

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY061021\
 Data File : VY005075.D
 Acq On : 10 Jun 2021 13:21
 Operator : SY/MD
 Sample : M2674-01
 Misc : 5.79G/5ML/MSVOA_Y/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 TMR-DS-01-060921

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\82Y060221S.M

Title : SW846 8260

Signal : TIC: VY005075.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.960	341	351	359	rBV	44937	126805	1.15%	0.090%
2	7.728	958	969	974	rBV	126867	301280	2.74%	0.214%
3	7.795	974	980	990	rVB	212444	474597	4.31%	0.337%
4	8.149	1028	1038	1050	rBV2	142671	326172	2.96%	0.232%
5	8.557	1094	1105	1118	rBV2	68664	164847	1.50%	0.117%
6	8.691	1118	1127	1135	rBV	355531	732698	6.66%	0.521%
7	10.185	1362	1372	1377	rBV	581193	1050308	9.55%	0.747%
8	10.246	1377	1382	1387	rVV	1174587	2007318	18.24%	1.427%
9	10.289	1387	1389	1395	rVV2	120736	186349	1.69%	0.133%
10	10.374	1395	1403	1409	rVB2	178523	326124	2.96%	0.232%
11	11.490	1578	1586	1593	rBV	546748	948299	8.62%	0.674%
12	11.593	1596	1603	1609	rBV	2180534	3448655	31.34%	2.452%
13	11.660	1609	1614	1615	rVV	325595	402513	3.66%	0.286%
14	11.703	1615	1621	1633	rVB	5808710	10654342	96.83%	7.576%
15	11.801	1633	1637	1644	rBV4	76314	137511	1.25%	0.098%
16	12.032	1665	1675	1683	rBV	3693710	5848053	53.15%	4.158%
17	12.136	1689	1692	1696	rVB2	87803	125650	1.14%	0.089%
18	12.239	1704	1709	1713	rBV2	215828	357094	3.25%	0.254%
19	12.331	1719	1724	1729	rBV3	385066	645029	5.86%	0.459%
20	12.386	1729	1733	1740	rVB2	137049	242081	2.20%	0.172%
21	12.489	1740	1750	1759	rBV3	995158	1934609	17.58%	1.376%
22	12.672	1775	1780	1785	rBV	1079083	1536452	13.96%	1.093%
23	12.739	1785	1791	1793	rBV	3023570	4649500	42.26%	3.306%
24	12.770	1793	1796	1799	rVV	1727910	2363907	21.48%	1.681%
25	12.813	1799	1803	1809	rVB	1833603	2929024	26.62%	2.083%
26	12.873	1809	1813	1821	rVB8	59678	119619	1.09%	0.085%
27	12.971	1821	1829	1838	rVB	2133242	3578352	32.52%	2.544%
28	13.123	1838	1854	1859	rBV	5249503	7980727	72.53%	5.675%
29	13.172	1859	1862	1867	rVB3	393015	587407	5.34%	0.418%
30	13.239	1867	1873	1877	rBV6	131508	349349	3.18%	0.248%
31	13.294	1878	1882	1883	rBV2	196433	249919	2.27%	0.178%
32	13.343	1883	1890	1900	rVB2	1542206	3506324	31.87%	2.493%
33	13.459	1900	1909	1913	rBV2	2549620	4268624	38.80%	3.035%
34	13.501	1913	1916	1926	rVB3	463884	861011	7.83%	0.612%
35	13.593	1926	1931	1932	rBV2	502383	639588	5.81%	0.455%
36	13.629	1932	1937	1940	rBV	2252839	3112841	28.29%	2.213%

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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_Y\methods\82Y060221S.M
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37	13.666	1940	1943	1947	rBV2	613951	787139	7.15%	0.560%
38	13.757	1955	1958	1962	rVB2	302081	378963	3.44%	0.269%
39	13.818	1962	1968	1974	rVB2	999473	2068257	18.80%	1.471%
40	13.886	1974	1979	1981	rBV	676290	976980	8.88%	0.695%

41	13.916	1981	1984	1989	rVV	1041488	1469474	13.36%	1.045%
42	13.971	1989	1993	1998	rVB	1220960	1779401	16.17%	1.265%
43	14.032	1998	2003	2006	rBV	318955	433041	3.94%	0.308%
44	14.081	2006	2011	2022	rVB2	1294413	2409585	21.90%	1.713%
45	14.209	2022	2032	2037	rVB4	623219	1526761	13.88%	1.086%

46	14.270	2037	2042	2046	rBV	1370318	2263548	20.57%	1.610%
47	14.392	2050	2062	2066	rVV2	4132833	11002701	100.00%	7.824%
48	14.428	2066	2068	2074	rVV3	756033	1213902	11.03%	0.863%
49	14.477	2074	2076	2080	rVV4	146916	249970	2.27%	0.178%
50	14.520	2081	2083	2086	rVV	167965	202271	1.84%	0.144%

