

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD102021\
 Data File : VD070778.D
 Acq On : 20 Oct 2021 16:20
 Operator : VA/SY
 Sample : M4284-03
 Misc : 5.01G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 30852

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D101521S.M
 Title : SW846 8260

Signal : TIC: VD070778.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.156	372	380	391	rBV	57065	157027	1.71%	0.107%
2	4.956	506	516	529	rBV3	47498	164270	1.79%	0.111%
3	7.914	1008	1019	1024	rBV	343839	849553	9.23%	0.577%
4	7.979	1024	1030	1042	rVB	552707	1247276	13.56%	0.847%
5	8.326	1080	1089	1103	rVB	330509	754005	8.20%	0.512%
6	8.861	1170	1180	1191	rBV	951167	1893690	20.58%	1.285%
7	10.332	1422	1430	1437	rBV	1456411	2731704	29.69%	1.854%
8	10.397	1437	1441	1447	rVB	87293	162139	1.76%	0.110%
9	10.520	1454	1462	1468	rVB3	76925	153005	1.66%	0.104%
10	11.238	1579	1584	1590	rVB2	72988	137812	1.50%	0.094%
11	11.420	1607	1615	1627	rVB5	140625	399771	4.35%	0.271%
12	11.520	1627	1632	1638	rBV	144164	235909	2.56%	0.160%
13	11.638	1645	1652	1662	rBV	1325827	2310545	25.11%	1.568%
14	11.738	1663	1669	1673	rBV	178509	310929	3.38%	0.211%
15	11.849	1678	1688	1697	rBV3	1105268	2388800	25.96%	1.621%
16	11.920	1697	1700	1706	rVB4	105048	192012	2.09%	0.130%
17	12.173	1730	1743	1751	rBV2	600210	1545045	16.79%	1.049%
18	12.261	1753	1758	1762	rVB	228649	341470	3.71%	0.232%
19	12.343	1767	1772	1773	rBV2	145247	172113	1.87%	0.117%
20	12.385	1775	1779	1788	rVB5	230817	552261	6.00%	0.375%
21	12.467	1788	1793	1798	rBV	188948	363987	3.96%	0.247%
22	12.549	1804	1807	1809	rBV2	109156	125005	1.36%	0.085%
23	12.573	1809	1811	1814	rVB2	140834	150220	1.63%	0.102%
24	12.620	1814	1819	1824	rBV	942723	1557381	16.93%	1.057%
25	12.708	1830	1834	1838	rVB4	72094	101484	1.10%	0.069%
26	12.808	1843	1851	1856	rBV	474035	895692	9.74%	0.608%
27	12.873	1856	1862	1866	rBV	1880248	3237252	35.19%	2.197%
28	12.908	1866	1868	1870	rVV	806874	1030344	11.20%	0.699%
29	12.943	1870	1874	1880	rVV3	2253977	4145498	45.06%	2.814%
30	12.996	1880	1883	1889	rVB2	210502	323390	3.52%	0.219%
31	13.108	1889	1902	1909	rBV	1125820	2390055	25.98%	1.622%
32	13.173	1909	1913	1920	rVB3	437879	802075	8.72%	0.544%
33	13.255	1921	1927	1933	rBV	4008582	6350230	69.02%	4.310%
34	13.391	1945	1950	1954	rVB2	375813	571844	6.22%	0.388%
35	13.449	1954	1960	1965	rBV2	896972	1553598	16.89%	1.054%
36	13.502	1965	1969	1974	rBV2	457346	703726	7.65%	0.478%

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Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D101521S.M
 Title : SW846 8260

37	13.561	1974	1979	1982	rBV	1268564	2069406	22.49%	1.405%
38	13.596	1982	1985	1986	rVV	1290403	1613628	17.54%	1.095%
39	13.626	1986	1990	1999	rVB	3795399	6350989	69.03%	4.311%
40	13.761	2003	2013	2016	rBV2	2097202	4124307	44.83%	2.799%
41	13.796	2016	2019	2024	rVB2	1062092	1553438	16.88%	1.054%
42	13.885	2029	2034	2039	rBV2	2212493	3458155	37.59%	2.347%
43	13.955	2041	2046	2052	rVV	940823	1629650	17.71%	1.106%
44	14.020	2052	2057	2059	rVV	1134052	1804617	19.62%	1.225%
45	14.049	2059	2062	2066	rVV	1269603	1884180	20.48%	1.279%
46	14.102	2066	2071	2076	rVV	1766405	2616249	28.44%	1.776%
47	14.167	2076	2082	2085	rVV3	260254	445008	4.84%	0.302%
48	14.214	2085	2090	2095	rVB	1969305	3165400	34.41%	2.148%
49	14.290	2095	2103	2107	rBV4	262689	644458	7.00%	0.437%
50	14.343	2107	2112	2116	rBV3	750797	1363982	14.83%	0.926%
51	14.396	2116	2121	2123	rVV3	934263	1743638	18.95%	1.183%
52	14.426	2123	2126	2129	rVV	1178950	1914117	20.81%	1.299%
53	14.467	2129	2133	2137	rVV	1566609	2298659	24.99%	1.560%
54	14.508	2137	2140	2144	rVV2	1042667	1626250	17.68%	1.104%
55	14.549	2144	2147	2154	rVV4	1262531	1989279	21.62%	1.350%
56	14.614	2154	2158	2159	rVV3	285237	341369	3.71%	0.232%
57	14.649	2160	2164	2168	rVV2	955676	1479735	16.08%	1.004%
58	14.696	2168	2172	2176	rVV2	1955857	3816931	41.49%	2.591%
59	14.737	2176	2179	2184	rVB2	1795250	2710995	29.47%	1.840%
60	14.814	2184	2192	2198	rBV4	3899086	9200152	100.00%	6.245%
61	14.873	2198	2202	2209	rVV2	1394948	2687253	29.21%	1.824%
62	14.967	2212	2218	2225	rVB	2745336	4672327	50.79%	3.171%
63	15.079	2232	2237	2241	rBV2	1559437	2641138	28.71%	1.793%
64	15.196	2249	2257	2260	rBV3	1468943	3389755	36.84%	2.301%
65	15.349	2277	2283	2286	rBV5	608288	1244479	13.53%	0.845%
66	15.443	2293	2299	2303	rBV	1363194	2190910	23.81%	1.487%
67	15.496	2303	2308	2310	rBV2	895545	1529243	16.62%	1.038%
68	15.579	2319	2322	2331	rVB3	1008392	1874042	20.37%	1.272%
69	15.679	2331	2339	2347	rBV3	1243782	2761239	30.01%	1.874%
70	15.755	2347	2352	2358	rVV3	428994	774527	8.42%	0.526%
71	15.890	2360	2375	2384	rVB	3327970	8312287	90.35%	5.642%
72	15.979	2386	2390	2395	rVV5	266486	508486	5.53%	0.345%
73	16.061	2397	2404	2419	rVB2	833173	2747453	29.86%	1.865%
74	16.208	2420	2429	2446	rBV2	1743590	4535358	49.30%	3.078%
75	16.326	2447	2449	2454	rBV5	74582	108344	1.18%	0.074%
76	16.408	2455	2463	2469	rBV	1502801	3203011	34.81%	2.174%

