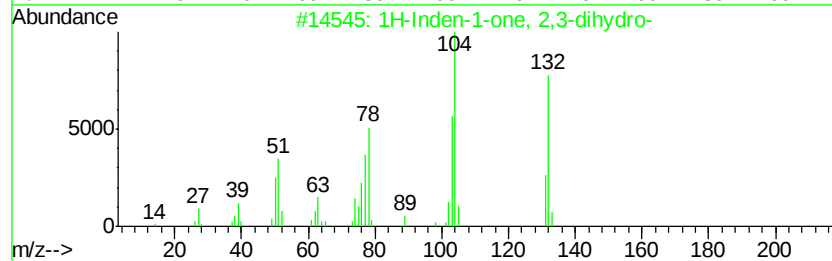
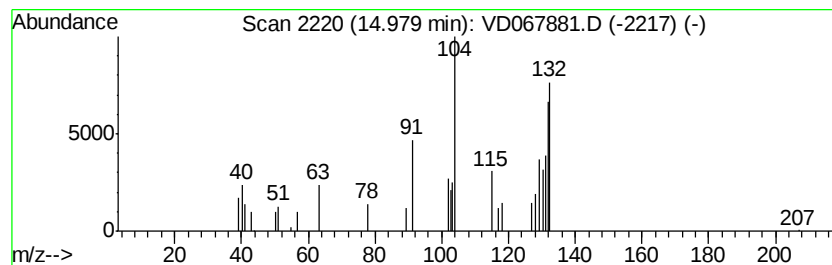
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Quant Title : SW846 8260

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

Peak Number 6 unknown14.98 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.98	6.15 ug/l	20490	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	46
2		Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	43
3		2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	42
4		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	42
5		1-Amino-indan-2-ol	149	C9H11NO	074165-73-4	30



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD120420\
Data File : VD067881.D
Acq On : 04 Dec 2020 12:57
Operator : VA/SY
Sample : L4950-01
Misc : 5.71G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
PL-02-120420

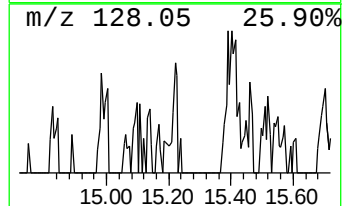
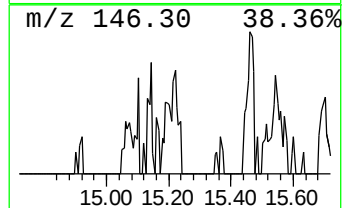
