

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041521\
 Data File : VY004437.D
 Acq On : 15 Apr 2021 13:25
 Operator : SY/MD
 Sample : M2047-05
 Misc : 6.66G/5ML/MSVOA_Y/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 130-SB3-(1.5-2)

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_Y\methods\82Y041221S.M
 Title : SW846 8260

Signal : TIC: VY004437.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.906	173	178	188	rVB2	28222	62514	2.44%	0.126%
2	3.259	229	236	252	rVB2	24270	66001	2.58%	0.133%
3	4.777	462	485	515	rBV2	107236	493923	19.27%	0.995%
4	5.247	548	562	576	rBV	77550	280818	10.96%	0.566%
5	5.759	634	646	657	rBV2	52535	168596	6.58%	0.340%
6	6.801	807	817	827	rBV2	73635	229355	8.95%	0.462%
7	7.533	930	937	947	rVB2	25173	64670	2.52%	0.130%
8	7.728	957	969	973	rBV	184157	461296	18.00%	0.929%
9	7.795	973	980	988	rVV2	383291	1025749	40.02%	2.066%
10	7.880	988	994	1004	rVB3	45854	122821	4.79%	0.247%
11	8.008	1006	1015	1022	rBV	123864	297582	11.61%	0.599%
12	8.149	1030	1038	1051	rVB	200296	440104	17.17%	0.887%
13	8.283	1051	1060	1064	rBV2	56520	144157	5.62%	0.290%
14	8.338	1064	1069	1078	rVV5	72152	234520	9.15%	0.472%
15	8.423	1078	1083	1094	rVB2	43013	102775	4.01%	0.207%
16	8.551	1094	1104	1118	rVB	86924	195689	7.64%	0.394%
17	8.697	1118	1128	1140	rBV	539337	1123740	43.84%	2.264%
18	9.185	1195	1208	1214	rBV	210708	532462	20.77%	1.073%
19	9.374	1227	1239	1245	rVV	38625	79294	3.09%	0.160%
20	9.465	1246	1254	1261	rVB	27268	61453	2.40%	0.124%
21	9.618	1273	1279	1289	rVB2	47082	94285	3.68%	0.190%
22	9.758	1298	1302	1307	rVB3	42610	78532	3.06%	0.158%
23	9.825	1307	1313	1316	rBV	91372	161037	6.28%	0.324%
24	9.959	1326	1335	1346	rBV	105612	246976	9.64%	0.498%
25	10.179	1365	1371	1386	rVV	887970	1636266	63.84%	3.296%
26	10.374	1395	1403	1409	rVB3	143904	292411	11.41%	0.589%
27	10.526	1423	1428	1436	rVB2	41135	68631	2.68%	0.138%
28	10.612	1436	1442	1454	rBV6	54535	124494	4.86%	0.251%
29	10.807	1463	1474	1479	rBV2	37297	73590	2.87%	0.148%
30	10.904	1479	1490	1496	rBV3	43505	122907	4.80%	0.248%
31	11.038	1508	1512	1516	rVB	58764	85290	3.33%	0.172%
32	11.093	1516	1521	1526	rVB	55058	91645	3.58%	0.185%
33	11.203	1534	1539	1543	rBV2	35797	61644	2.41%	0.124%
34	11.282	1543	1552	1563	rVB3	139311	338790	13.22%	0.682%
35	11.380	1563	1568	1578	rBV	121354	211640	8.26%	0.426%
36	11.489	1580	1586	1596	rBV	733728	1225364	47.81%	2.468%

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Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
 Title : SW846 8260

37	11.593	1597	1603	1611	rBV	468933	739817	28.86%	1.490%
38	11.709	1611	1622	1631	rBV2	377417	739276	28.84%	1.489%
39	12.008	1664	1671	1673	rBV5	71105	137790	5.38%	0.278%
40	12.124	1686	1690	1694	rVB	71023	90393	3.53%	0.182%
41	12.203	1698	1703	1706	rVV2	54648	93394	3.64%	0.188%
42	12.245	1706	1710	1714	rVB3	52427	95923	3.74%	0.193%
43	12.331	1719	1724	1728	rVV	100294	175472	6.85%	0.353%
44	12.386	1728	1733	1734	rVV5	67182	123546	4.82%	0.249%
45	12.416	1735	1738	1740	rVV	107909	168964	6.59%	0.340%
46	12.441	1740	1742	1744	rVV	110514	127383	4.97%	0.257%
47	12.489	1744	1750	1760	rVB2	1379733	2563037	100.00%	5.163%
48	12.672	1772	1780	1785	rBV	371706	579603	22.61%	1.168%
49	12.733	1785	1790	1792	rVV2	66652	106033	4.14%	0.214%
50	12.770	1792	1796	1799	rVV	304691	504433	19.68%	1.016%
51	12.806	1799	1802	1808	rVV2	450002	688581	26.87%	1.387%
52	12.971	1821	1829	1837	rBV	488597	914521	35.68%	1.842%
53	13.038	1837	1840	1846	rVB	94855	124616	4.86%	0.251%
54	13.123	1847	1854	1859	rBV	1110420	1712938	66.83%	3.451%
55	13.166	1859	1861	1866	rVB3	49845	67564	2.64%	0.136%
56	13.233	1866	1872	1873	rBV2	75992	122047	4.76%	0.246%
57	13.318	1879	1886	1890	rBV3	127318	223828	8.73%	0.451%
58	13.367	1890	1894	1897	rBV2	135481	207903	8.11%	0.419%
59	13.428	1897	1904	1907	rVV	764688	1387026	54.12%	2.794%
60	13.459	1907	1909	1913	rVV	594289	683906	26.68%	1.378%
61	13.501	1913	1916	1923	rVB2	138029	204091	7.96%	0.411%
62	13.593	1926	1931	1933	rBV	268865	383958	14.98%	0.773%
63	13.629	1933	1937	1940	rVV2	644887	969085	37.81%	1.952%
64	13.672	1940	1944	1952	rVB3	460155	997603	38.92%	2.010%
65	13.751	1952	1957	1962	rBV2	521915	741229	28.92%	1.493%
66	13.824	1962	1969	1974	rBV	250204	388118	15.14%	0.782%
67	13.885	1974	1979	1981	rBV	295772	415981	16.23%	0.838%
68	13.916	1981	1984	1988	rVB	313368	427143	16.67%	0.860%
69	13.971	1988	1993	1998	rVB	1087906	1525235	59.51%	3.072%
70	14.032	1998	2003	2006	rBV2	136706	203645	7.95%	0.410%
71	14.080	2006	2011	2016	rVB	686320	1051062	41.01%	2.117%
72	14.148	2016	2022	2027	rBV6	87357	178018	6.95%	0.359%
73	14.215	2027	2033	2037	rBV3	224906	465326	18.16%	0.937%
74	14.294	2042	2046	2050	rVV	642253	1193537	46.57%	2.404%
75	14.337	2050	2053	2057	rVV	512168	718347	28.03%	1.447%
76	14.385	2057	2061	2064	rVV	342379	483687	18.87%	0.974%