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ClientSampleId :
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Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D101521S.M
Title : SW846 8260

77	16.473	2469	2474	2480	rBV4	802001	1656813	18.01%	1.125%
78	16.684	2502	2510	2516	rBV4	551859	1649497	17.93%	1.120%

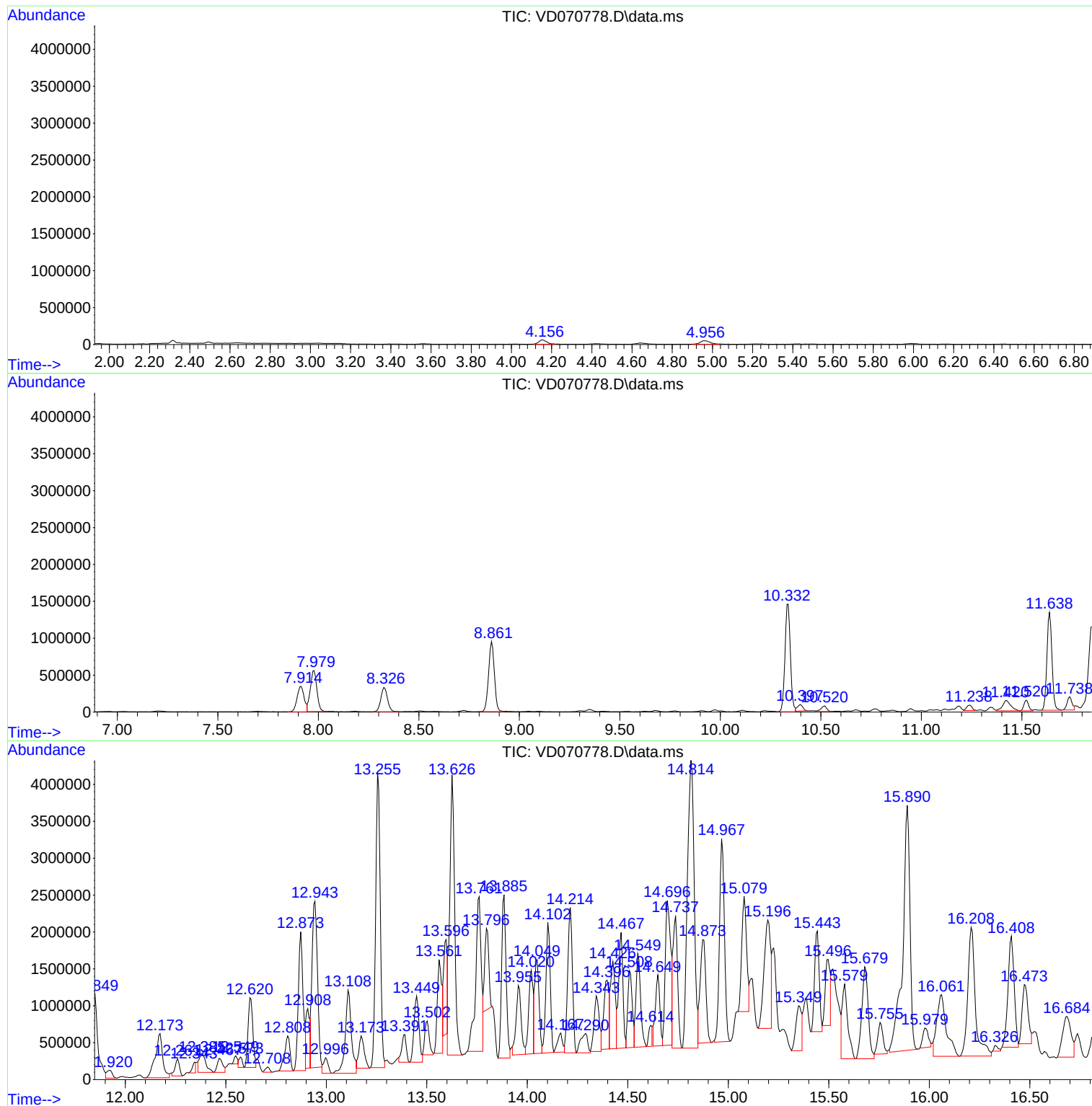
Sum of corrected areas: 147331871

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD102021\
Data File : VD070778.D
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Operator : VA/SY
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Misc : 5.01G/5.00ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
30852