51	14.568	2086	2091	2103	rVB3	1495816	3810235	34.63%	2.709%
52	14.684	2103	2110	2116	rBV3	1785445	4022465	36.56%	2.860%
53	14.751	2116	2121	2126	rVV2	905356	1428622	12.98%	1.016%
54	14.837	2130	2135	2143	rBV	931602	1449548	13.17%	1.031%
55	14.910	2143	2147	2148	rBV2	252475	320043	2.91%	0.228%

56	14.946	2148	2153	2163	rVV5	744376	2093001	19.02%	1.488%
57	15.056	2164	2171	2178	rVB4	375044	849976	7.73%	0.604%
58	15.196	2184	2194	2197	rBV	2148209	5259118	47.80%	3.740%
59	15.233	2197	2200	2209	rVV2	1852000	4276314	38.87%	3.041%
60	15.318	2209	2214	2224	rVV	3778155	7117208	64.69%	5.061%

61	15.404	2224	2228	2232	rVV	919069	1448725	13.17%	1.030%
62	15.452	2232	2236	2242	rVB3	327586	706097	6.42%	0.502%
63	15.532	2242	2249	2256	rBV4	333723	806388	7.33%	0.573%
64	15.611	2256	2262	2267	rBV3	221488	487890	4.43%	0.347%
65	15.666	2267	2271	2274	rBV4	128234	234126	2.13%	0.166%

66	15.721	2277	2280	2283	rVV3	122164	185638	1.69%	0.132%
67	15.763	2284	2287	2291	rVB	252731	396197	3.60%	0.282%
68	15.855	2291	2302	2315	rVB3	520506	2365375	21.50%	1.682%
69	15.964	2315	2320	2322	rBV	150079	259540	2.36%	0.185%
70	16.001	2322	2326	2329	rBV	125709	195040	1.77%	0.139%

71	16.135	2343	2348	2354	rVB	529359	1050012	9.54%	0.747%
72	16.208	2354	2360	2366	rVV	693904	1685797	15.32%	1.199%
73	16.294	2366	2374	2381	rVB2	1382586	3154131	28.67%	2.243%
74	16.428	2390	2396	2401	rBV2	417514	926740	8.42%	0.659%
75	16.489	2401	2406	2416	rVB	745353	1724616	15.67%	1.226%

76	16.696	2434	2440	2446	rVB	202368	445321	4.05%	0.317%
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ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TMR-DS-01-060921

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\82Y060221S.M
Title : SW846 8260

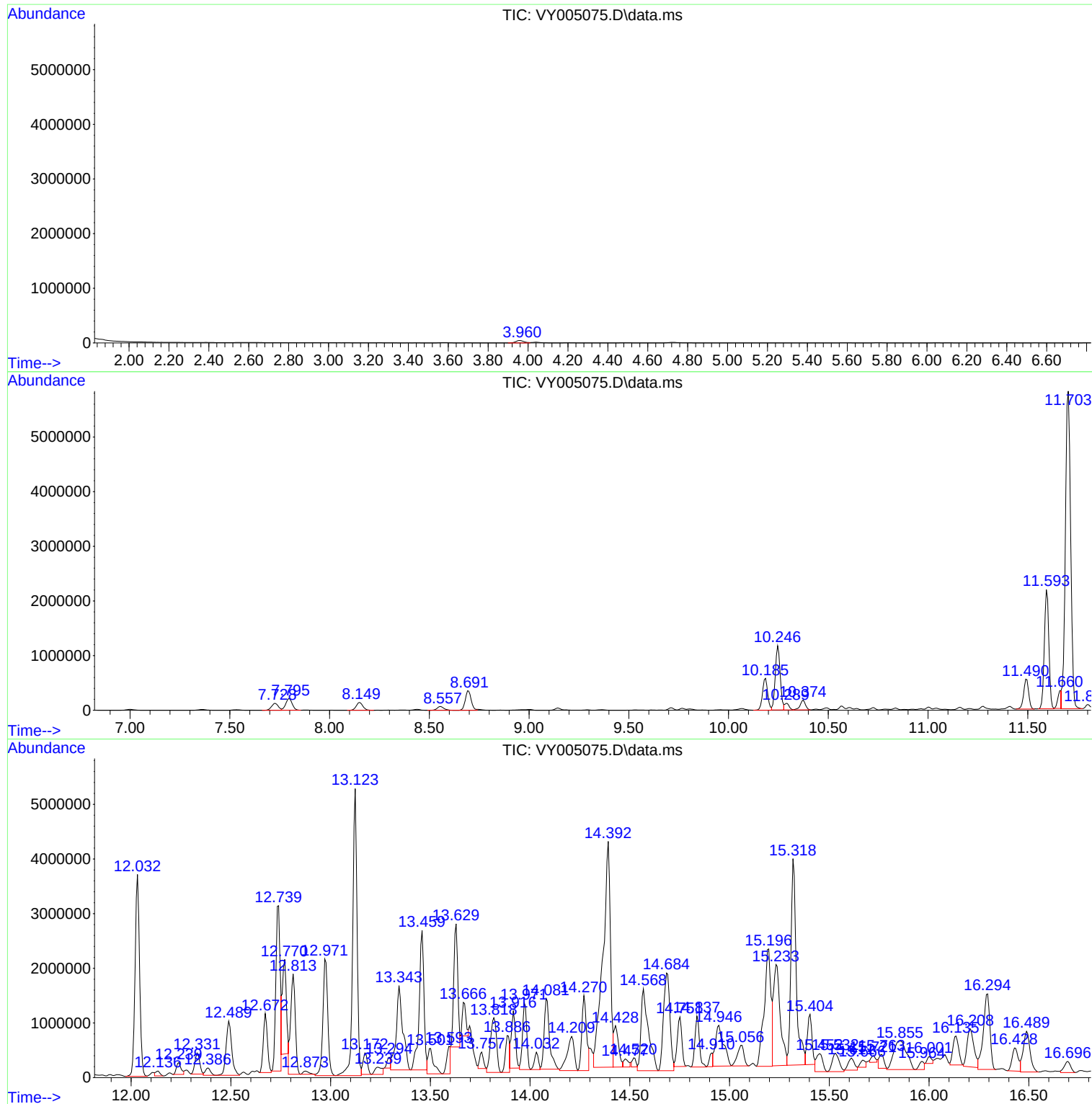
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Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY061021\
Data File : VY005075.D
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Instrument :
MSVOA_Y
ClientSampleId :
TMR-DS-01-060921