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77	14.422	2064	2067	2070	rVV2	165975	223518	8.72%	0.450%
78	14.452	2070	2072	2075	rVV	74708	102206	3.99%	0.206%
79	14.519	2079	2083	2087	rVV	198247	278471	10.86%	0.561%
80	14.568	2087	2091	2095	rVV2	685016	1157150	45.15%	2.331%
81	14.611	2095	2098	2103	rVB2	604747	829161	32.35%	1.670%
82	14.684	2103	2110	2115	rBV3	1075677	2418329	94.35%	4.872%
83	14.739	2115	2119	2125	rVB2	348696	612391	23.89%	1.234%
84	14.836	2131	2135	2138	rVB	85849	101336	3.95%	0.204%
85	14.910	2143	2147	2148	rBV2	128761	174603	6.81%	0.352%
86	14.946	2148	2153	2157	rBV3	463898	772330	30.13%	1.556%
87	15.056	2165	2171	2177	rVB2	430593	861505	33.61%	1.735%
88	15.135	2181	2184	2190	rVB4	98230	169608	6.62%	0.342%
89	15.239	2190	2201	2207	rBV	904107	1828860	71.36%	3.684%
90	15.349	2213	2219	2223	rBV	260059	445724	17.39%	0.898%
91	15.397	2223	2227	2230	rBV	114379	173254	6.76%	0.349%
92	15.434	2230	2233	2242	rVB3	307431	504139	19.67%	1.016%
93	15.525	2242	2248	2255	rBV2	310134	640569	24.99%	1.290%
94	15.605	2258	2261	2265	rVB	68781	100025	3.90%	0.201%
95	15.690	2266	2275	2285	rBV3	169475	579828	22.62%	1.168%
96	15.818	2291	2296	2301	rVV2	106253	178153	6.95%	0.359%
97	15.885	2301	2307	2313	rVB3	143765	311190	12.14%	0.627%
98	16.025	2325	2330	2336	rBV4	59045	136083	5.31%	0.274%
99	16.293	2367	2374	2381	rBV	1122885	2151703	83.95%	4.334%
100	16.489	2399	2406	2418	rVB	497873	1066847	41.62%	2.149%