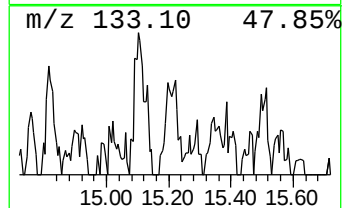
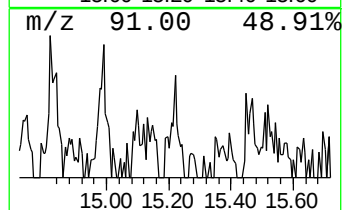
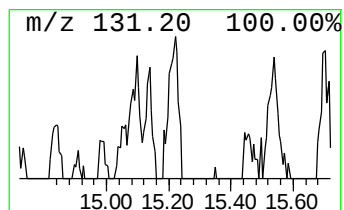
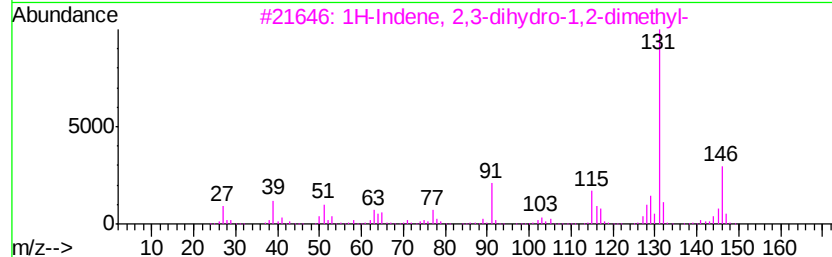
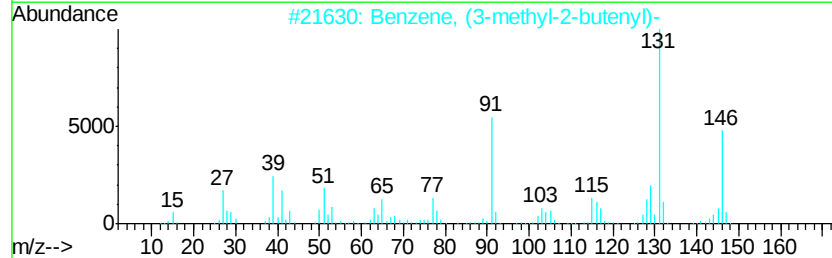
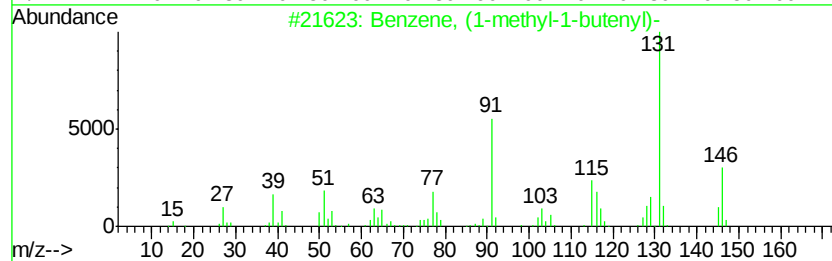
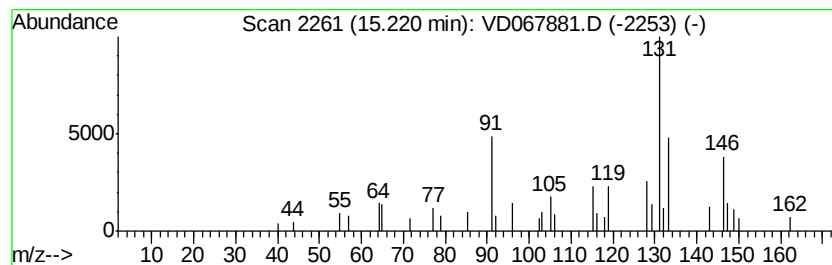
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, (1-methyl-1-butenyl)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.22	7.22 ug/l	24044	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	60
2		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	49
3		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	49
4		Naphthalene, 1,2,3,4-tetrahydro...	146	C11H14	001559-81-5	49
5		1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD120420\
Data File : VD067881.D
Acq On : 04 Dec 2020 12:57
Operator : VA/SY
Sample : L4950-01
Misc : 5.71G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