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D101521S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD102021\
Data File : VD070778.D
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Sample : M4284-03
Misc : 5.01G/5.00ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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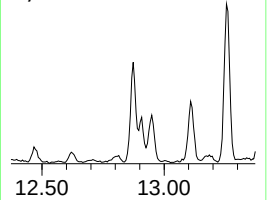
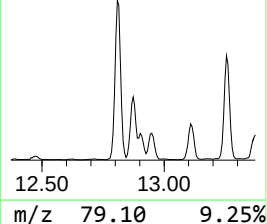
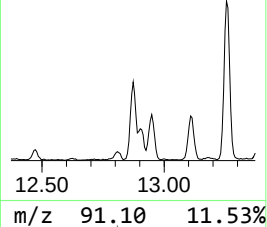
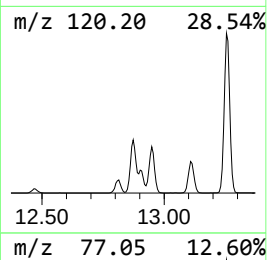
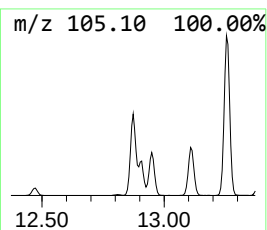
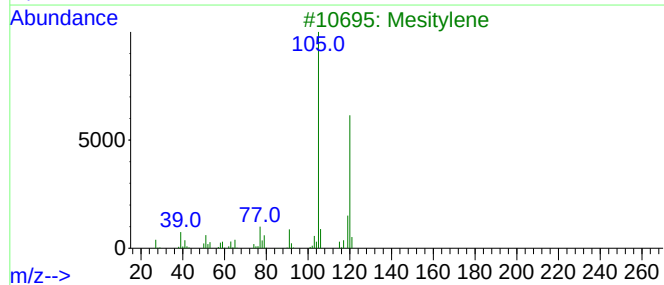
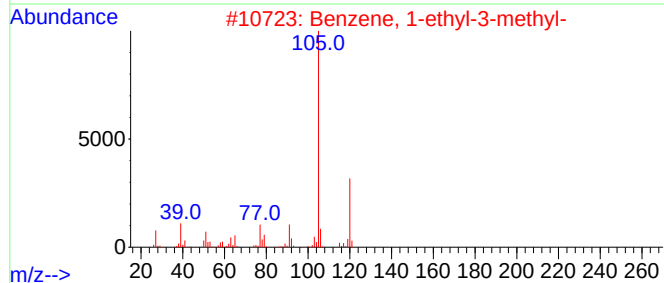
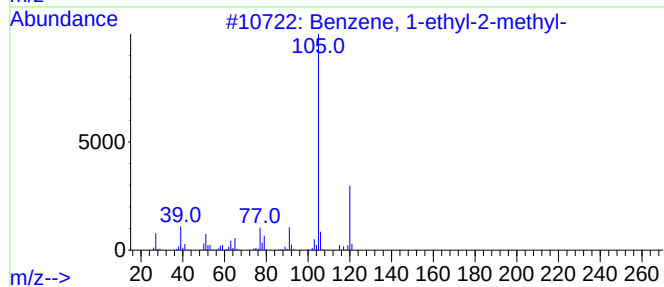
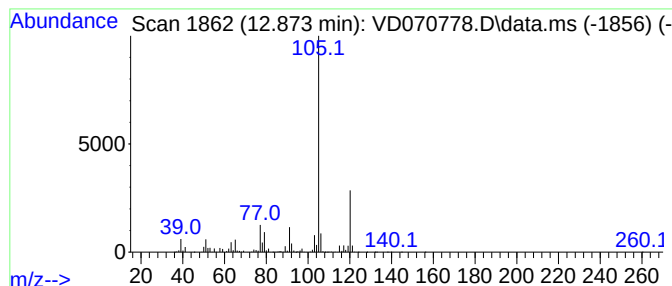
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D101521S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.873	78.22 ug/l	3237250	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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Misc : 5.01G/5.00ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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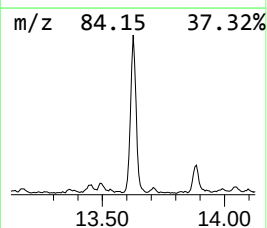
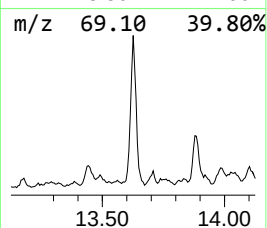
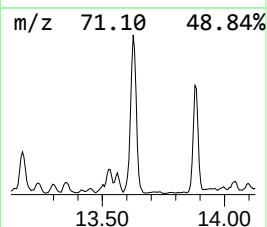
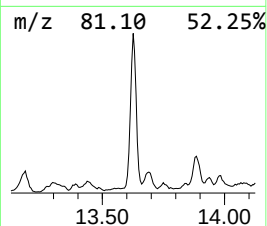
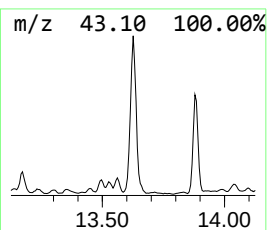
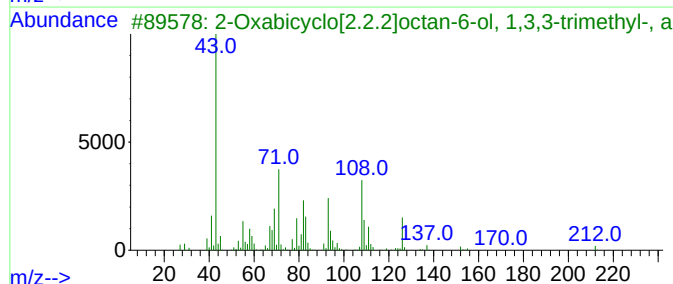
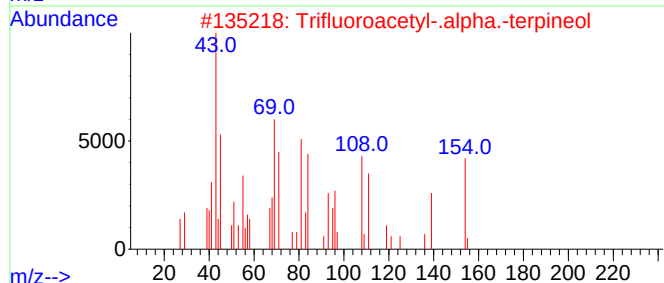
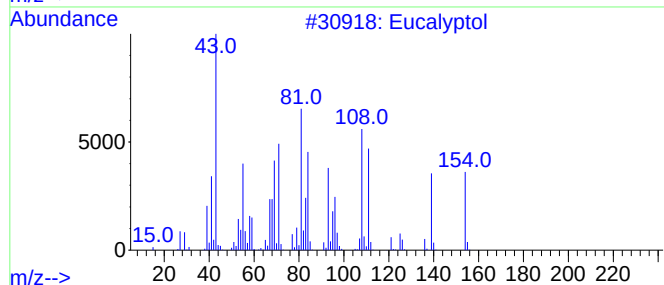
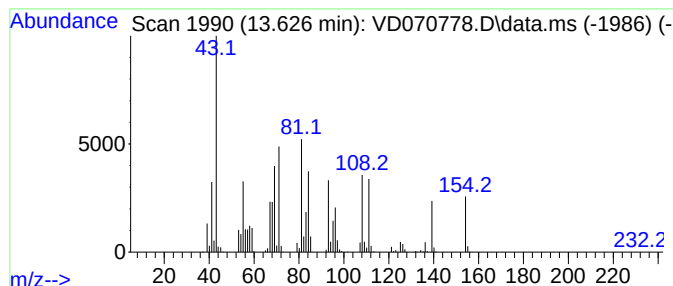
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D101521S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Eucalyptol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.626	153.45 ug/l	6350990	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eucalyptol	154	C10H18O	000470-82-6	97
2			Trifluoroacetyl-.alpha.-terpineol	250	C12H17F3O2	1000058-17-6	64
3			2-Oxabicyclo[2.2.2]octan-6-ol, 1...	212	C12H20O3	057709-95-2	38
4			5-Hepten-2-one, 6-methyl-	126	C8H14O	000110-93-0	30
5			3-Hepten-2-one, 3-propyl-	154	C10H18O	032064-69-0	25



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ALS Vial : 14 Sample Multiplier: 1