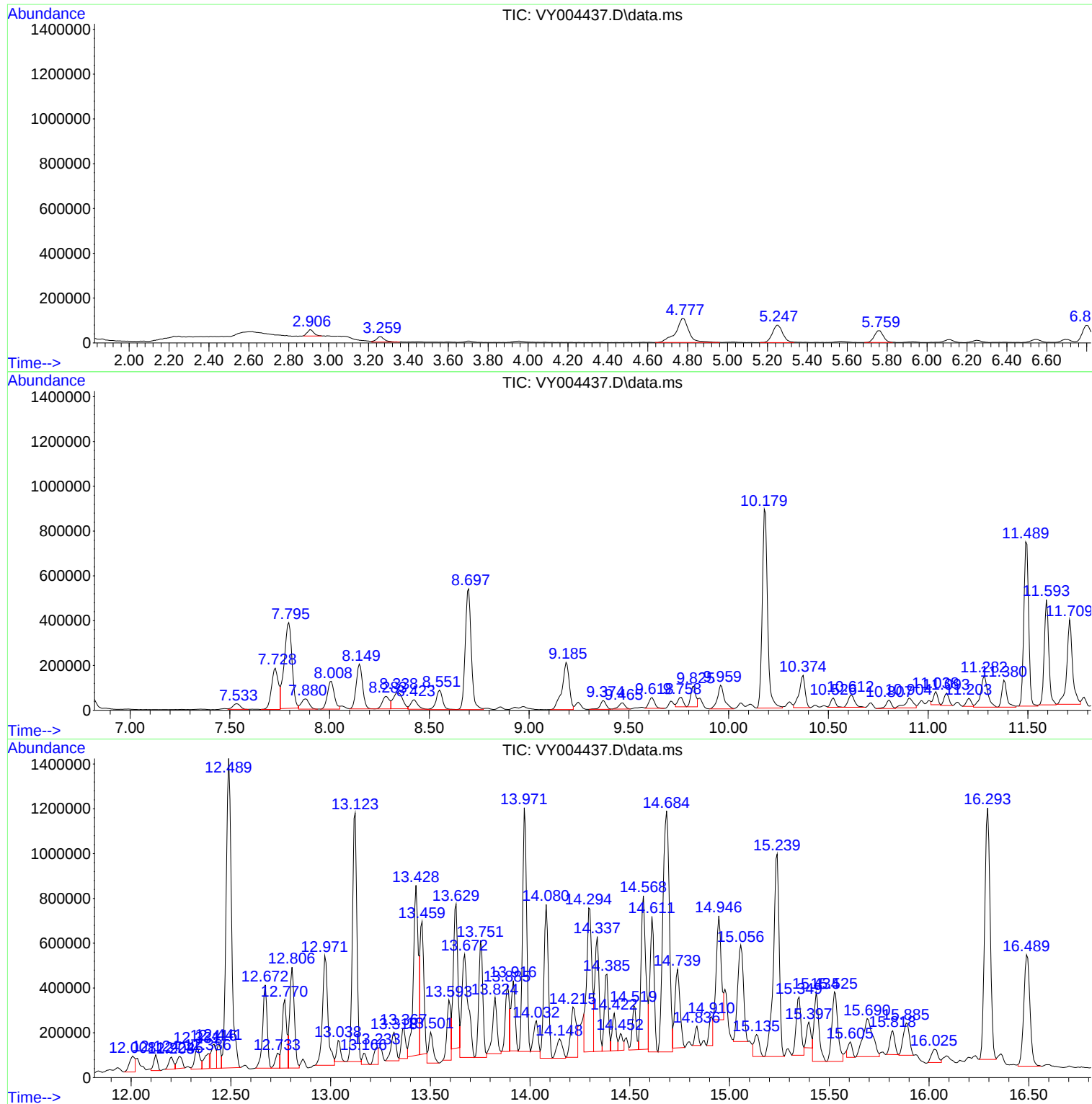
Sum of corrected areas: 49642063

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ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
130-SB3-(1.5-2)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_Y
ClientSampleId :
130-SB3-(1.5-2)

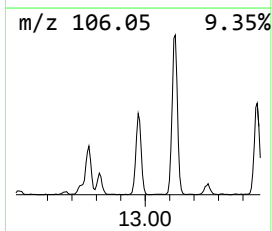
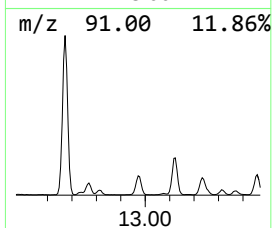
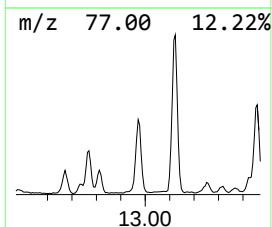
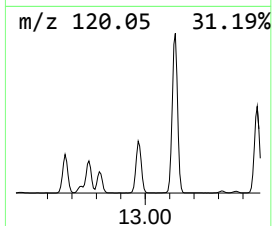
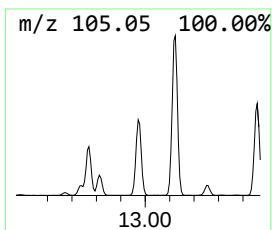
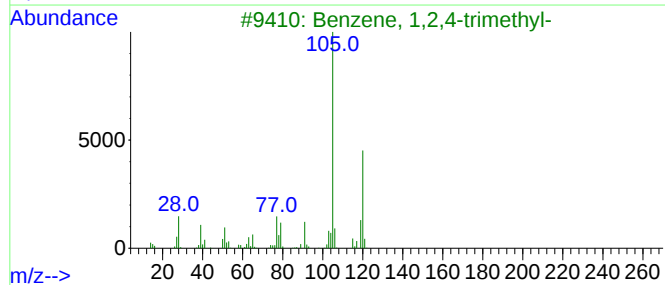
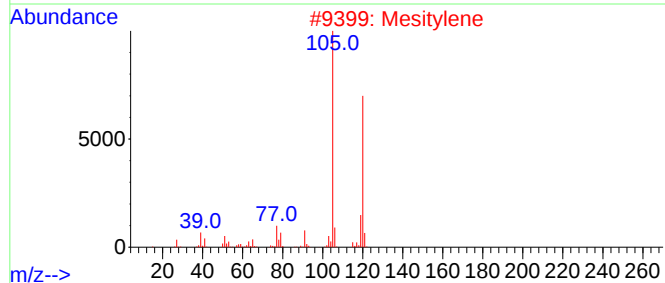
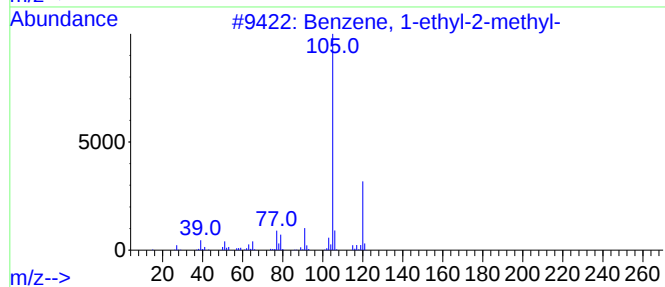
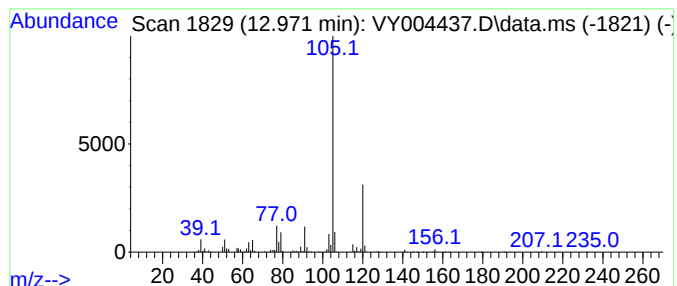
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.971	32.97 ug/l	914521	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Mesitylene	120	C9H12	000108-67-8	91
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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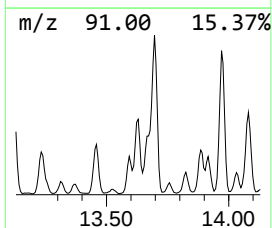
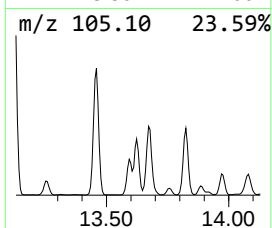
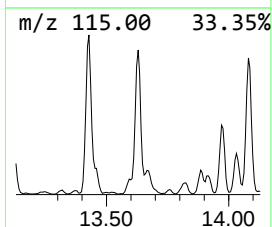
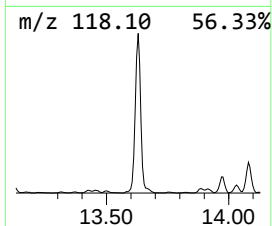
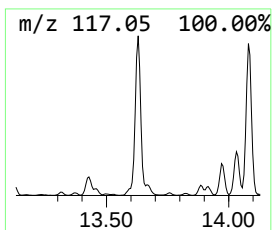
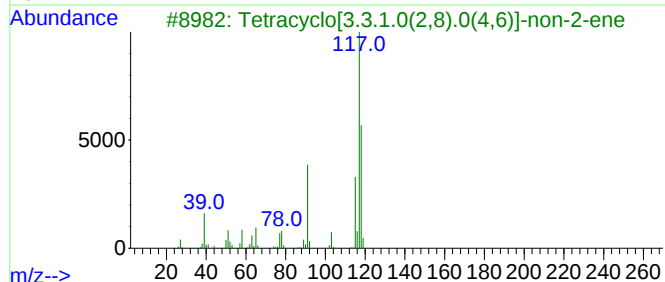
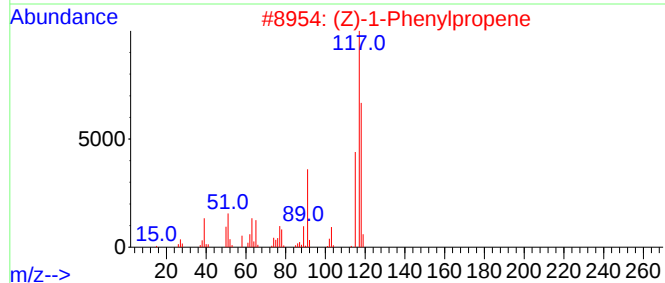
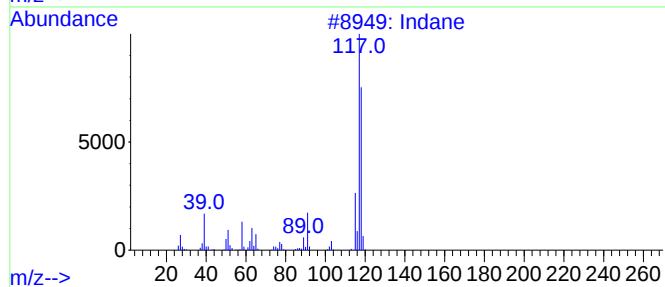
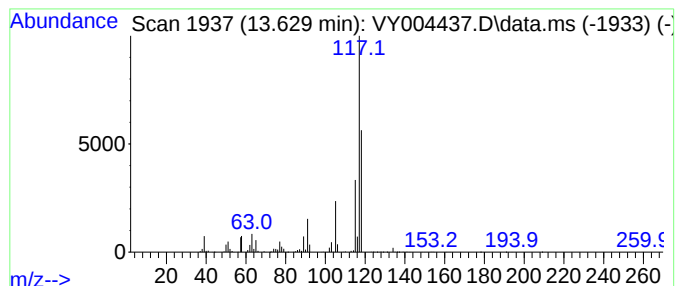
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Indane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.629	34.93 ug/l	969085	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	76
2			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	62
3			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	59
4			Benzene, 2-propenyl-	118	C9H10	000300-57-2	59
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	58