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y060221S.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY061021\
Data File : VY005075.D
Acq On : 10 Jun 2021 13:21
Operator : SY/MD
Sample : M2674-01
Misc : 5.79G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TMR-DS-01-060921

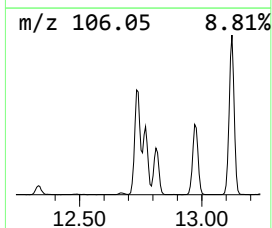
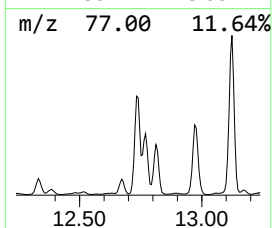
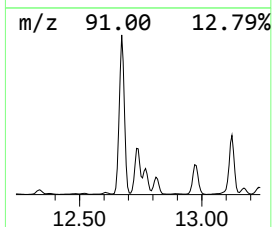
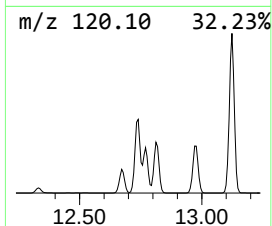
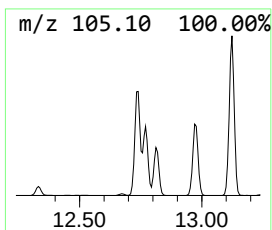
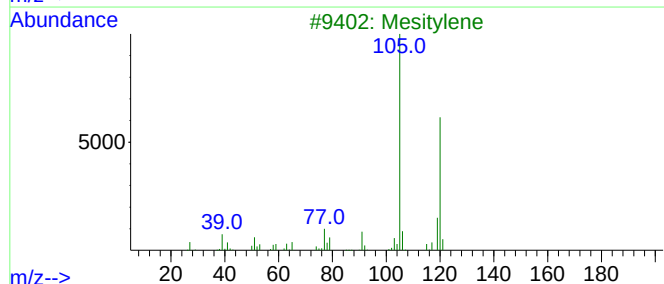
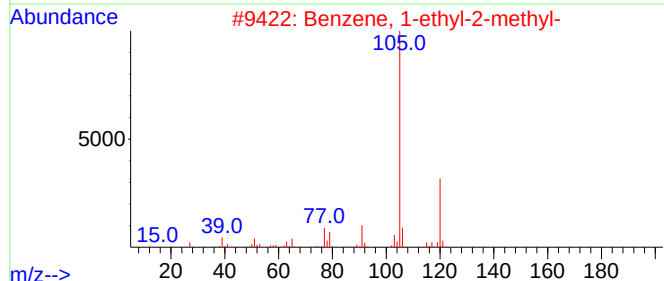
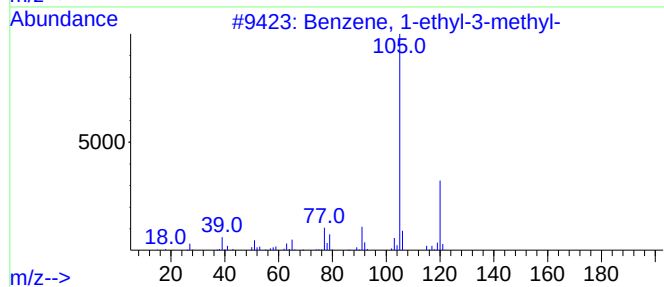
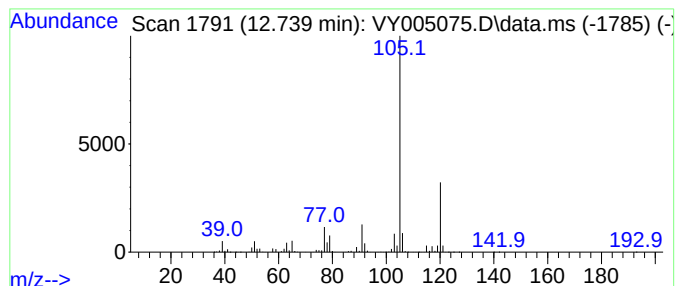
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y060221S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.739	54.46 ug/l	4649500	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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ALS Vial : 9 Sample Multiplier: 1

