

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY042321\
 Data File : VY004603.D
 Acq On : 23 Apr 2021 14:25
 Operator : SY/MD
 Sample : M2142-01
 Misc : 7.05G/5ML/MSVOA_Y/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 EO-04-042221

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: VY004603.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.174	47	58	59	rBV3	5692	14560	1.06%	0.103%
2	2.564	113	122	123	rBV2	11348	26953	1.96%	0.191%
3	4.735	469	478	492	rVB2	10226	32097	2.34%	0.227%
4	7.734	960	970	975	rBV	104799	249015	18.13%	1.764%
5	7.801	975	981	991	rVB	180712	401521	29.23%	2.845%
6	8.155	1028	1039	1050	rBV	115628	253480	18.45%	1.796%
7	8.697	1120	1128	1138	rBV	286143	556820	40.53%	3.946%
8	10.185	1364	1372	1381	rBV	469027	805657	58.65%	5.709%
9	10.374	1395	1403	1411	rVB3	12637	24051	1.75%	0.170%
10	11.045	1509	1513	1517	rVV	14529	22878	1.67%	0.162%
11	11.093	1517	1521	1527	rVB2	15100	25399	1.85%	0.180%
12	11.288	1543	1553	1563	rVB5	26642	72375	5.27%	0.513%
13	11.386	1563	1569	1573	rBV	22037	35220	2.56%	0.250%
14	11.496	1580	1587	1597	rVB	420564	670512	48.81%	4.751%
15	11.593	1597	1603	1607	rBV2	14456	28840	2.10%	0.204%
16	11.715	1612	1623	1631	rVV3	102491	247905	18.05%	1.757%
17	11.782	1631	1634	1640	rVB2	20663	33100	2.41%	0.235%
18	12.032	1663	1675	1683	rBV3	72073	196897	14.33%	1.395%
19	12.124	1687	1690	1694	rVV	22451	30811	2.24%	0.218%
20	12.203	1699	1703	1706	rVV4	23127	41014	2.99%	0.291%
21	12.252	1706	1711	1720	rVV3	41706	95425	6.95%	0.676%
22	12.337	1720	1725	1729	rVV3	14891	32443	2.36%	0.230%
23	12.392	1729	1734	1735	rVV2	15998	31499	2.29%	0.223%
24	12.416	1735	1738	1740	rVV	29194	42078	3.06%	0.298%
25	12.441	1740	1742	1745	rVV	25626	35066	2.55%	0.248%
26	12.489	1745	1750	1760	rVB3	754575	1373721	100.00%	9.734%
27	12.648	1772	1776	1778	rBV2	24315	38112	2.77%	0.270%
28	12.672	1778	1780	1784	rVB	32219	39785	2.90%	0.282%
29	12.739	1784	1791	1794	rBV	99095	169778	12.36%	1.203%
30	12.770	1794	1796	1798	rVV	65973	84065	6.12%	0.596%
31	12.813	1798	1803	1808	rVV2	293456	503623	36.66%	3.569%
32	12.861	1808	1811	1817	rVB3	27745	42320	3.08%	0.300%
33	12.977	1822	1830	1837	rBV	107300	214871	15.64%	1.523%
34	13.044	1837	1841	1848	rVB2	66896	102532	7.46%	0.727%
35	13.123	1848	1854	1859	rBV	185726	273188	19.89%	1.936%
36	13.172	1859	1862	1866	rVB2	12486	16549	1.20%	0.117%

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Stop Thrs : 0

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Peak separation: 5

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37	13.251	1867	1875	1879	rVB5	31246	66262	4.82%	0.470%
38	13.319	1879	1886	1890	rBV2	97866	175182	12.75%	1.241%
39	13.367	1890	1894	1897	rVV3	53343	79901	5.82%	0.566%
40	13.428	1897	1904	1907	rVV	352438	591671	43.07%	4.193%
41	13.459	1907	1909	1914	rVV	193895	248099	18.06%	1.758%
42	13.508	1914	1917	1923	rVB2	25772	32261	2.35%	0.229%
43	13.599	1928	1932	1933	rBV	26994	33801	2.46%	0.240%
44	13.629	1933	1937	1940	rVV2	155130	224003	16.31%	1.587%
45	13.666	1940	1943	1952	rVB3	162735	313650	22.83%	2.222%
46	13.757	1952	1958	1962	rBV2	230793	351858	25.61%	2.493%
47	13.825	1962	1969	1975	rBV	87837	146838	10.69%	1.040%
48	13.892	1975	1980	1982	rBV	73532	112320	8.18%	0.796%
49	13.916	1982	1984	1989	rVB	101491	134047	9.76%	0.950%
50	13.977	1989	1994	1998	rVB2	173040	266825	19.42%	1.891%
51	14.026	1998	2002	2006	rVB2	22837	38298	2.79%	0.271%
52	14.081	2006	2011	2016	rVB	165231	261920	19.07%	1.856%
53	14.142	2016	2021	2022	rBV2	29899	46829	3.41%	0.332%
54	14.221	2028	2034	2037	rBV3	67102	126932	9.24%	0.899%
55	14.263	2037	2041	2044	rVV5	65953	137789	10.03%	0.976%
56	14.300	2044	2047	2050	rVV	87936	149075	10.85%	1.056%
57	14.337	2050	2053	2057	rVV	151441	213375	15.53%	1.512%
58	14.385	2057	2061	2064	rVV	76616	101047	7.36%	0.716%
59	14.422	2064	2067	2071	rVV2	64756	91103	6.63%	0.646%
60	14.526	2079	2084	2087	rVV2	57091	90851	6.61%	0.644%
61	14.568	2087	2091	2095	rVV2	75913	138735	10.10%	0.983%
62	14.617	2095	2099	2103	rVB	166552	235835	17.17%	1.671%
63	14.678	2103	2109	2116	rBV3	263532	601813	43.81%	4.264%
64	14.739	2116	2119	2125	rVV2	65446	107376	7.82%	0.761%
65	14.794	2125	2128	2130	rVV4	11439	16307	1.19%	0.116%
66	14.837	2130	2135	2141	rVV	183567	269708	19.63%	1.911%
67	14.910	2144	2147	2149	rVV2	23130	31558	2.30%	0.224%
68	14.946	2149	2153	2156	rVV2	80510	140767	10.25%	0.997%
69	14.983	2156	2159	2165	rVV2	54450	99002	7.21%	0.702%
70	15.056	2165	2171	2177	rVB2	89052	190528	13.87%	1.350%
71	15.141	2182	2185	2191	rVB3	26959	47994	3.49%	0.340%
72	15.215	2191	2197	2201	rBV5	36293	87648	6.38%	0.621%
73	15.294	2206	2210	2215	rBV	71826	117232	8.53%	0.831%
74	15.355	2215	2220	2222	rBV2	43200	81089	5.90%	0.575%
75	15.440	2230	2234	2242	rVB4	63391	114492	8.33%	0.811%
76	15.532	2242	2249	2256	rBV3	45295	109136	7.94%	0.773%