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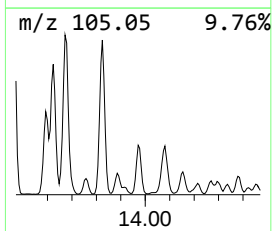
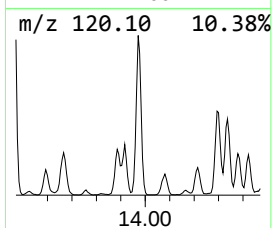
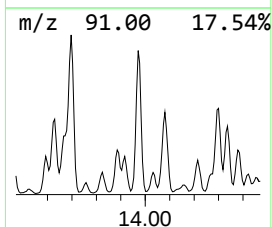
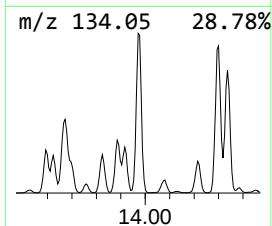
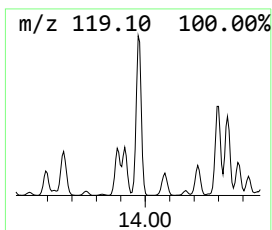
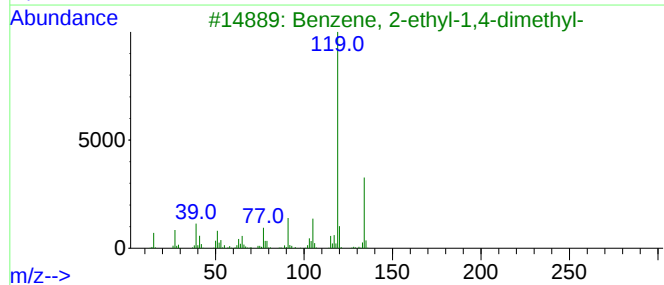
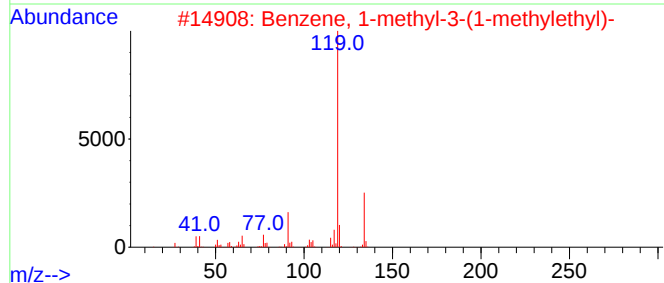
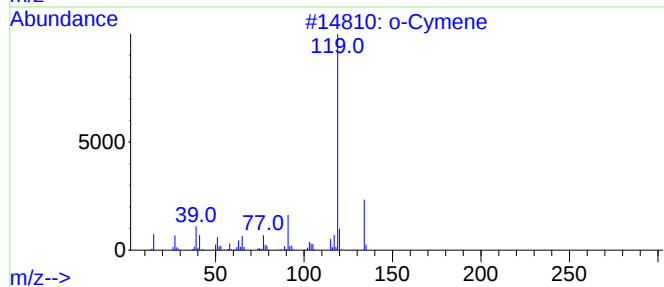
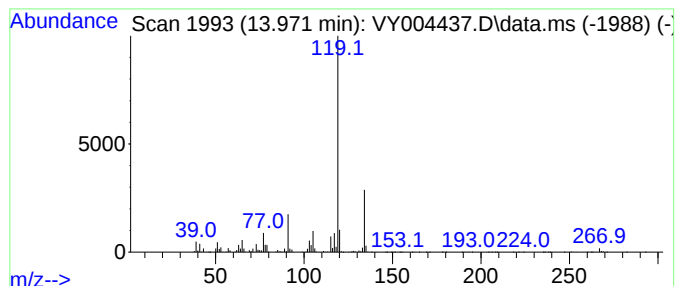
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Quant Title : SW846 8260