Instrument :
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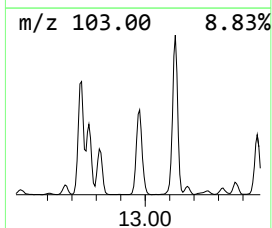
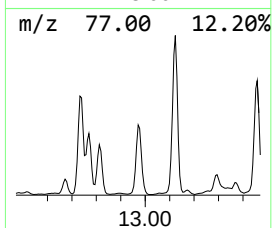
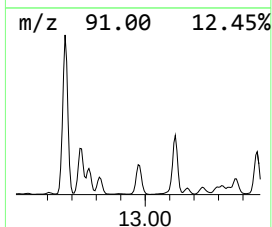
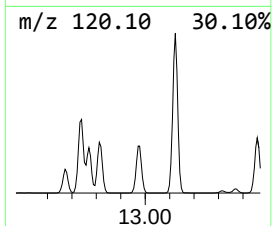
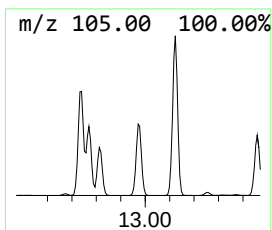
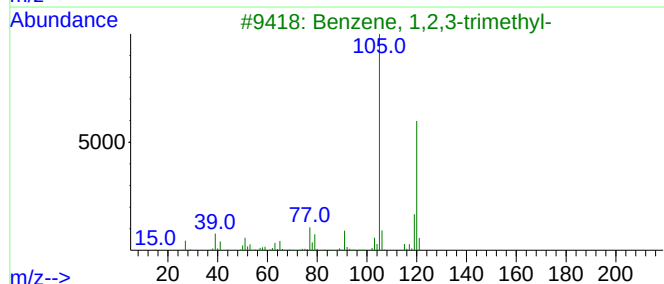
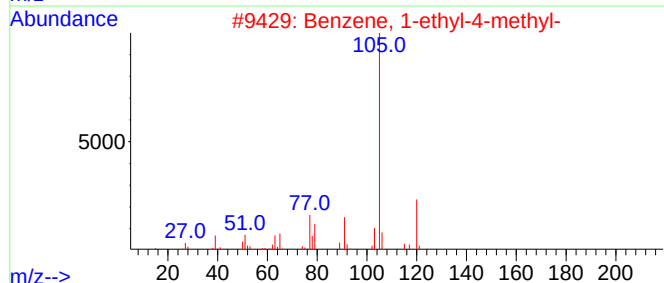
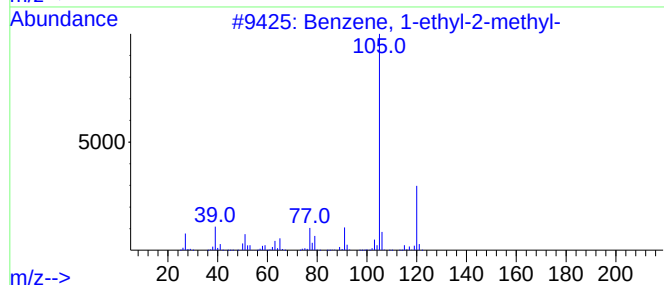
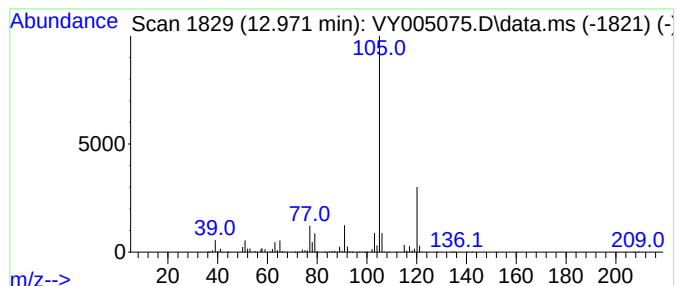
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y060221S.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.971	41.91 ug/l	3578350	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	93
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
4			Mesitylene	120	C9H12	000108-67-8	90
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90