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77	15.611	2256	2262	2266	rVV5	17009	32808	2.39%	0.232%
78	15.696	2266	2276	2277	rBV5	34432	76381	5.56%	0.541%
79	15.733	2277	2282	2289	rVB	125052	224298	16.33%	1.589%
80	15.818	2292	2296	2302	rVB4	15704	28321	2.06%	0.201%
81	15.891	2302	2308	2313	rBV4	26478	60234	4.38%	0.427%
82	16.038	2325	2332	2338	rBV	71630	154738	11.26%	1.096%
83	16.233	2356	2364	2370	rBV3	49306	104513	7.61%	0.741%
84	16.300	2370	2375	2380	rBV3	28144	53210	3.87%	0.377%
85	16.495	2400	2407	2413	rBV5	23670	61644	4.49%	0.437%
86	16.629	2425	2429	2438	rVB4	11146	29051	2.11%	0.206%

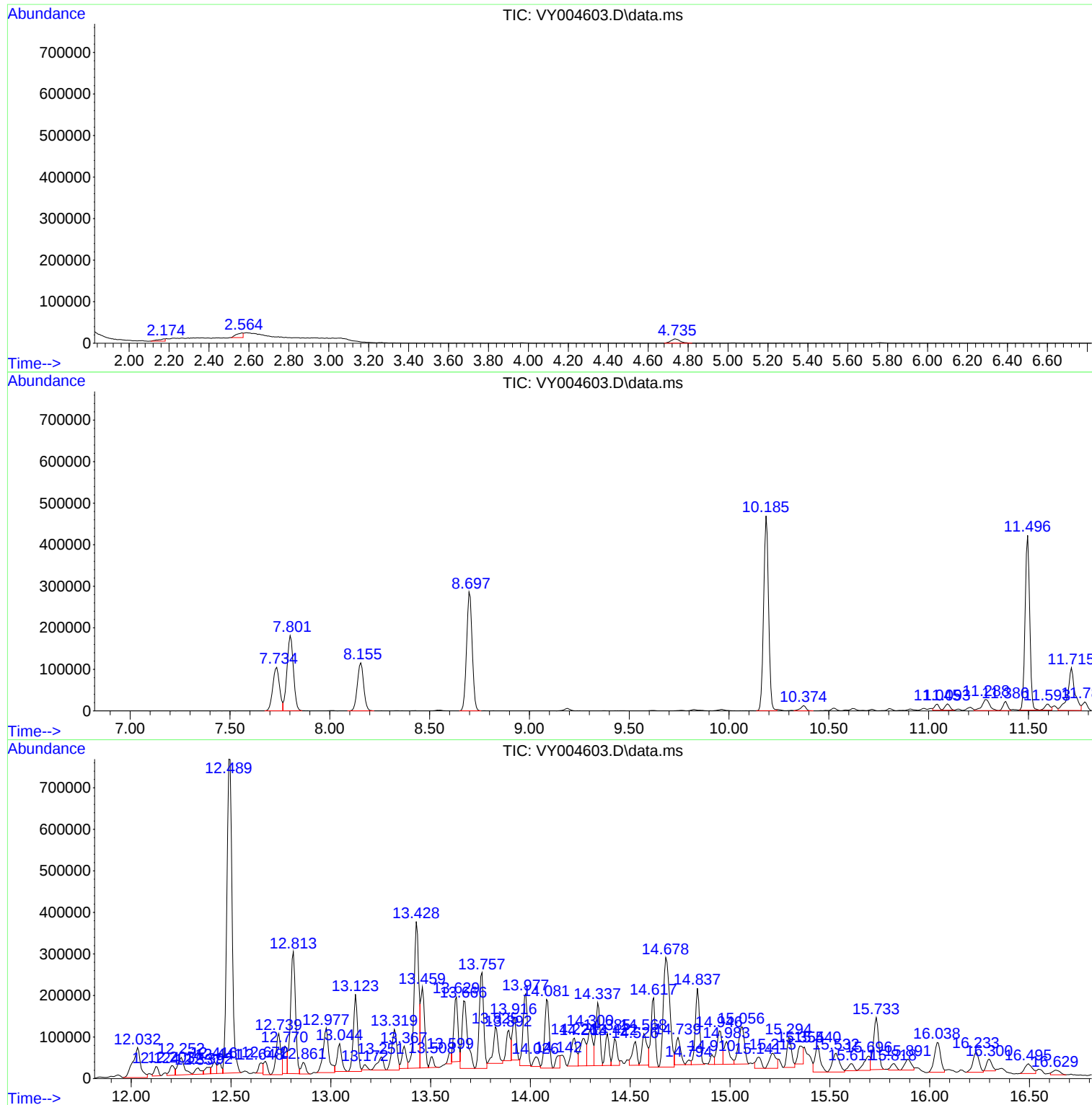
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

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



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ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EO-04-042221