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

Peak Number 3 o-Cymene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.971	54.98 ug/l	1525240	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	95
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5			p-Cymene	134	C10H14	000099-87-6	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041521\
Data File : VY004437.D
Acq On : 15 Apr 2021 13:25
Operator : SY/MD
Sample : M2047-05
Misc : 6.66G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
130-SB3-(1.5-2)

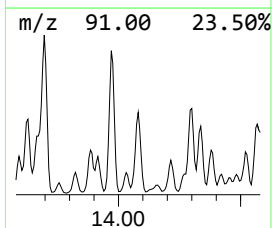
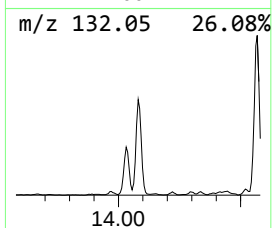
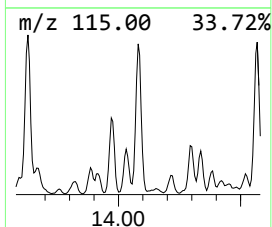
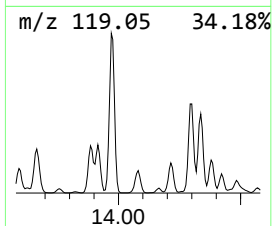
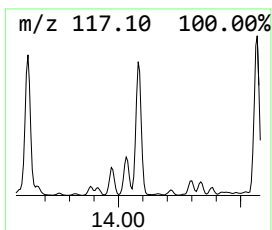
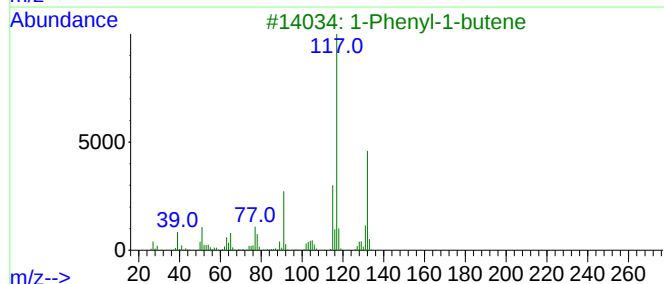
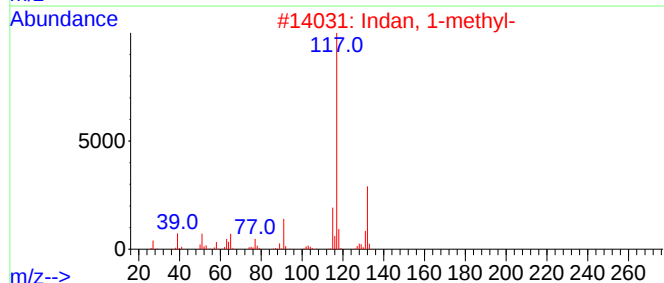
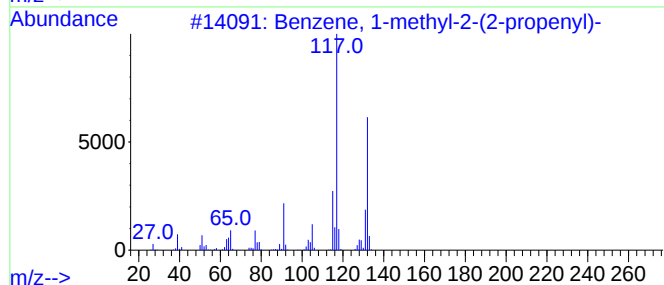
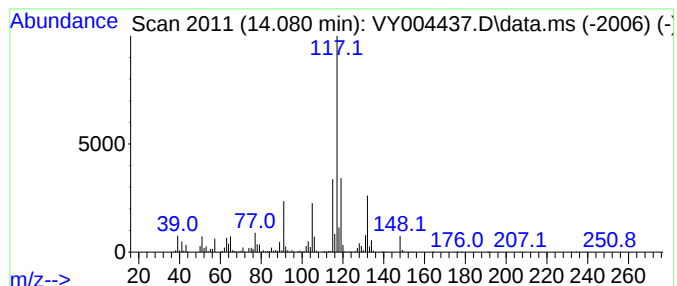
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-methyl-2-(2-prop... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.081	37.89 ug/l	1051060	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	76
2			Indan, 1-methyl-	132	C10H12	000767-58-8	70
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	70
4			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	70
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041521\
Data File : VY004437.D
Acq On : 15 Apr 2021 13:25
Operator : SY/MD
Sample : M2047-05
Misc : 6.66G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
130-SB3-(1.5-2)