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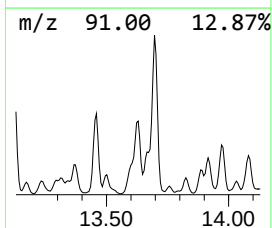
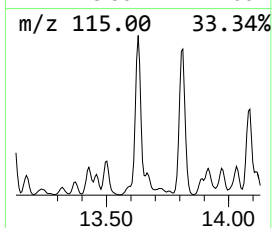
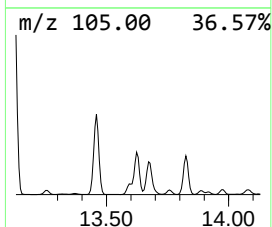
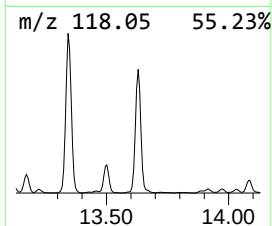
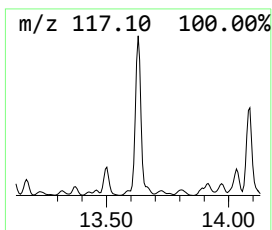
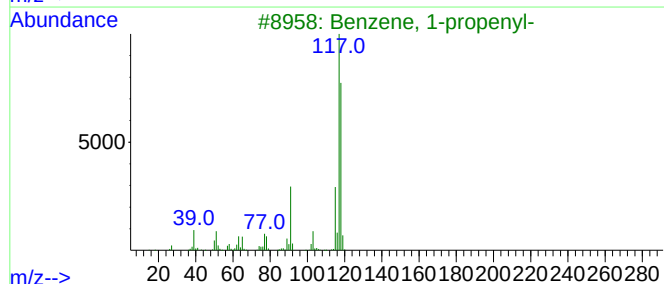
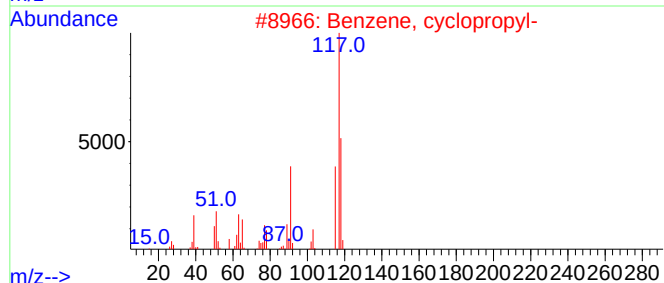
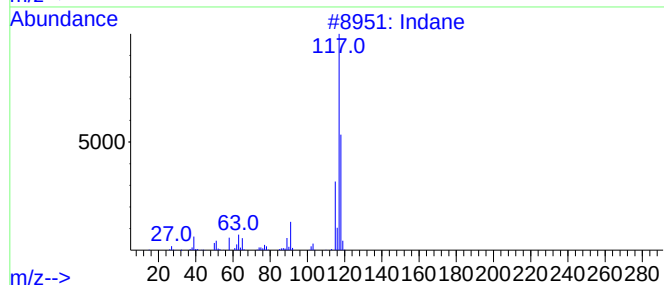
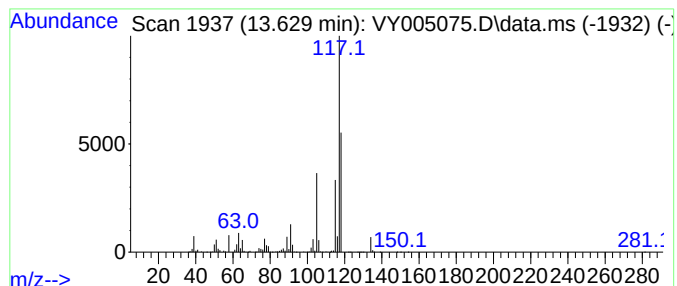
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y060221S.M
Quant Title : SW846 8260

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

Peak Number 3 Indane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.630	36.46 ug/l	3112840	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	70
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
3			Benzene, 1-propenyl-	118	C9H10	000637-50-3	62
4			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	62
5			Deltacyclene	118	C9H10	007785-10-6	58



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ALS Vial : 9 Sample Multiplier: 1