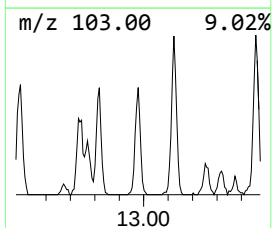
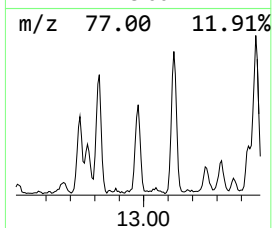
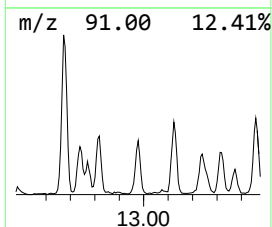
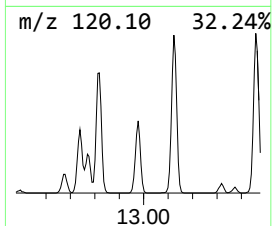
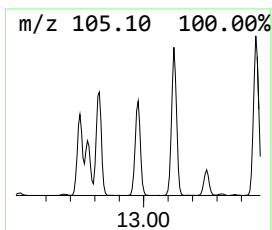
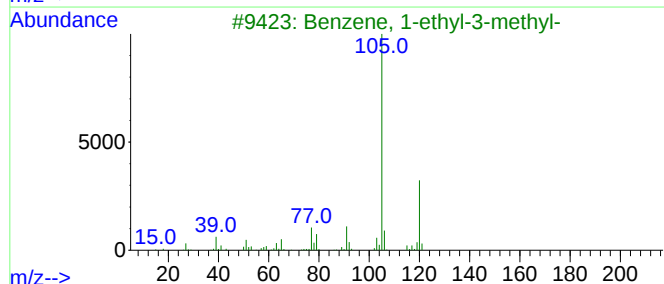
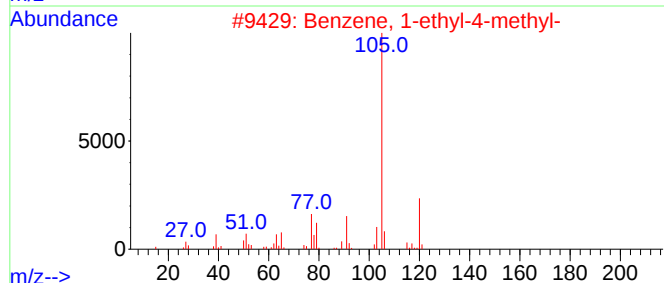
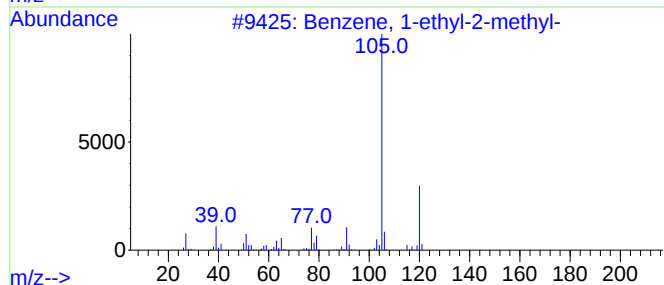
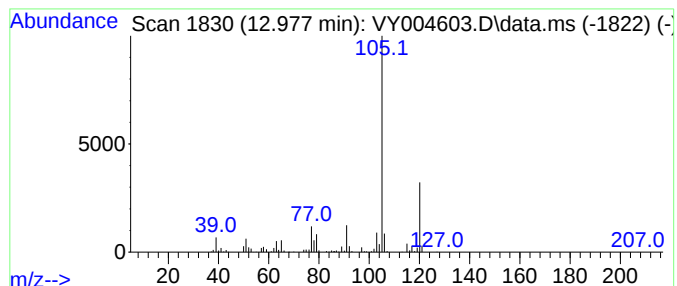
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

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.
12.977	18.16 ug/l	214871	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	95
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5			Mesitylene	120	C9H12	000108-67-8	91



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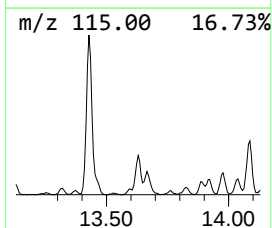
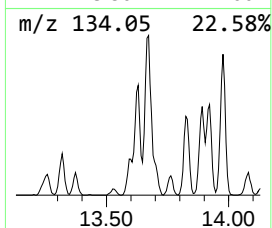
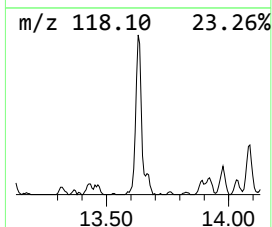
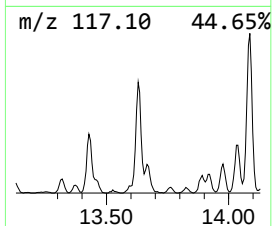
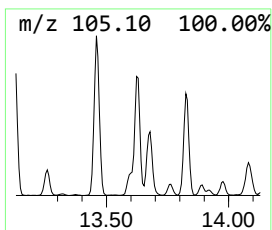
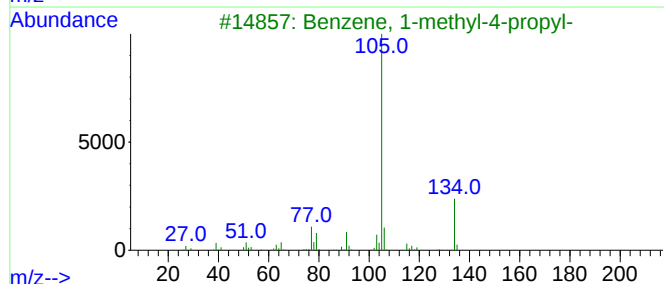
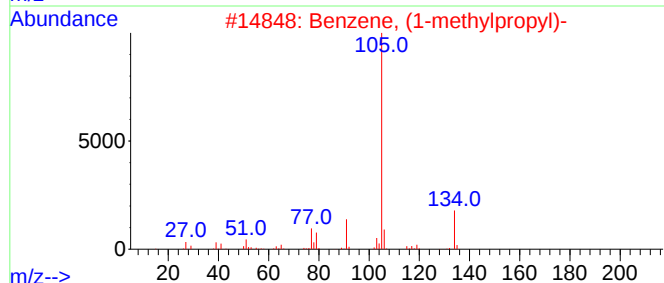
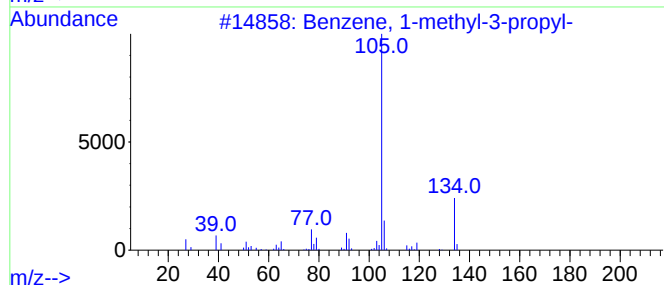
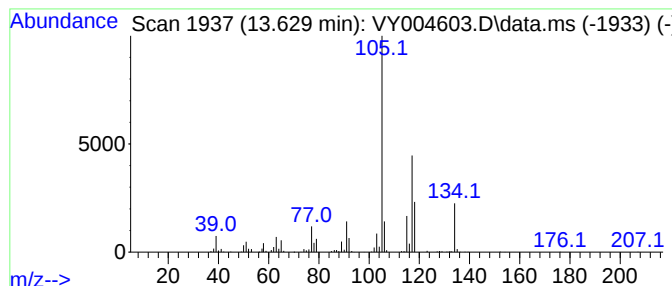
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1-methyl-3-propyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.629	18.93 ug/l	224003	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	50
2			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	43
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	38
4			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	38
5			2-Tolylloxirane	134	C9H10O	002783-26-8	35