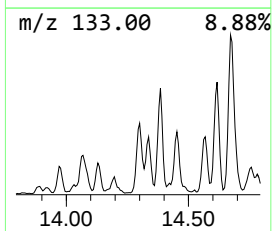
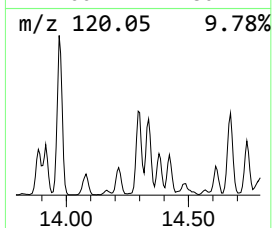
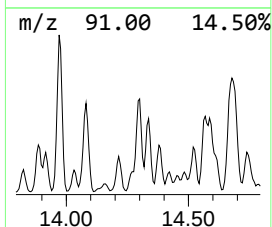
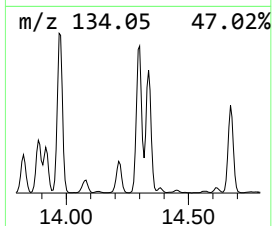
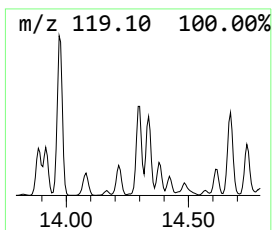
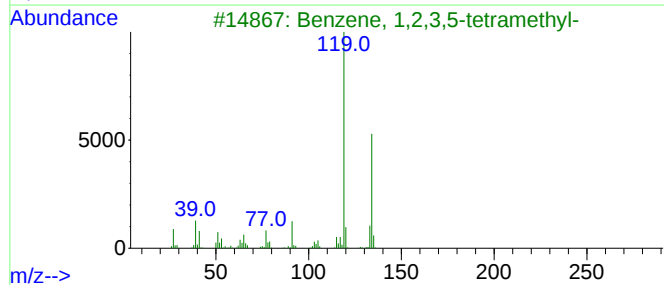
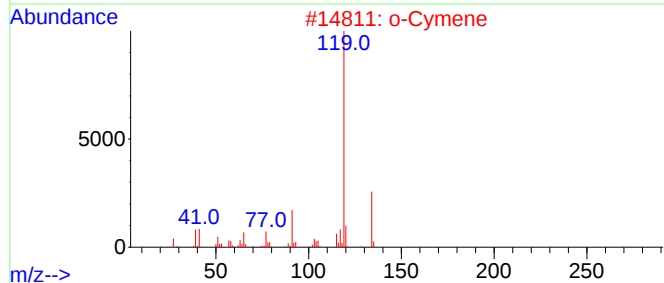
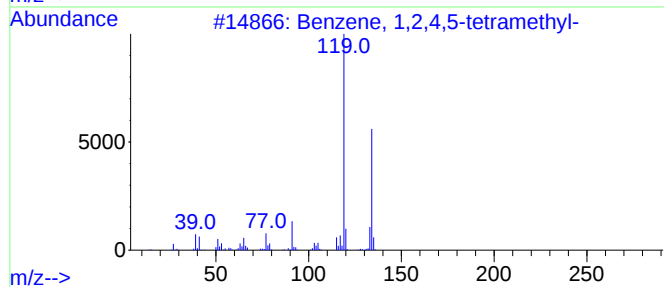
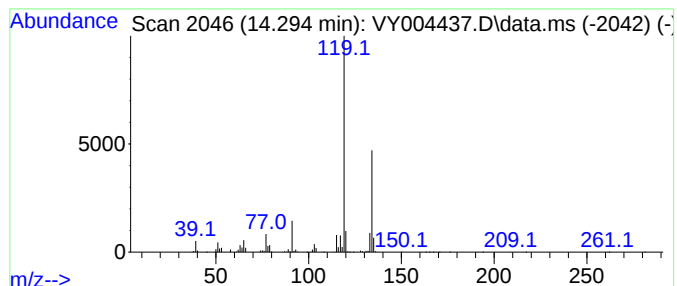
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.294	43.02 ug/l	1193540	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2			o-Cymene	134	C10H14	000527-84-4	95
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041521\
Data File : VY004437.D
Acq On : 15 Apr 2021 13:25
Operator : SY/MD
Sample : M2047-05
Misc : 6.66G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
130-SB3-(1.5-2)