Instrument :
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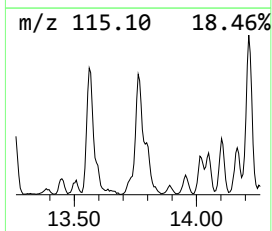
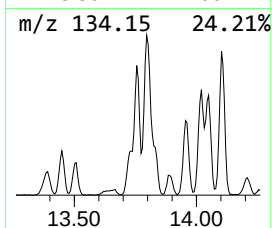
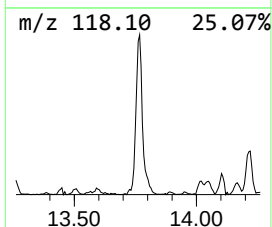
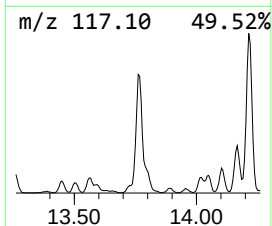
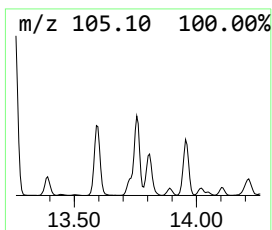
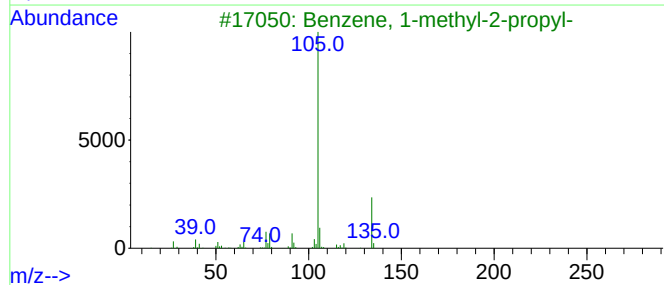
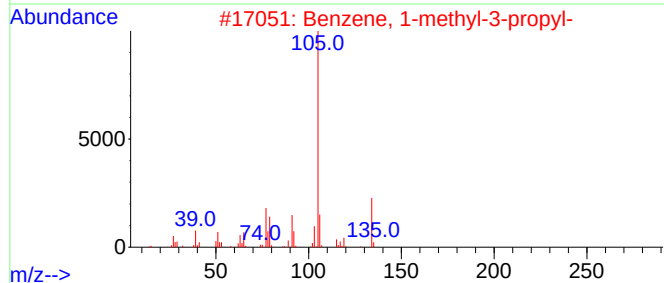
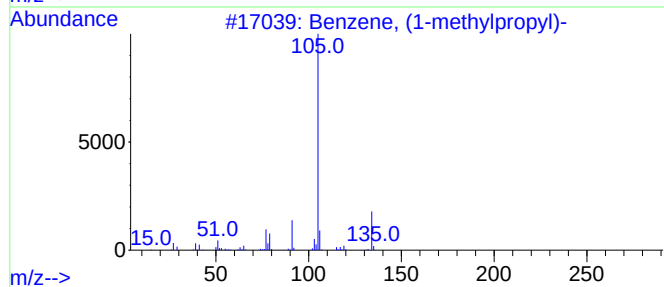
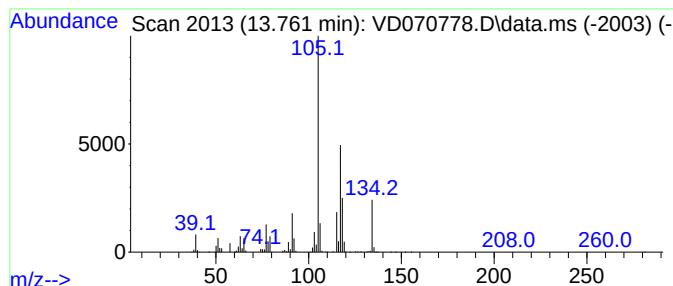
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D101521S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, (1-methylpropyl)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.761	99.65 ug/l	4124310	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	50
2			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	46
3			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	38
4			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	38
5			2-Indanol	134	C9H10O	004254-29-9	27



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ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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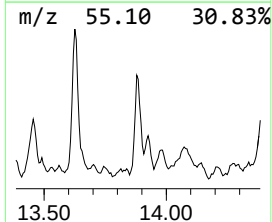
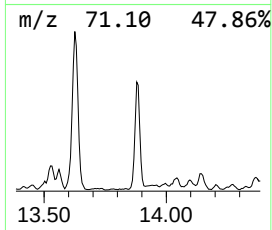
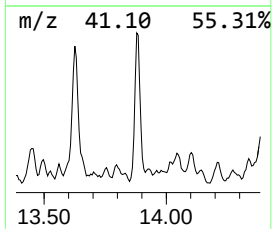
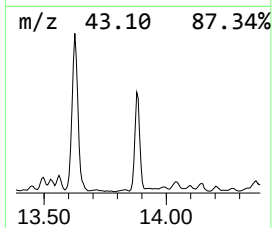
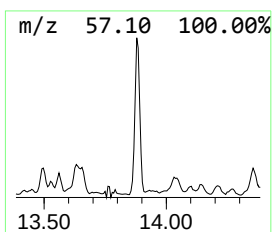
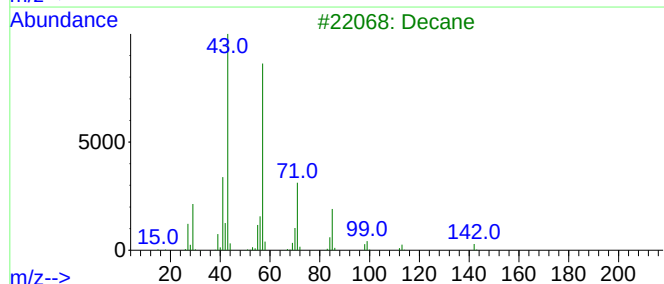
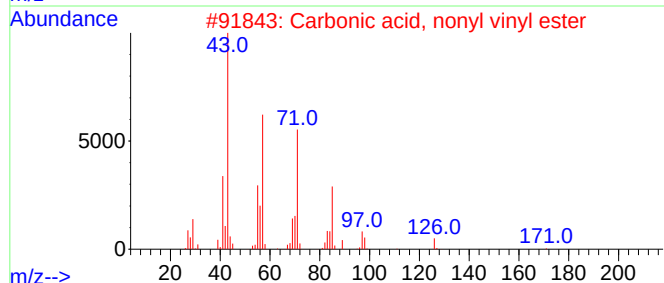
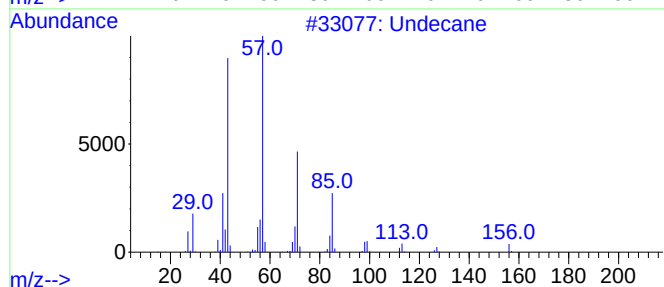
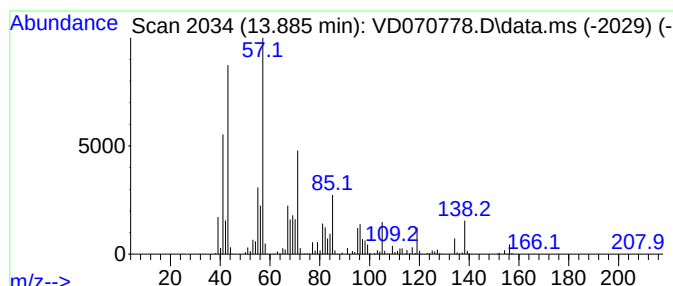
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Undecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.885	83.55 ug/l	3458160	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	60
2	Carbonic acid, nonyl vinyl ester			214	C12H22O3	1000383-25-6	43
3	Decane			142	C10H22	000124-18-5	43
4	Tetradecane			198	C14H30	000629-59-4	43
5	Hexadecane			226	C16H34	000544-76-3	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD102021\
Data File : VD070778.D
Acq On : 20 Oct 2021 16:20
Operator : VA/SY
Sample : M4284-03
Misc : 5.01G/5.00ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
30852