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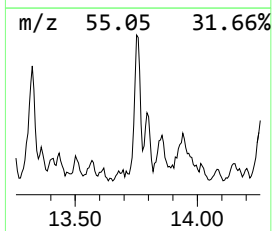
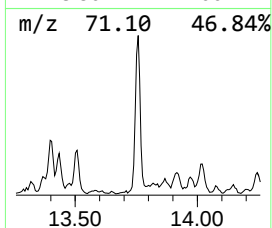
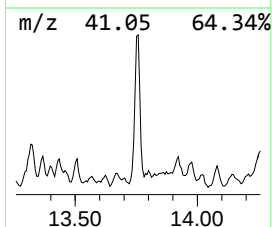
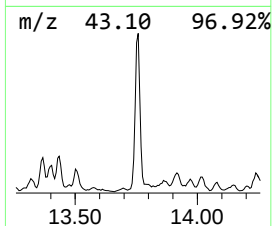
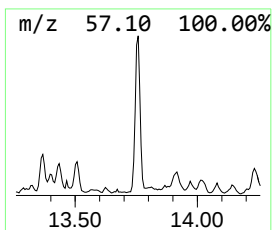
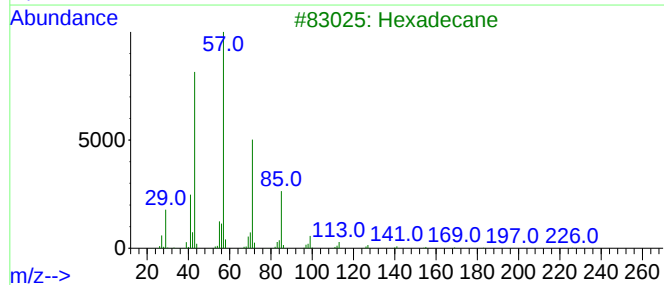
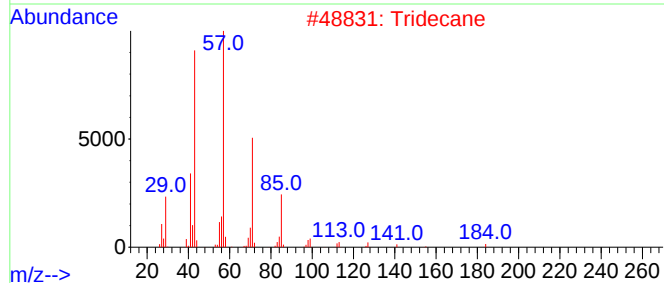
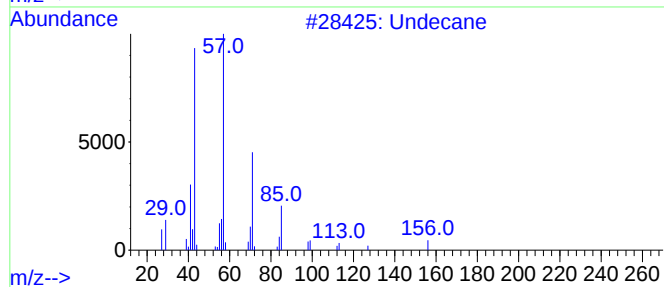
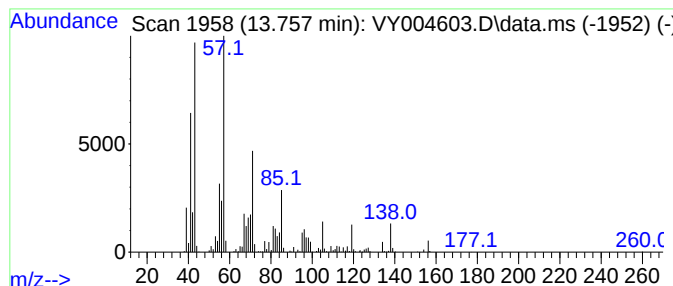
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

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

Peak Number 3 Undecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.758	29.73 ug/l	351858	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	60
2			Tridecane	184	C13H28	000629-50-5	50
3			Hexadecane	226	C16H34	000544-76-3	47
4			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	47
5			Octane, 1,1'-oxybis-	242	C16H34O	000629-82-3	47



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EO-04-042221

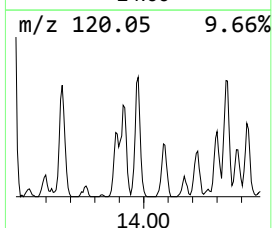
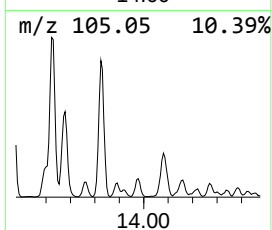
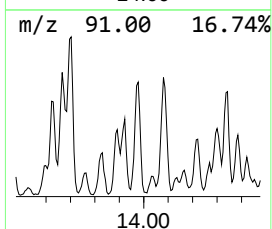
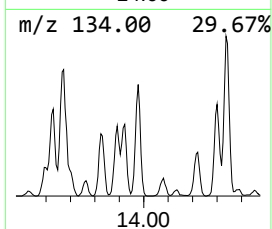
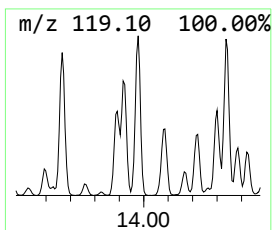
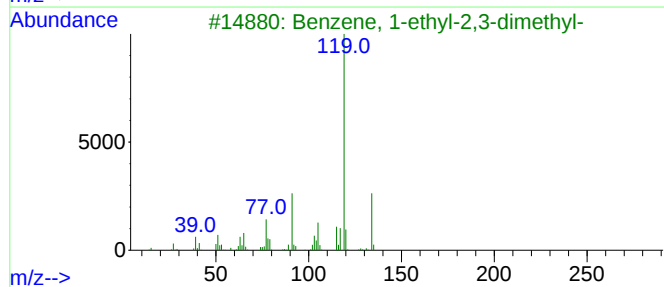
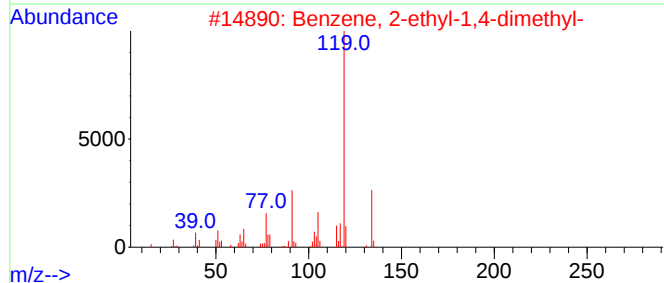
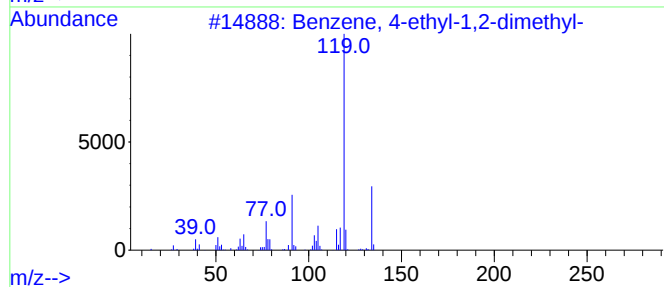
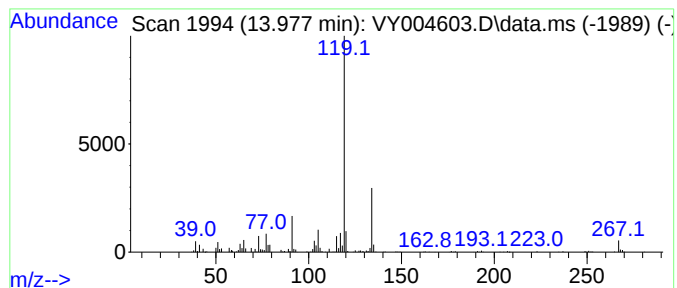
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.977	22.55 ug/l	266825	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5			o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY042321\
Data File : VY004603.D
Acq On : 23 Apr 2021 14:25
Operator : SY/MD
Sample : M2142-01
Misc : 7.05G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-04-042221