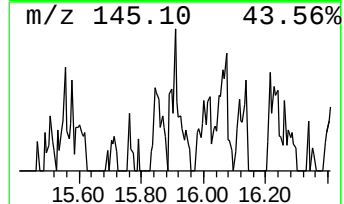
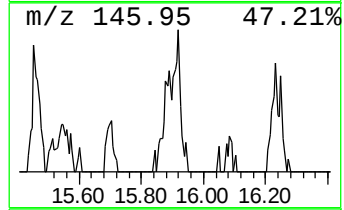
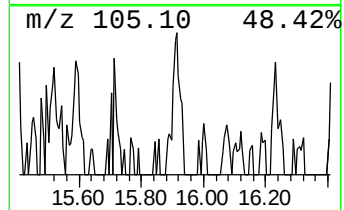
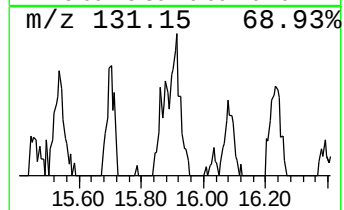
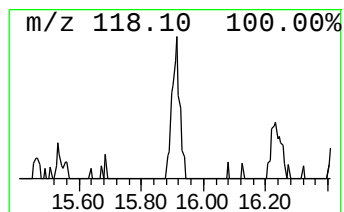
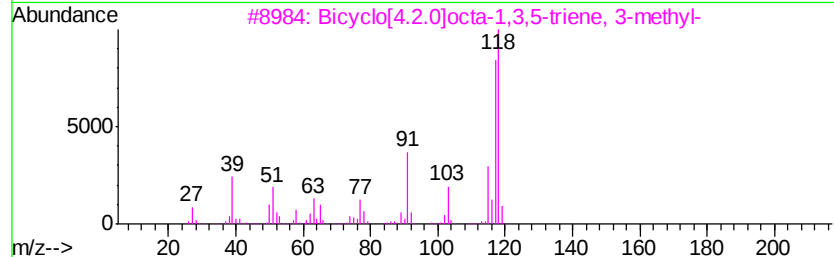
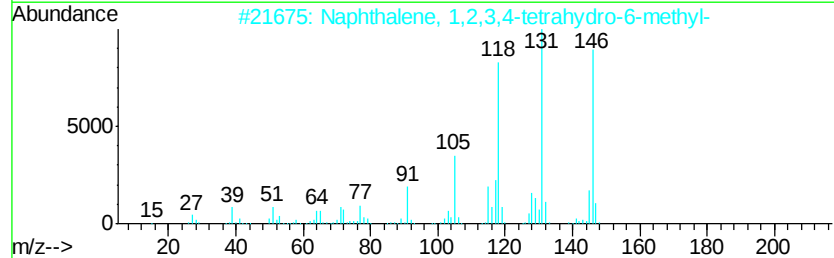
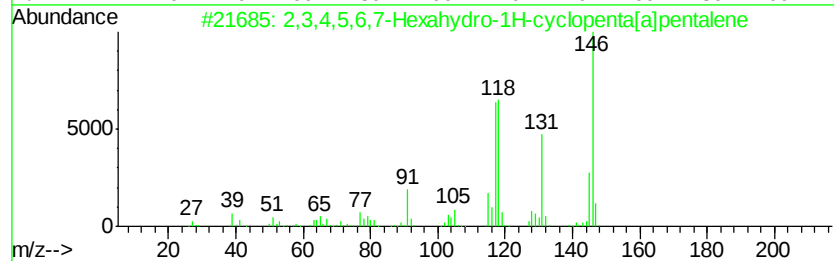
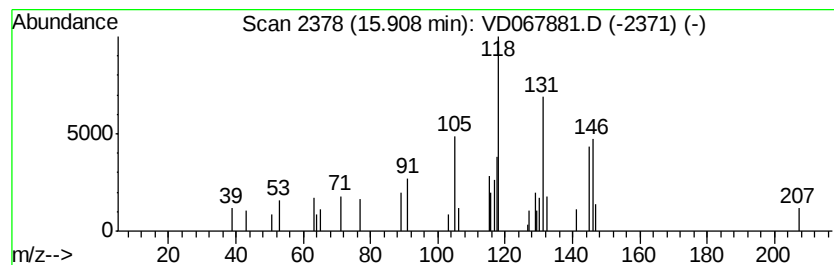
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MSVOA_D
ClientSampleId :
PL-02-120420