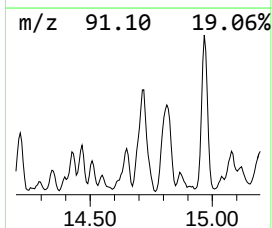
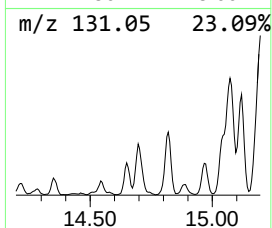
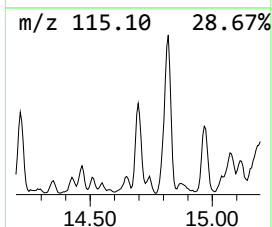
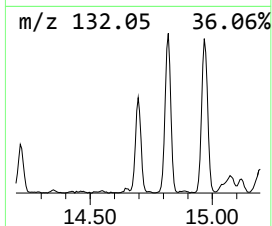
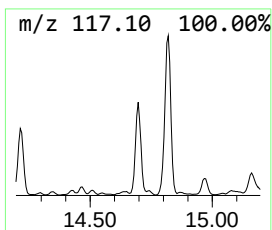
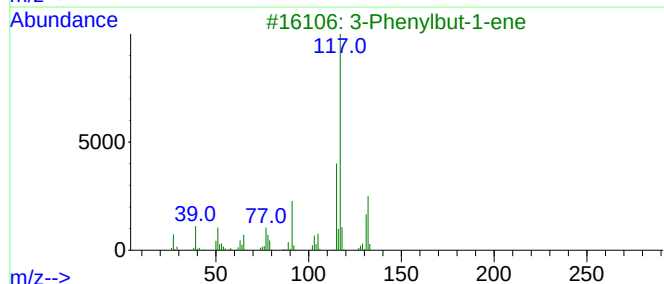
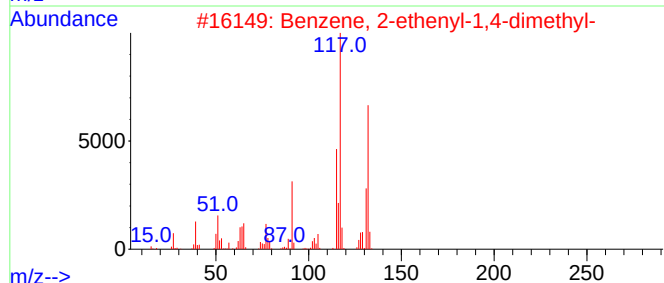
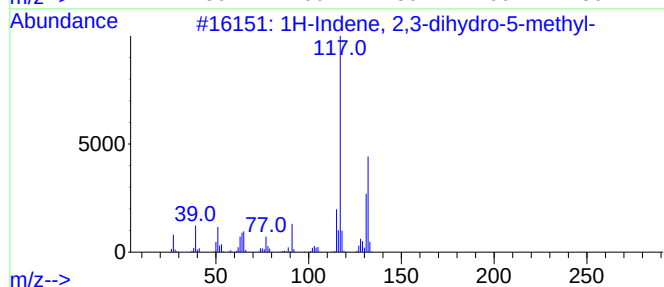
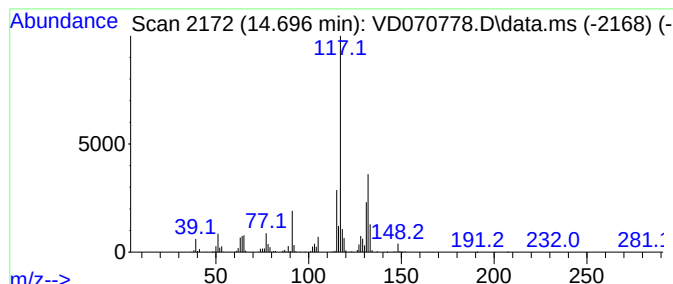
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.696	92.22 ug/l	3816930	1,4-Dichlorobenzene-d4	13.561