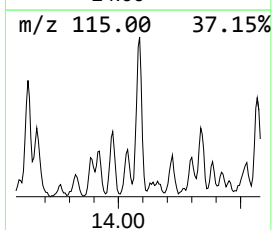
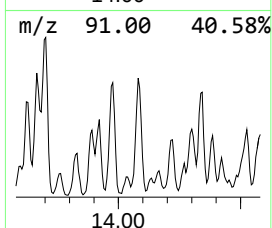
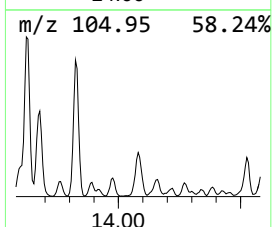
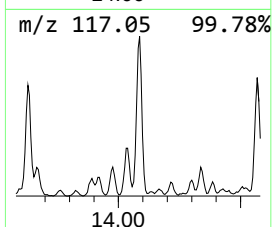
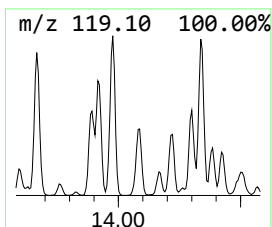
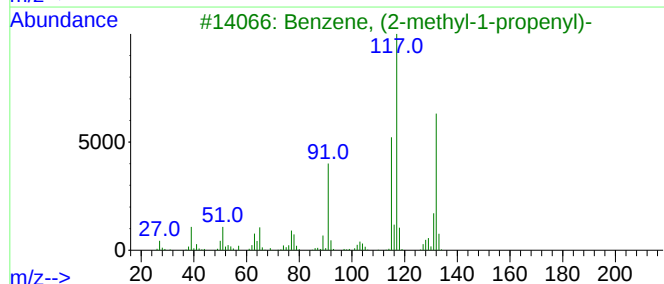
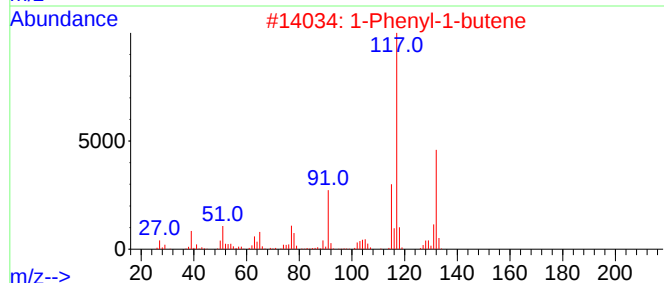
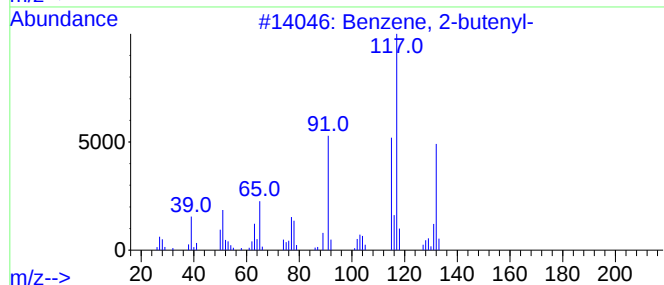
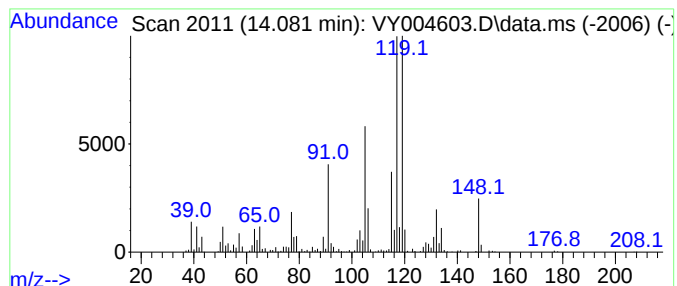
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 2-butenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.081	22.13 ug/l	261920	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-butenyl-	132	C10H12	001560-06-1	74
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	64
3			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	50
4			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	50
5			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY042321\
Data File : VY004603.D
Acq On : 23 Apr 2021 14:25
Operator : SY/MD
Sample : M2142-01
Misc : 7.05G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-04-042221

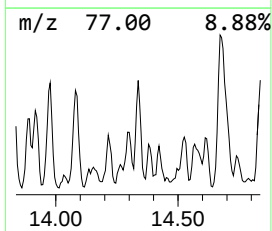
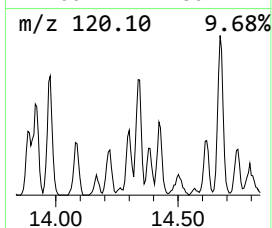
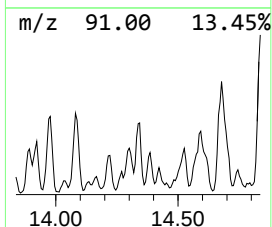
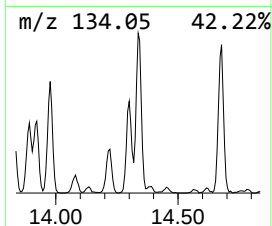
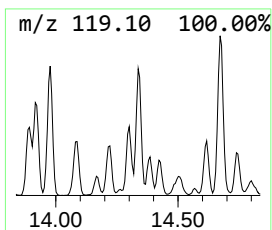
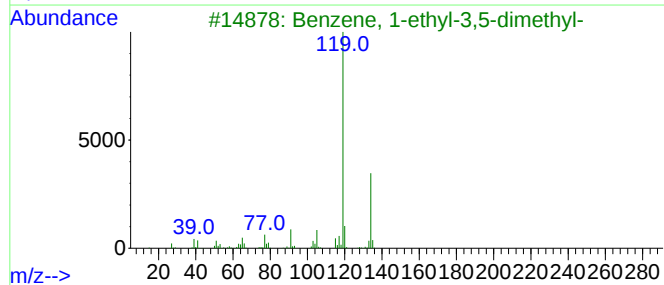
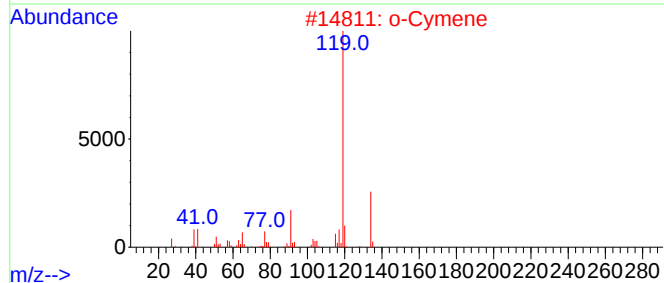
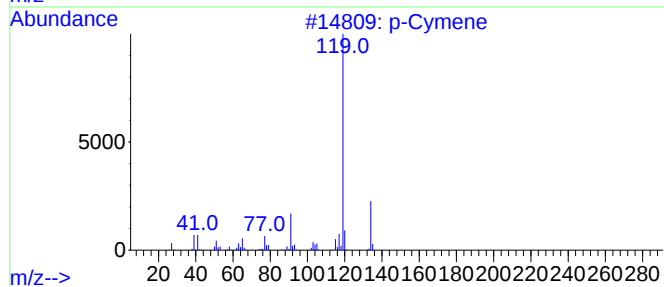
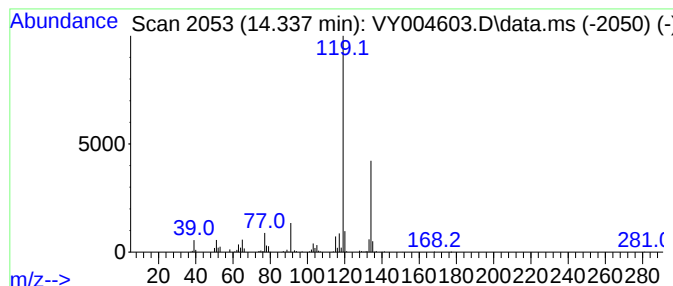
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Quant Title : SW846 8260