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ClientSampleId :
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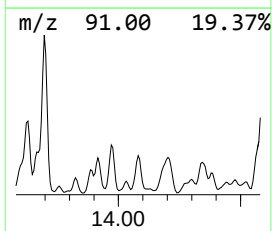
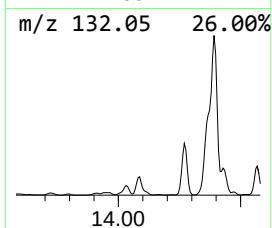
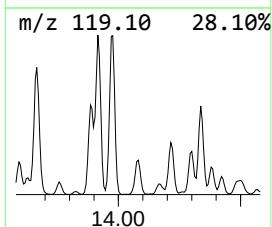
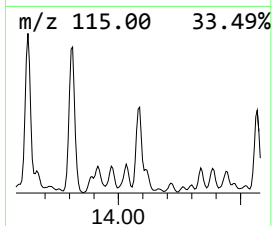
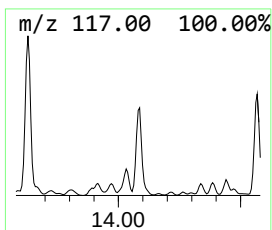
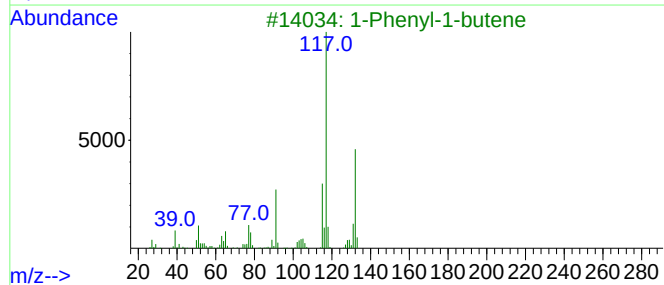
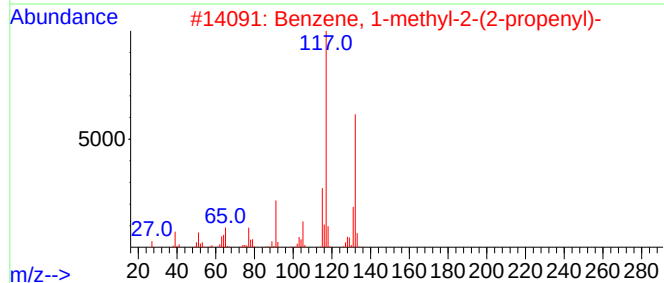
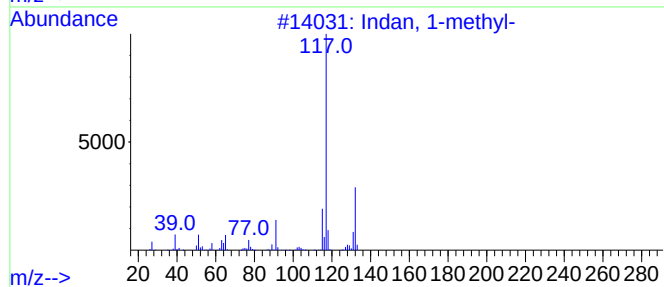
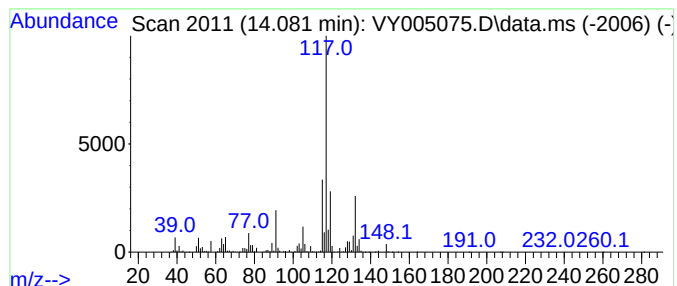
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Quant Title : SW846 8260

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

Peak Number 4 Indan, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.081	28.22 ug/l	2409590	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	81
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	81
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	81
4			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	74
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY061021\
Data File : VY005075.D
Acq On : 10 Jun 2021 13:21
Operator : SY/MD
Sample : M2674-01
Misc : 5.79G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TMR-DS-01-060921

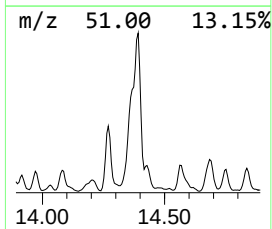
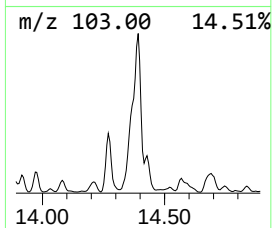
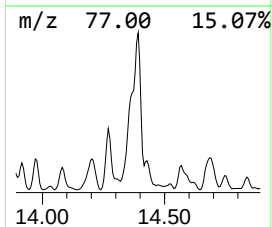
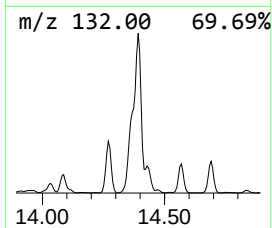
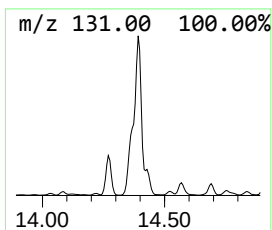
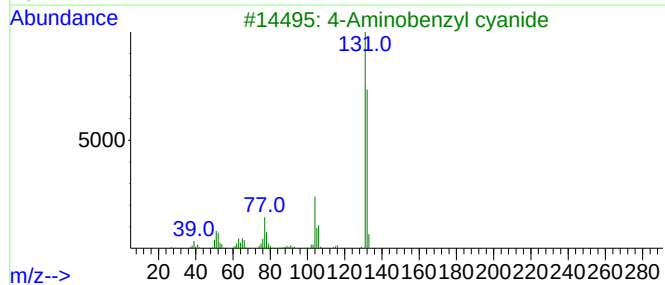
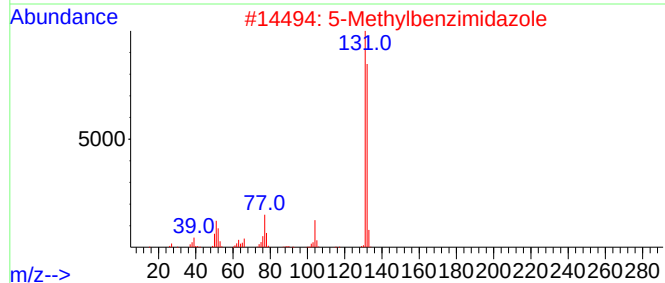
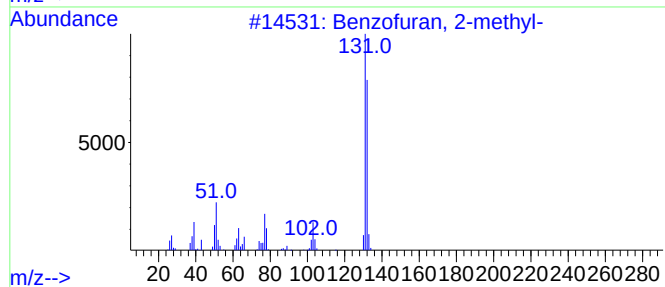
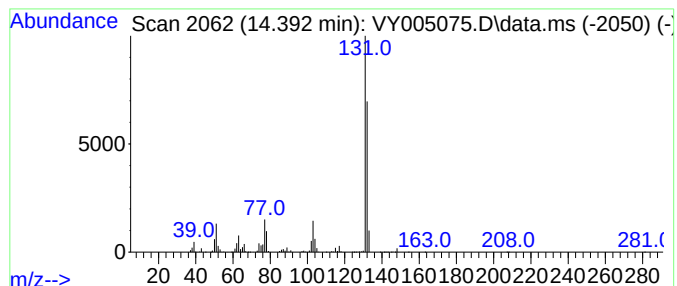
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Quant Title : SW846 8260