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
5		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD102021\
Data File : VD070778.D
Acq On : 20 Oct 2021 16:20
Operator : VA/SY
Sample : M4284-03
Misc : 5.01G/5.00ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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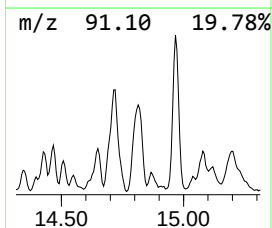
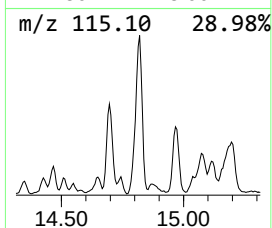
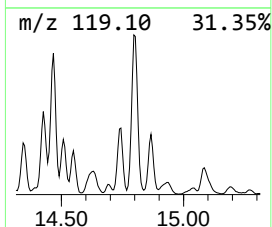
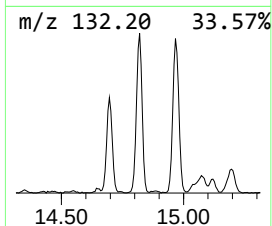
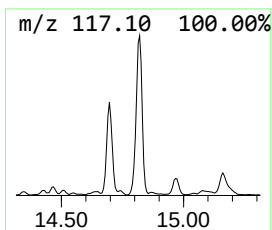
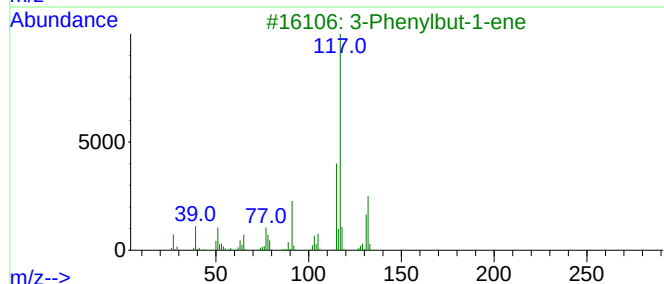
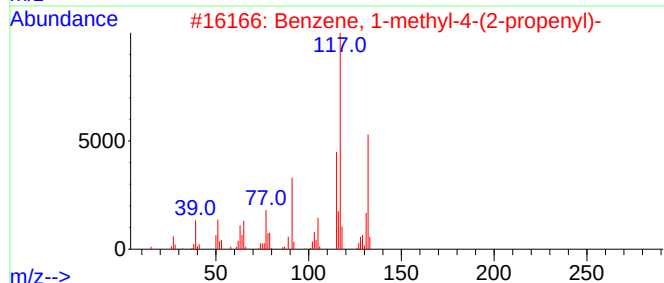
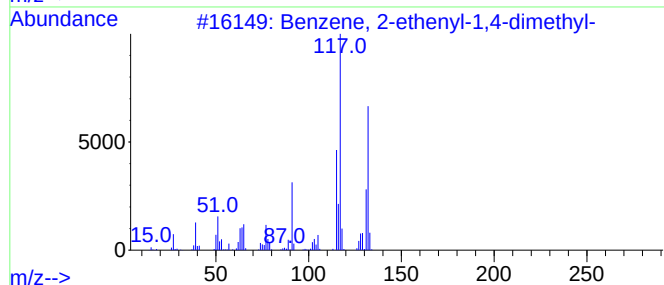
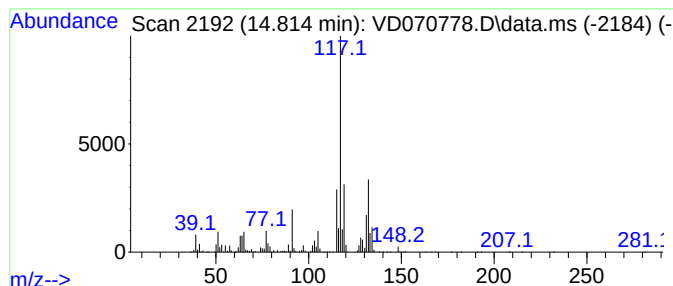
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.814	222.29 ug/l	9200150	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	90
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
4			Indan, 1-methyl-	132	C10H12	000767-58-8	76
5			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD102021\
Data File : VD070778.D
Acq On : 20 Oct 2021 16:20
Operator : VA/SY
Sample : M4284-03
Misc : 5.01G/5.00ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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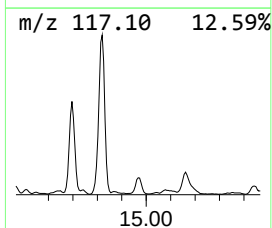
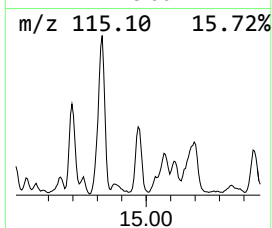
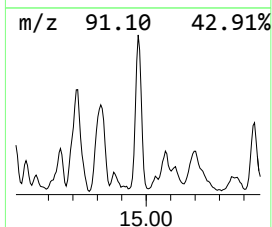
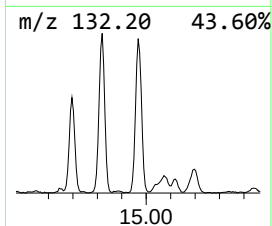
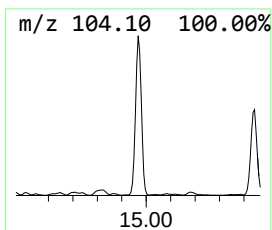
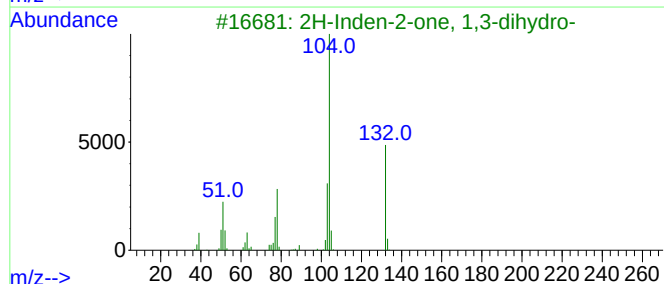
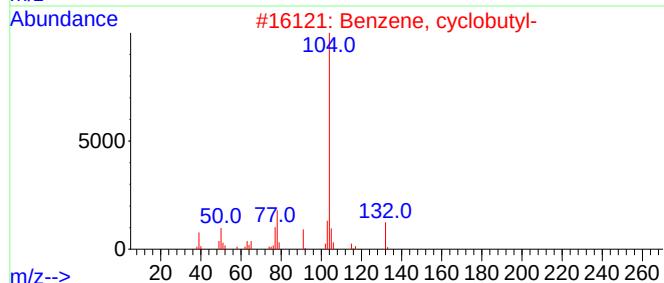
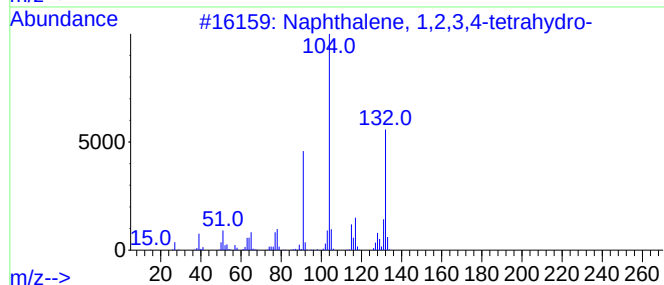
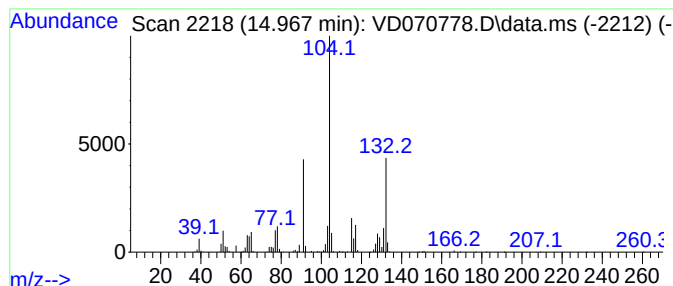
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.967	112.89 ug/l	4672330	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2			Benzene, cyclobutyl-	132	C10H12	004392-30-7	59
3			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	58
4			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	49