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

Peak Number 6 p-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.337	18.03 ug/l	213375	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	95
2			o-Cymene	134	C10H14	000527-84-4	95
3			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY042321\
Data File : VY004603.D
Acq On : 23 Apr 2021 14:25
Operator : SY/MD
Sample : M2142-01
Misc : 7.05G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-04-042221

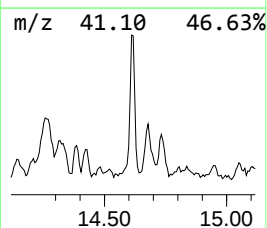
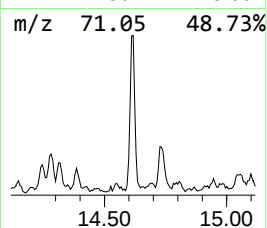
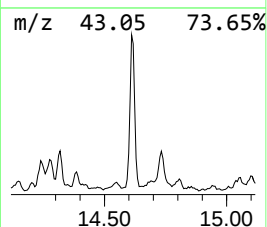
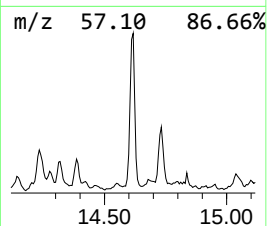
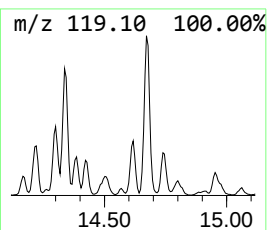
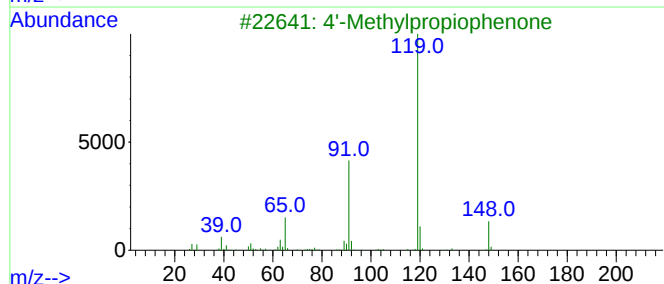
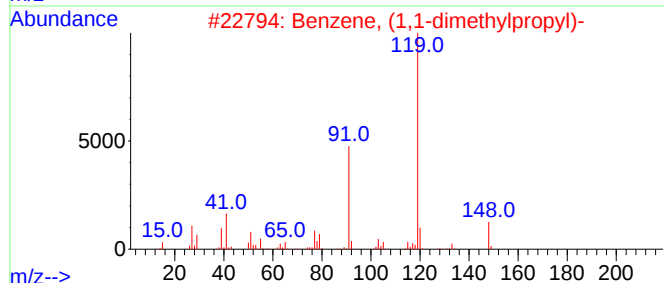
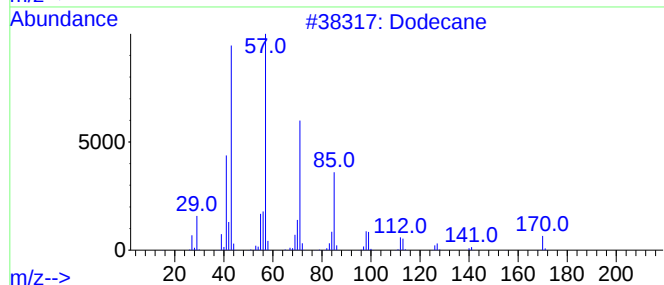
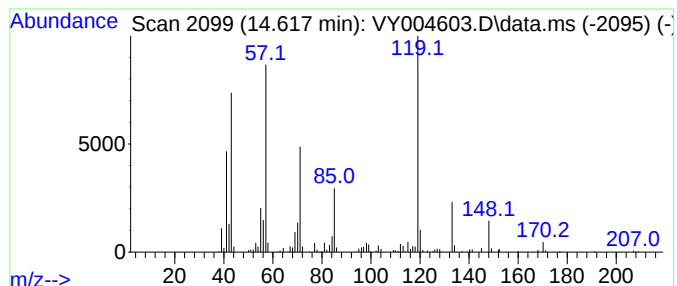
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Quant Title : SW846 8260

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

Peak Number 7 unknown14.617 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.617	19.93 ug/l	235835	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	38
2			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	38
3			4'-Methylpropiophenone	148	C10H12O	005337-93-9	35
4			Tetradecane	198	C14H30	000629-59-4	30
5			Pentadecane	212	C15H32	000629-62-9	27



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY042321\
Data File : VY004603.D
Acq On : 23 Apr 2021 14:25
Operator : SY/MD
Sample : M2142-01
Misc : 7.05G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-04-042221