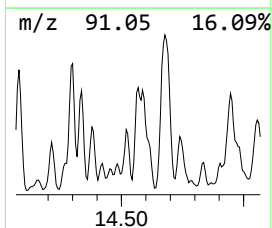
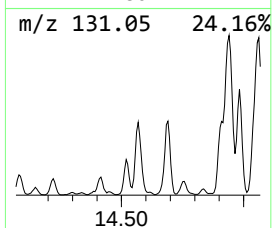
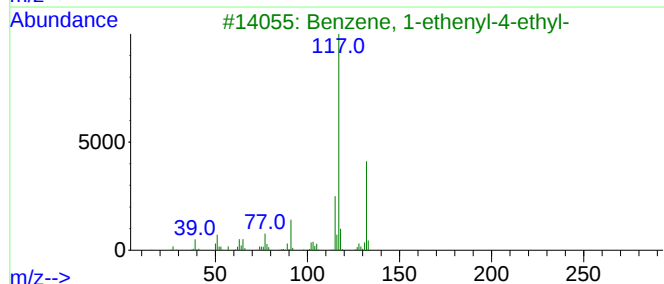
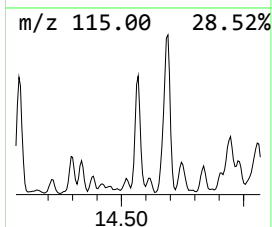
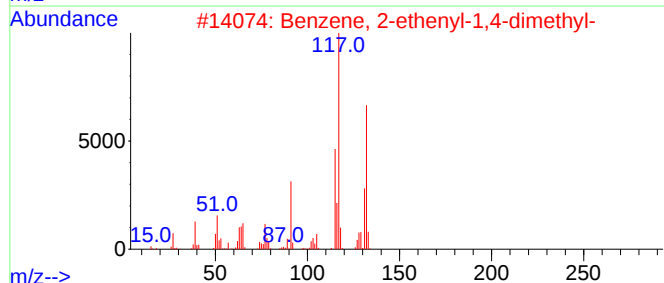
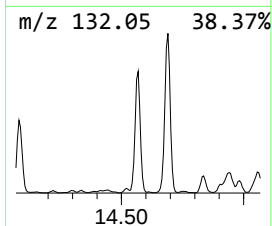
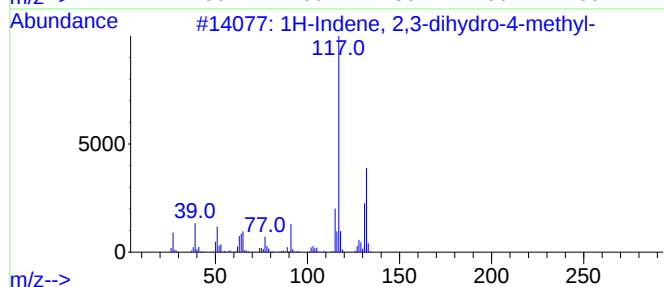
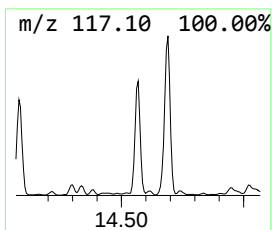
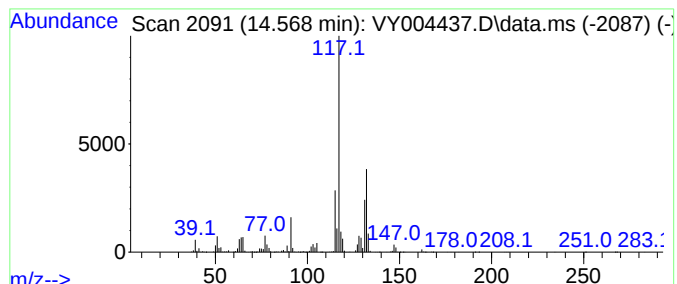
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.568	41.71 ug/l	1157150	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-			132	C10H12	000824-22-6	94
2	Benzene, 2-ethenyl-1,4-dimethyl-			132	C10H12	002039-89-6	93
3	Benzene, 1-ethenyl-4-ethyl-			132	C10H12	003454-07-7	91
4	3-Phenylbut-1-ene			132	C10H12	000934-10-1	91
5	Indan, 1-methyl-			132	C10H12	000767-58-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041521\
Data File : VY004437.D
Acq On : 15 Apr 2021 13:25
Operator : SY/MD
Sample : M2047-05
Misc : 6.66G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
130-SB3-(1.5-2)

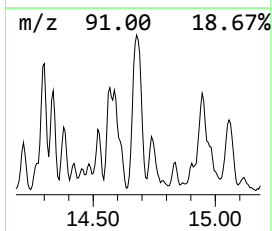
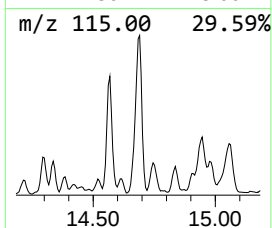
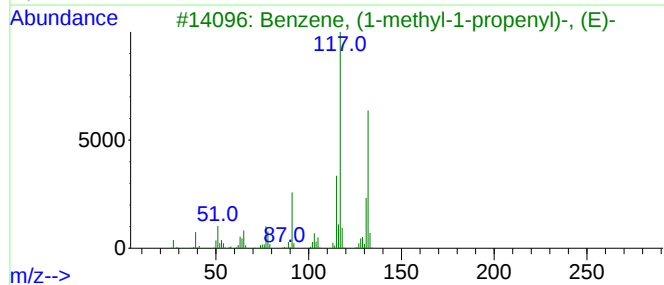
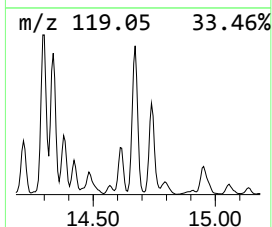
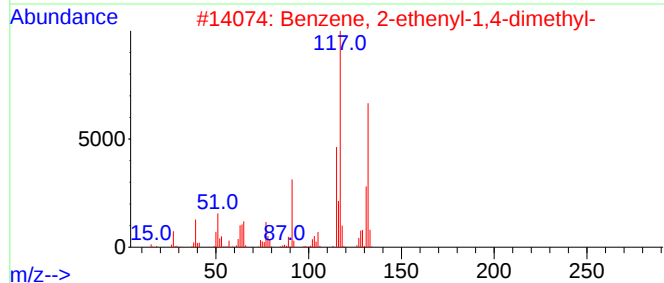
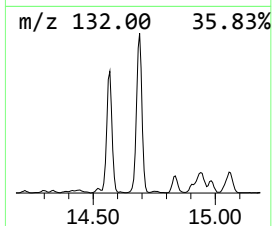
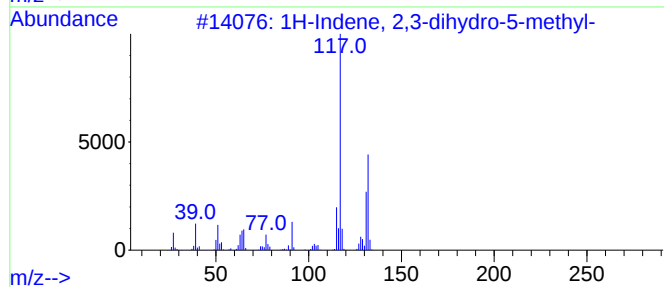
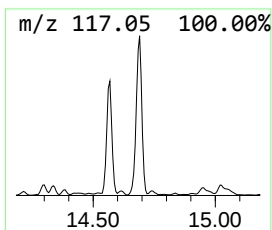
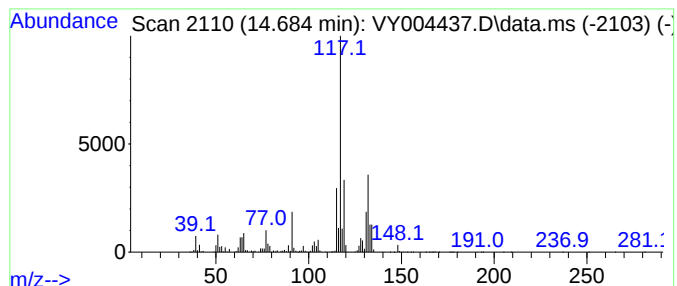
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	87.18 ug/l	2418330	1,4-Dichlorobenzene-d4	13.428