5			Benzene, 3-cyclohexen-1-yl-	158	C12H14	004994-16-5	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD102021\
Data File : VD070778.D
Acq On : 20 Oct 2021 16:20
Operator : VA/SY
Sample : M4284-03
Misc : 5.01G/5.00ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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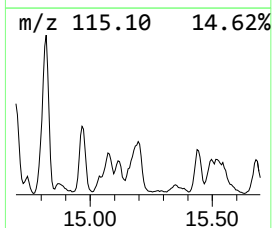
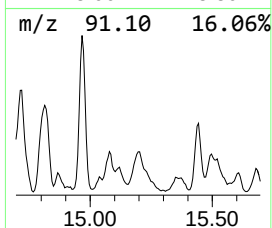
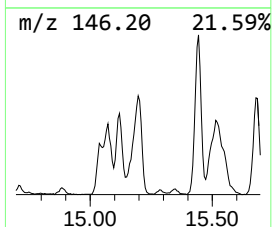
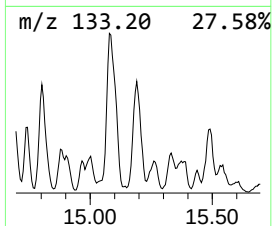
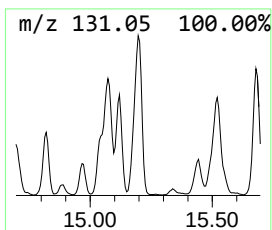
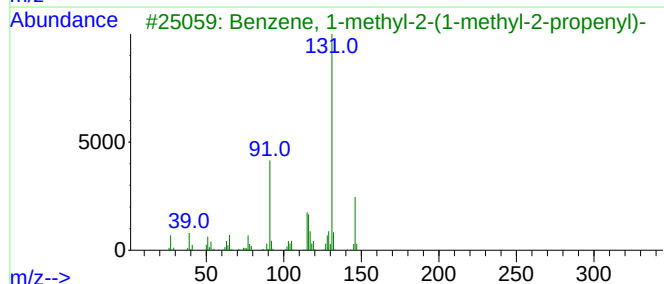
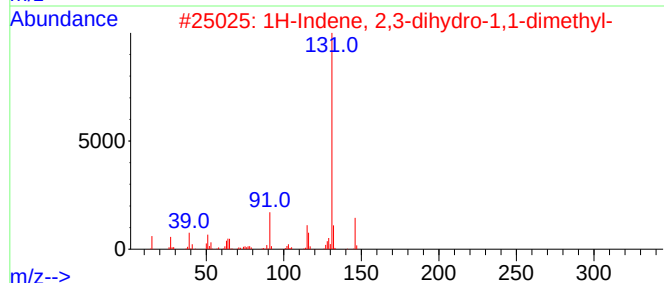
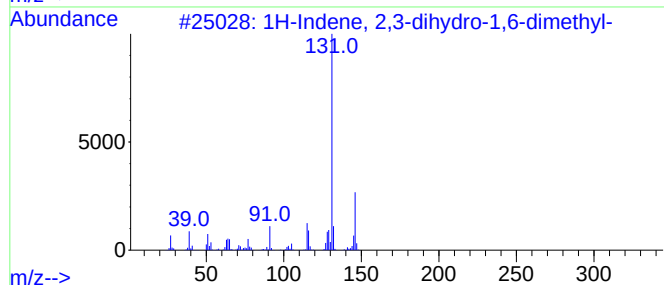
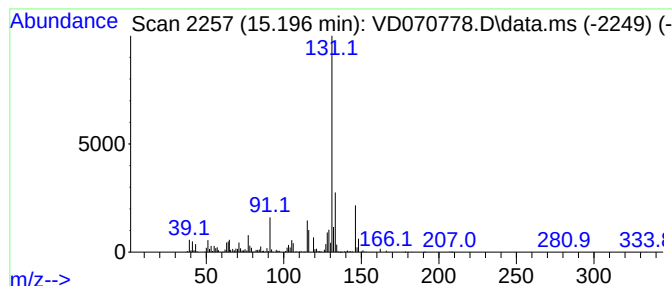
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.196	81.90 ug/l	3389760	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	93
2			1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	81
3			Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	81
4			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	81
5			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD102021\
Data File : VD070778.D
Acq On : 20 Oct 2021 16:20
Operator : VA/SY
Sample : M4284-03
Misc : 5.01G/5.00ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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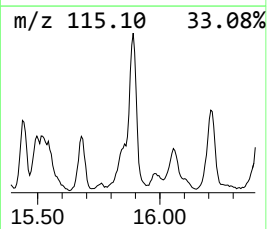
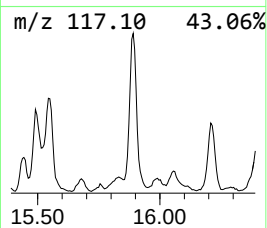
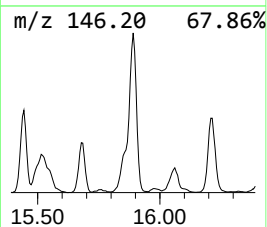
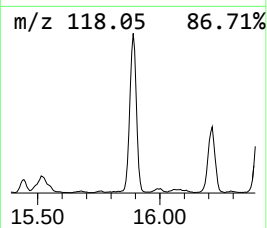
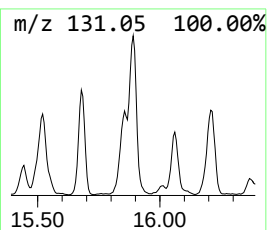
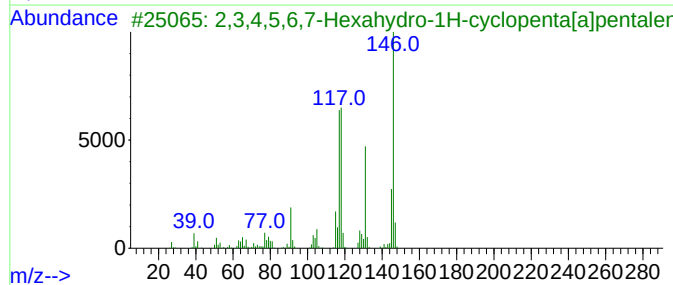
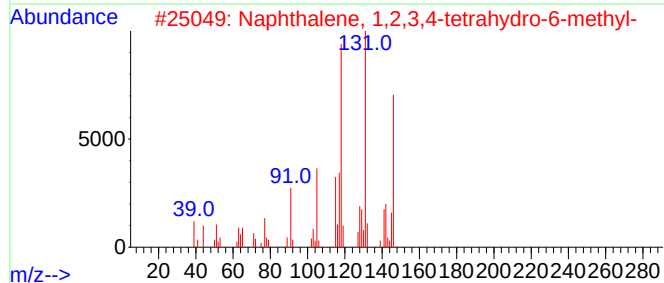
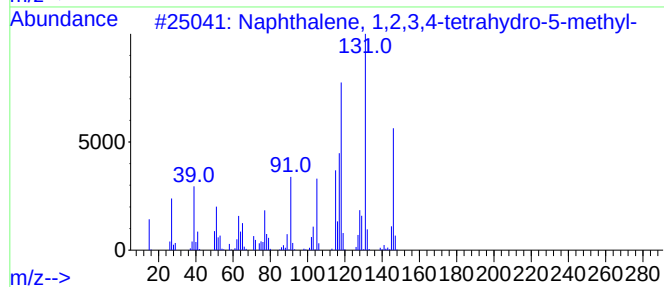
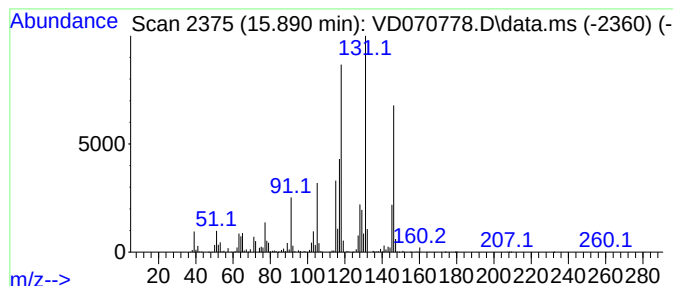
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.890	200.84 ug/l	8312290	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
3			2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1010189-31-0	74
4			1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	64