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

Peak Number 5 Benzofuran, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.392	128.88 ug/l	11002700	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2			5-Methylbenzimidazole	132	C8H8N2	000614-97-1	87
3			4-Aminobenzyl cyanide	132	C8H8N2	003544-25-0	86
4			3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	83
5			Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY061021\
Data File : VY005075.D
Acq On : 10 Jun 2021 13:21
Operator : SY/MD
Sample : M2674-01
Misc : 5.79G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TMR-DS-01-060921

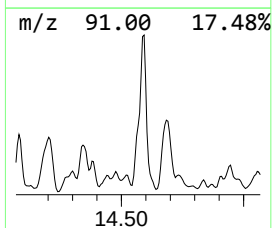
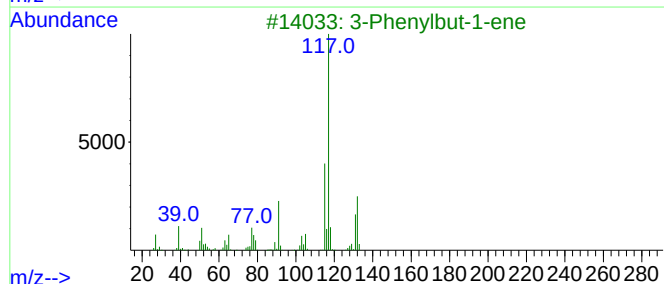
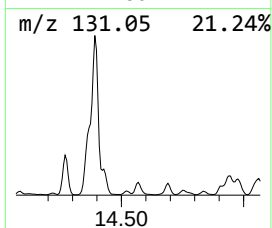
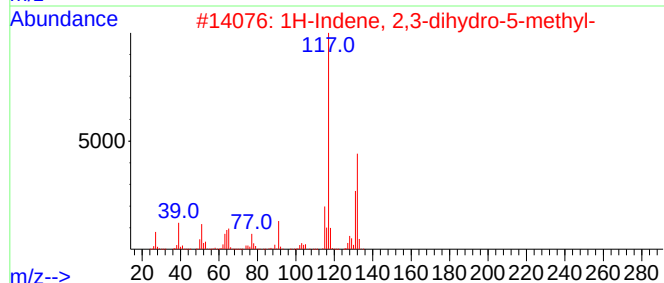
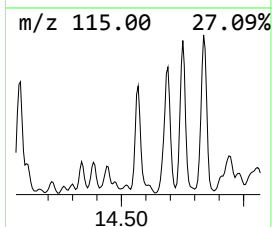
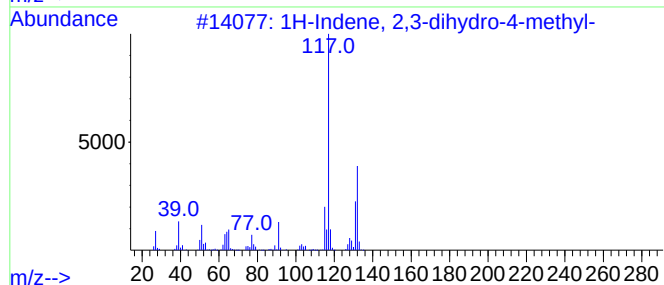
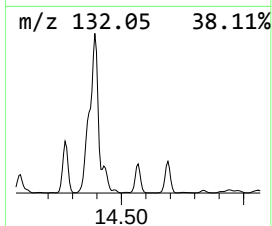
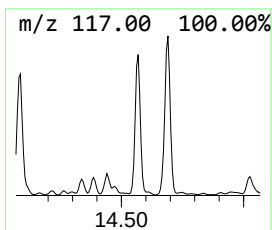
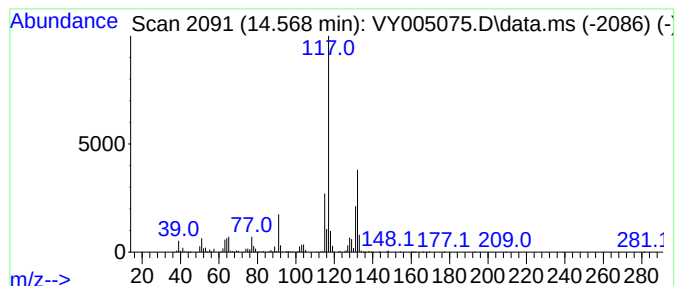
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Quant Title : SW846 8260