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.M
Quant Title : SW846 8260

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

Peak Number 8 2,3,4,5,6,7-Hexahydro-1H-cy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
15.91	8.88 ug/l	29560	1,4-Dichlorobenzene-d4	13.58		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14		1000189-31-0	59
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14		001680-51-9	47
3	Bicvclo[4.2.0]octa-1,3,5-triene,...	118	C9H10		022250-74-4	38
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14		002809-64-5	38
5	Heptacyclo[8.4.0(2,6).0(3,8).0(5...	184	C14H16		1000221-91-1	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD120420\
Data File : VD067881.D
Acq On : 04 Dec 2020 12:57
Operator : VA/SY
Sample : L4950-01
Misc : 5.71G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
PL-02-120420

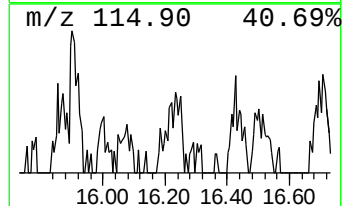
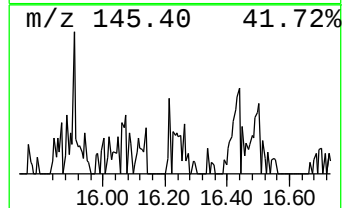
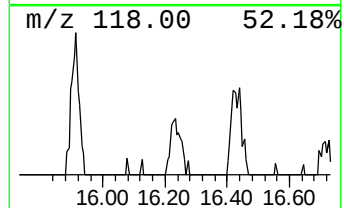
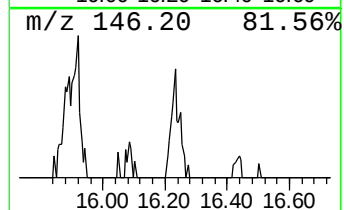
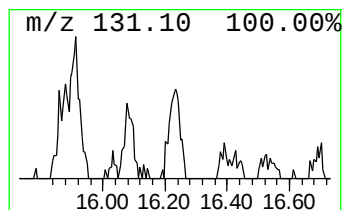
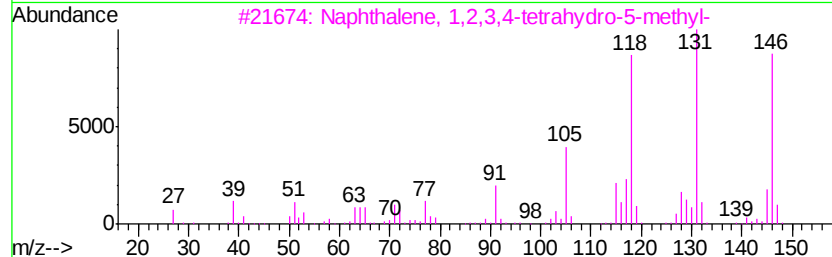
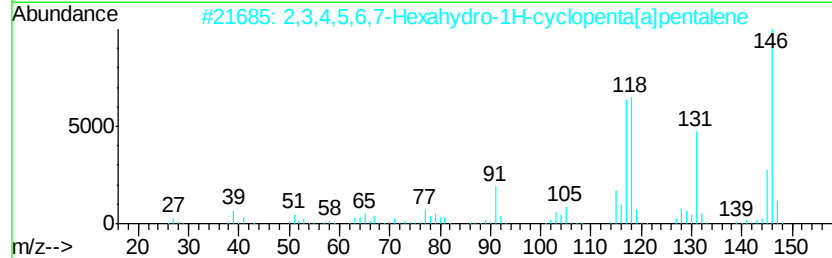
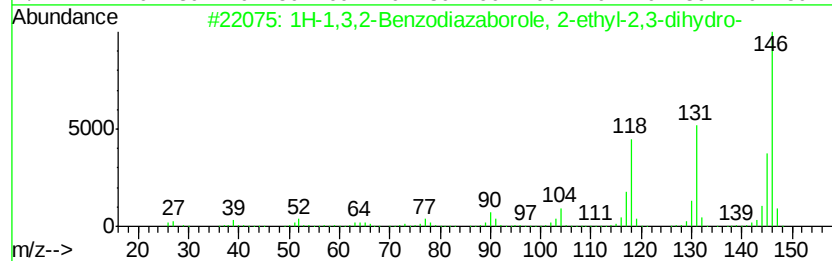
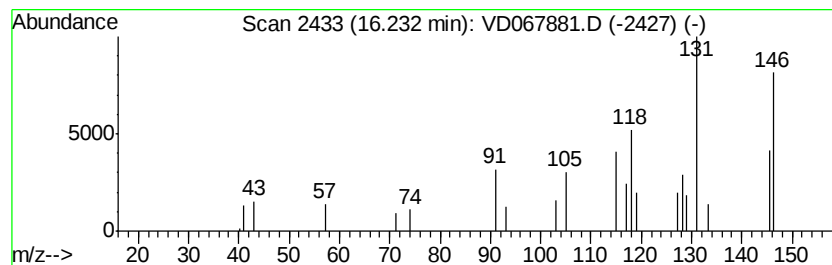
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.M
Quant Title : SW846 8260

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

Peak Number 9 1H-1,3,2-Benzodiazaborole, ... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.23	6.17 ug/l	20537	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	62
2		2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1000189-31-0	58
3		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	52
4		2-(2-Hydroxyphenyl)buta-1,3-diene	146	C10H10O	038865-47-3	47
5		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD120420\
Data File : VD067881.D
Acq On : 04 Dec 2020 12:57
Operator : VA/SY
Sample : L4950-01
Misc : 5.71G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
PL-02-120420