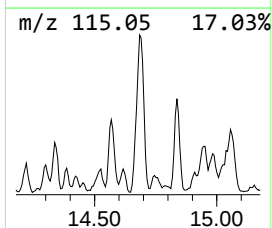
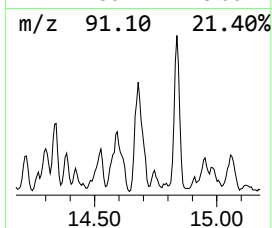
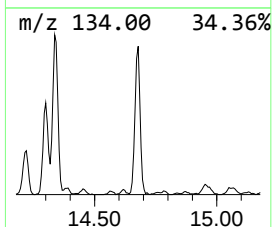
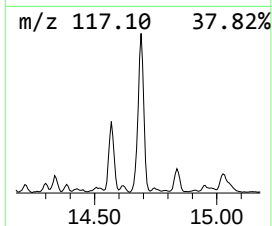
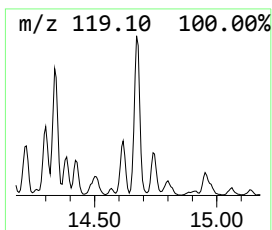
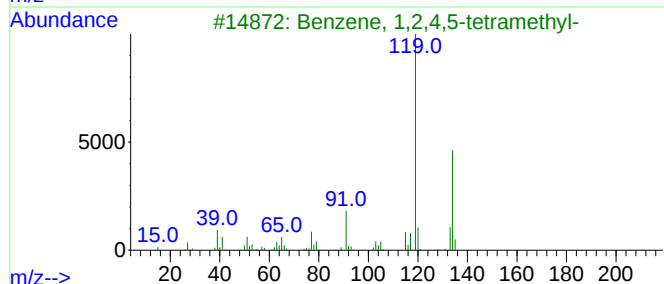
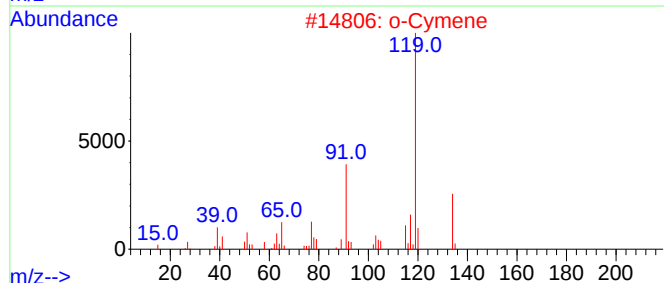
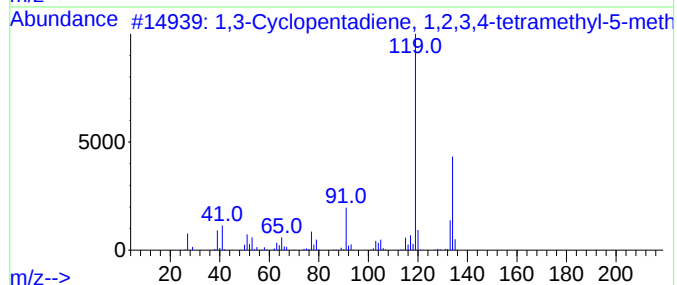
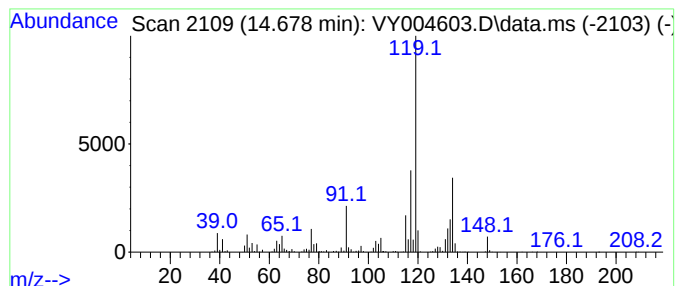
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Quant Title : SW846 8260

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

Peak Number 8 1,3-Cyclopentadiene, 1,2,3,... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.678	50.86 ug/l	601813	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	81
2			o-Cymene	134	C10H14	000527-84-4	76
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	76
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	70
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY042321\
Data File : VY004603.D
Acq On : 23 Apr 2021 14:25
Operator : SY/MD
Sample : M2142-01
Misc : 7.05G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-04-042221

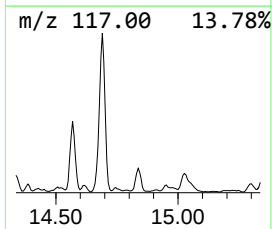
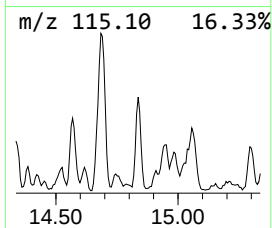
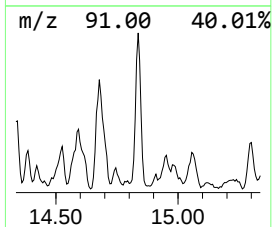
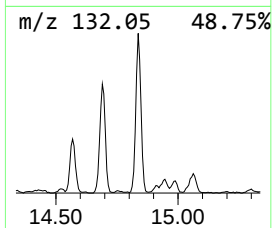
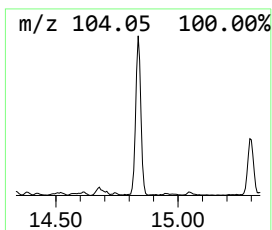
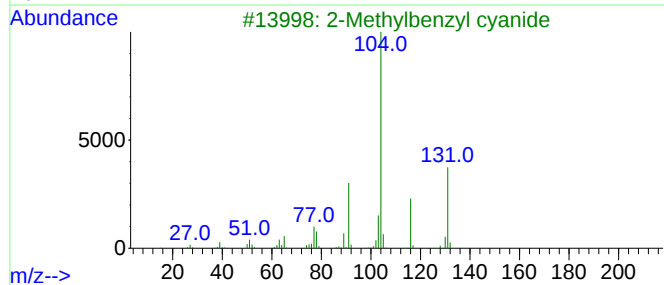
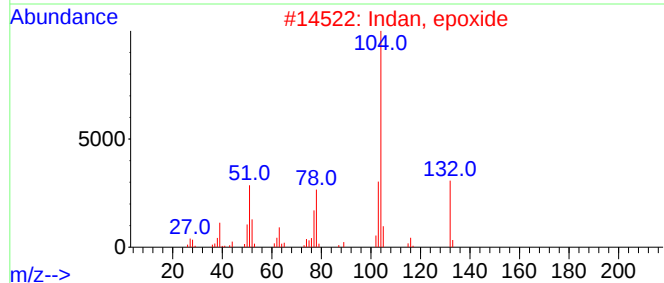
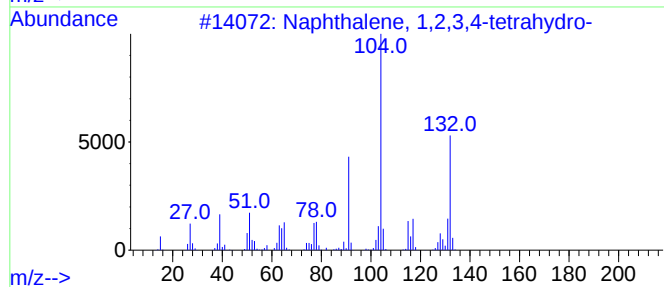
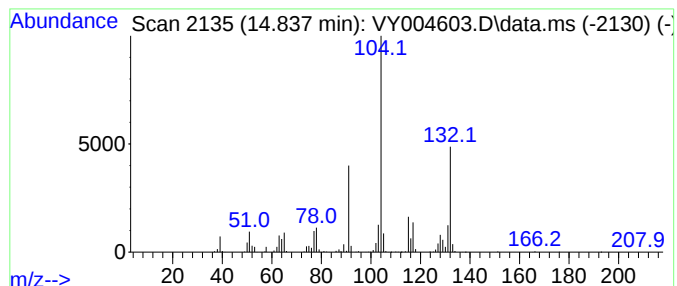
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.836	22.79 ug/l	269708	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2			Indan, epoxide	132	C9H8O	000768-22-9	58
3			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	46
4			Benzene, cyclobutyl-	132	C10H12	004392-30-7	45
5			Isoquinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000091-21-4	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY042321\
Data File : VY004603.D
Acq On : 23 Apr 2021 14:25
Operator : SY/MD
Sample : M2142-01
Misc : 7.05G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-04-042221