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	92
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041521\
Data File : VY004437.D
Acq On : 15 Apr 2021 13:25
Operator : SY/MD
Sample : M2047-05
Misc : 6.66G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
130-SB3-(1.5-2)

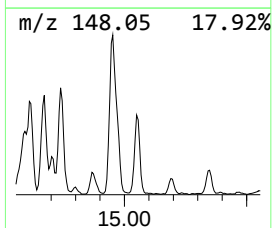
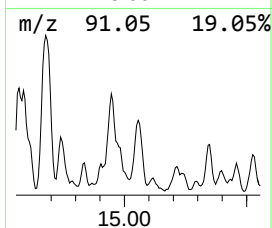
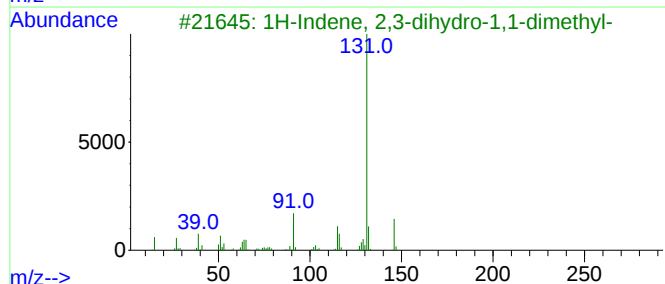
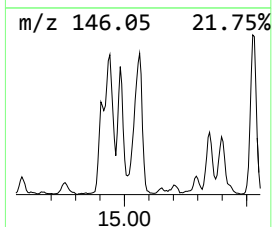
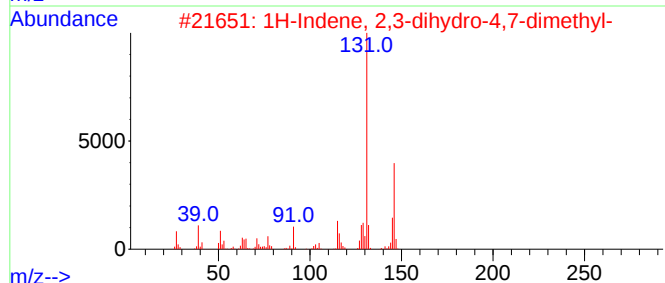
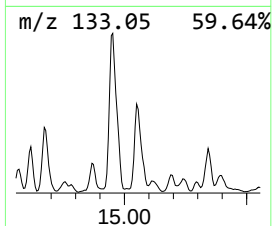
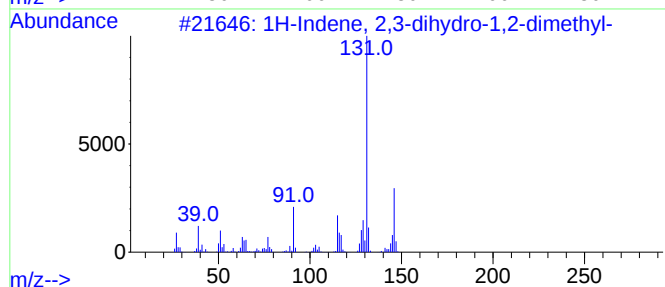
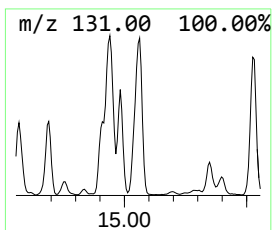
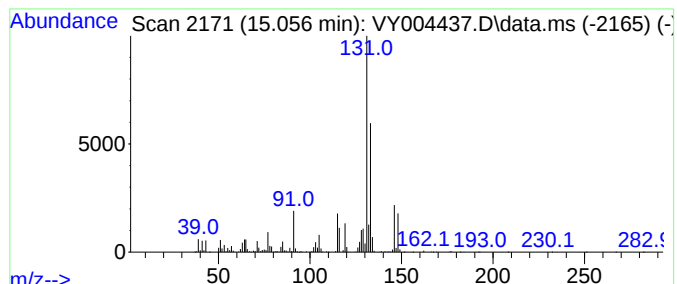
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.056	31.06 ug/l	861505	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	64
2			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	64
3			1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	64
4			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	64
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041521\
Data File : VY004437.D
Acq On : 15 Apr 2021 13:25
Operator : SY/MD
Sample : M2047-05
Misc : 6.66G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
130-SB3-(1.5-2)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.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
Benzene, 1-ethy...	12.971	33.0	ug/l	914521	4	13.428	1387030	50.0
Indane	13.629	34.9	ug/l	969085	4	13.428	1387030	50.0
o-Cymene	13.971	55.0	ug/l	1525240	4	13.428	1387030	50.0
Benzene, 1-meth...	14.081	37.9	ug/l	1051060	4	13.428	1387030	50.0
Benzene, 1,2,4,...	14.294	43.0	ug/l	1193540	4	13.428	1387030	50.0
1H-Indene, 2,3-...	14.568	41.7	ug/l	1157150	4	13.428	1387030	50.0
1H-Indene, 2,3-...	14.684	87.2	ug/l	2418330	4	13.428	1387030	50.0
1H-Indene, 2,3-...	15.056	31.1	ug/l	861505	4	13.428	1387030	50.0