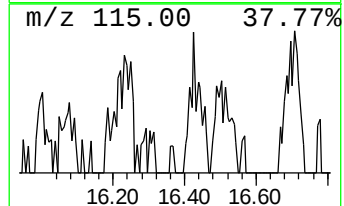
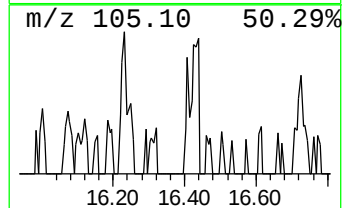
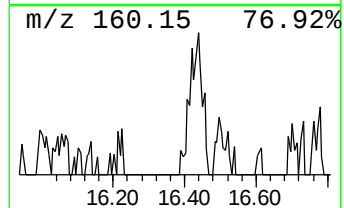
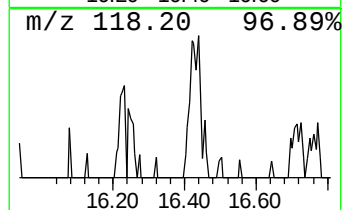
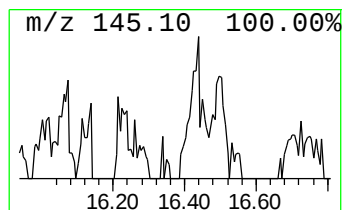
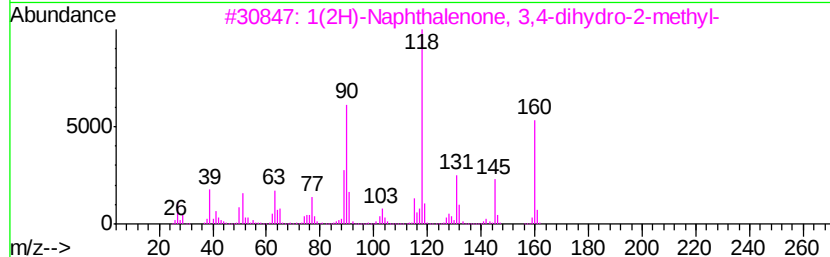
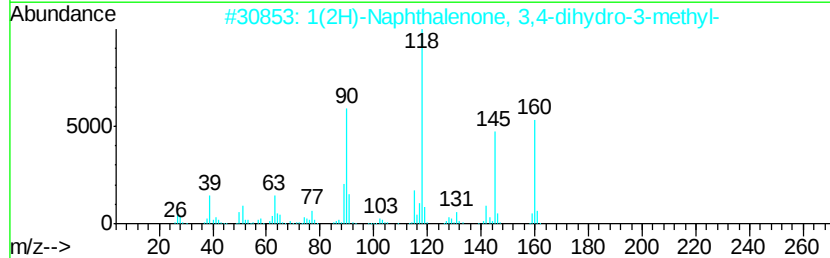
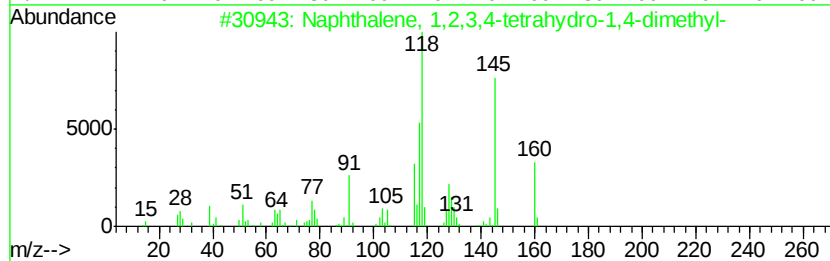
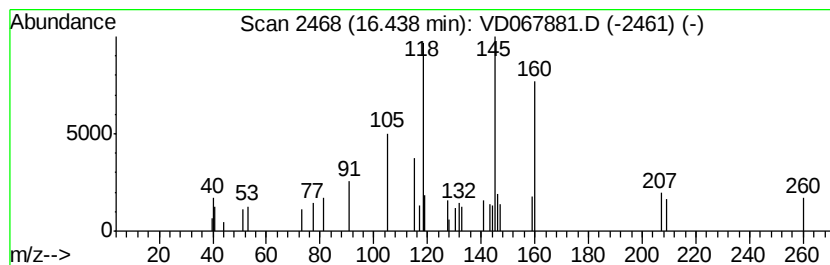
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.44	7.07 ug/l	23553	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	59
2		1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	014944-23-1	49
3		1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	001590-08-5	43
4		1,5-Naphthylidin-2-amine	145	C8H7N3	017965-80-9	43
5		1-Phthalanol, 1,3,3-trimethyl-	178	C11H14O2	001521-94-4	43



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD120420\
Data File : VD067881.D
Acq On : 04 Dec 2020 12:57
Operator : VA/SY
Sample : L4950-01
Misc : 5.71G/5.00ml/MSVOA_D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
PL-02-120420

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.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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unknown13.90	13.90	5.3	ug/l	17738	4	13.58	166477	50.0
unknown14.23	14.23	5.4	ug/l	18026	4	13.58	166477	50.0
unknown14.76	14.76	6.2	ug/l	20673	4	13.58	166477	50.0
Benzene, 1-methyl...	14.82	16.9	ug/l	56091	4	13.58	166477	50.0
unknown14.98	14.98	6.2	ug/l	20490	4	13.58	166477	50.0
Benzene, (1-methy...	15.22	7.2	ug/l	24044	4	13.58	166477	50.0
2,3,4,5,6,7-Hexah...	15.91	8.9	ug/l	29560	4	13.58	166477	50.0
1H-1,3,2-Benzodia...	16.23	6.2	ug/l	20537	4	13.58	166477	50.0
Naphthalene, 1,2,...	16.44	7.1	ug/l	23553	4	13.58	166477	50.0