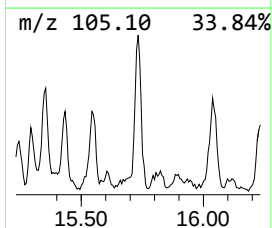
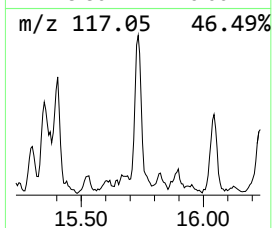
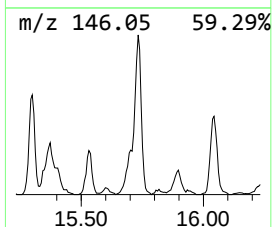
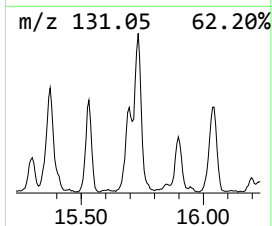
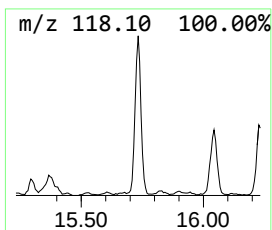
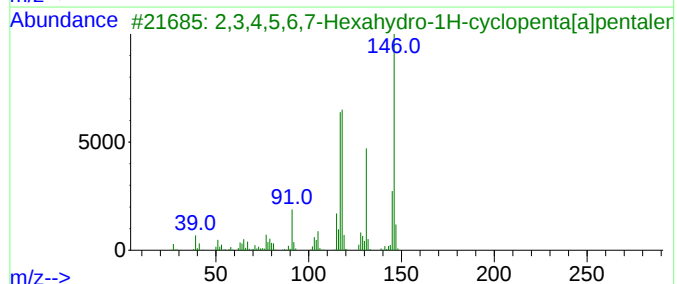
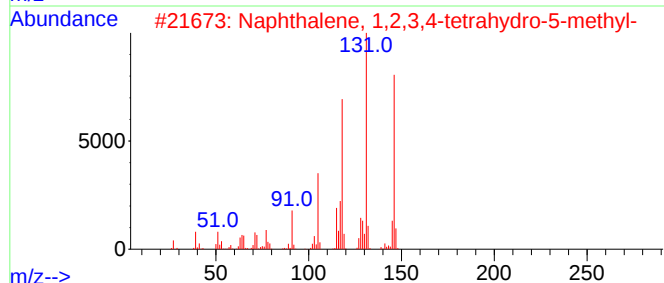
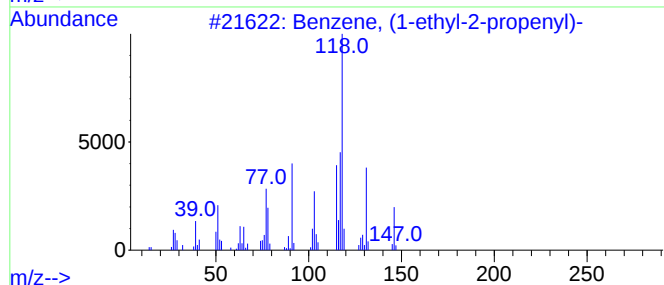
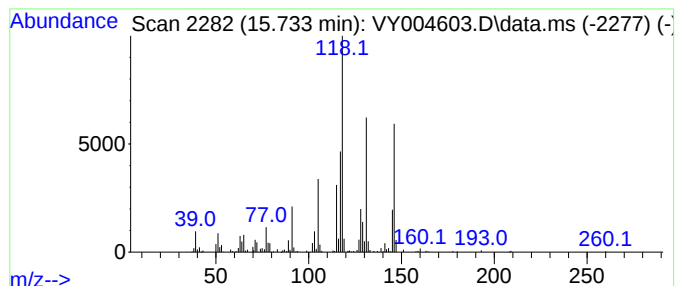
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, (1-ethyl-2-propenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.733	18.95 ug/l	224298	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-ethyl-2-propenyl)-	146	C11H14	019947-22-9	64
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	58
3			2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1000189-31-0	49
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	43
5			Pyridine, 3-(3,4-dihydro-2H-pyrr...	146	C9H10N2	000532-12-7	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY042321\
Data File : VY004603.D
Acq On : 23 Apr 2021 14:25
Operator : SY/MD
Sample : M2142-01
Misc : 7.05G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-04-042221

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...	12.977	18.2	ug/l	214871	4	13.428	591671	50.0
Benzene, 1-meth...	13.629	18.9	ug/l	224003	4	13.428	591671	50.0
Undecane	13.758	29.7	ug/l	351858	4	13.428	591671	50.0
Benzene, 4-ethy...	13.977	22.6	ug/l	266825	4	13.428	591671	50.0
Benzene, 2-bute...	14.081	22.1	ug/l	261920	4	13.428	591671	50.0
p-Cymene	14.337	18.0	ug/l	213375	4	13.428	591671	50.0
unknown14.617	14.617	19.9	ug/l	235835	4	13.428	591671	50.0
1,3-Cyclopentad...	14.678	50.9	ug/l	601813	4	13.428	591671	50.0
Naphthalene, 1,...	14.836	22.8	ug/l	269708	4	13.428	591671	50.0
Benzene, (1-eth...	15.733	18.9	ug/l	224298	4	13.428	591671	50.0