5			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD102021\
Data File : VD070778.D
Acq On : 20 Oct 2021 16:20
Operator : VA/SY
Sample : M4284-03
Misc : 5.01G/5.00ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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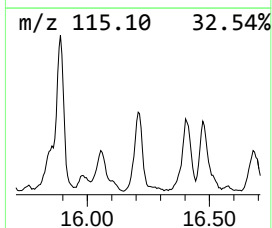
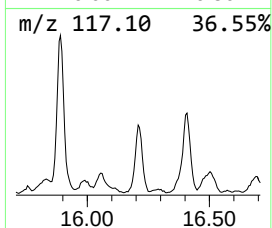
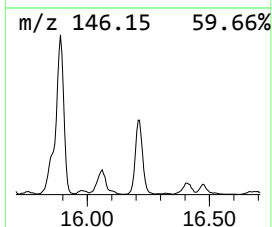
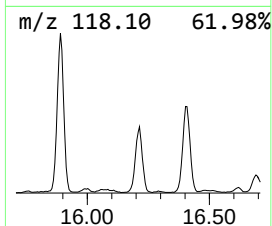
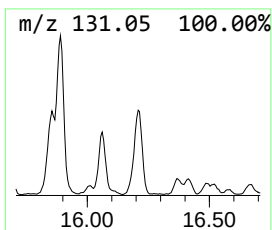
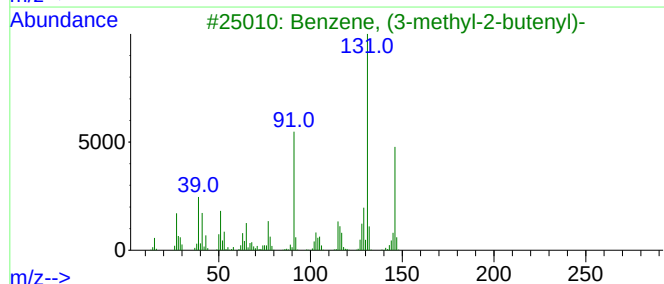
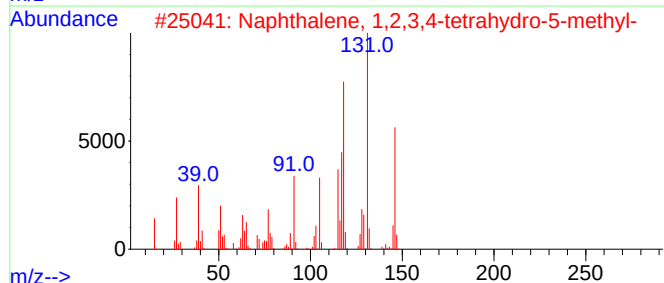
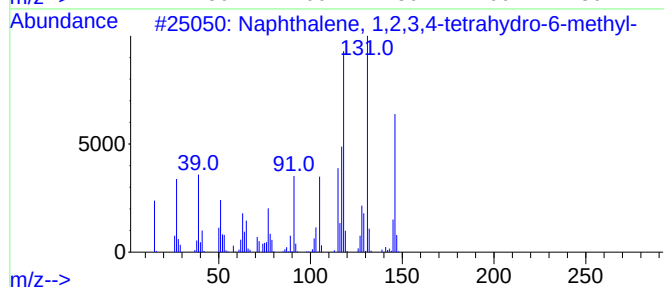
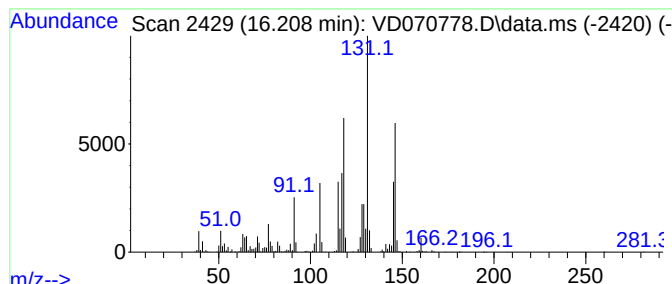
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.208	109.58 ug/l	4535360	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	93
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	90
3			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	58
4			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	58
5			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD102021\
Data File : VD070778.D
Acq On : 20 Oct 2021 16:20
Operator : VA/SY
Sample : M4284-03
Misc : 5.01G/5.00ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
30852

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D101521S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethy...	12.873	78.2	ug/l	3237250	4	13.561	2069410	50.0
Eucalyptol	13.626	153.4	ug/l	6350990	4	13.561	2069410	50.0
Benzene, (1-met...	13.761	99.7	ug/l	4124310	4	13.561	2069410	50.0
Undecane	13.885	83.5	ug/l	3458160	4	13.561	2069410	50.0
1H-Indene, 2,3-...	14.696	92.2	ug/l	3816930	4	13.561	2069410	50.0
Benzene, 2-ethe...	14.814	222.3	ug/l	9200150	4	13.561	2069410	50.0
Naphthalene, 1,...	14.967	112.9	ug/l	4672330	4	13.561	2069410	50.0
1H-Indene, 2,3-...	15.196	81.9	ug/l	3389760	4	13.561	2069410	50.0
Naphthalene, 1,...	15.890	200.8	ug/l	8312290	4	13.561	2069410	50.0
Naphthalene, 1,...	16.208	109.6	ug/l	4535360	4	13.561	2069410	50.0