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	44.63 ug/l	3810240	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			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4			Indan, 1-methyl-	132	C10H12	000767-58-8	83
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY061021\
Data File : VY005075.D
Acq On : 10 Jun 2021 13:21
Operator : SY/MD
Sample : M2674-01
Misc : 5.79G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TMR-DS-01-060921

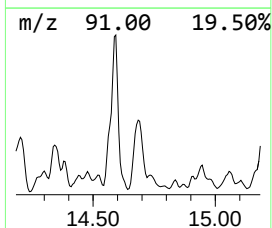
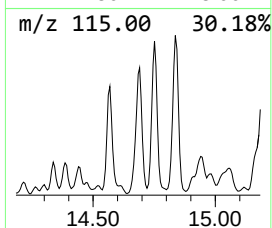
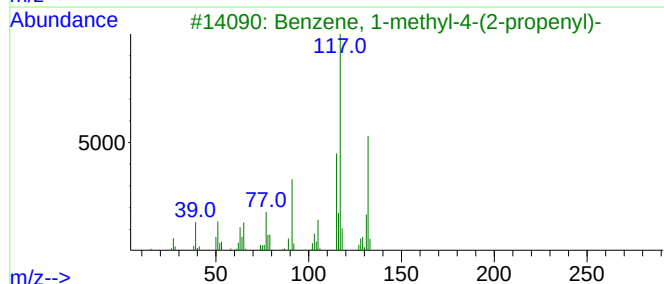
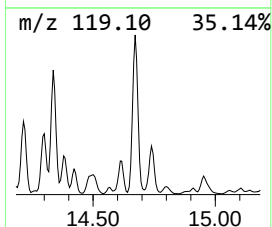
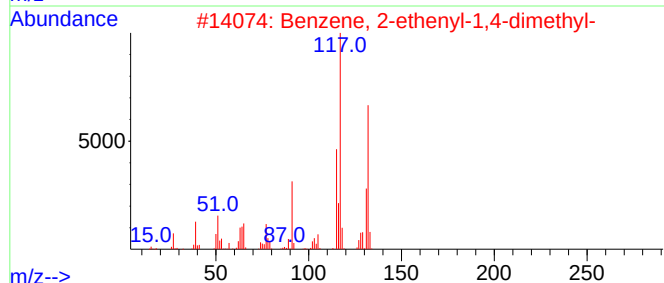
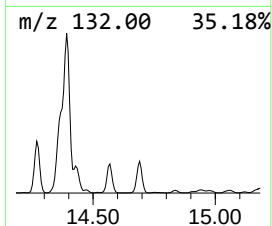
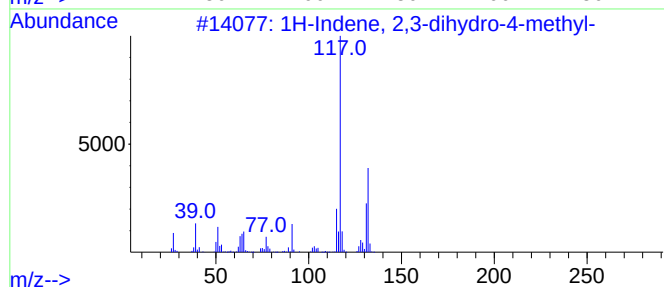
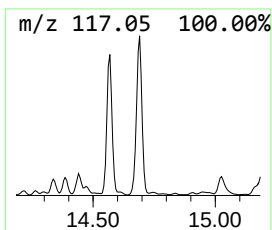
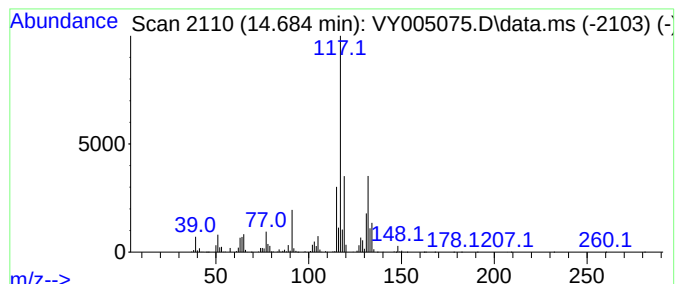
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	47.12 ug/l	4022470	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	93
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
3		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	91
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\VY061021\
Data File : VY005075.D
Acq On : 10 Jun 2021 13:21
Operator : SY/MD
Sample : M2674-01
Misc : 5.79G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TMR-DS-01-060921

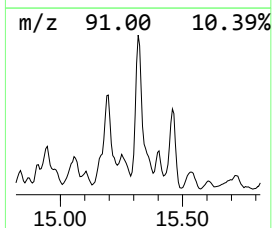
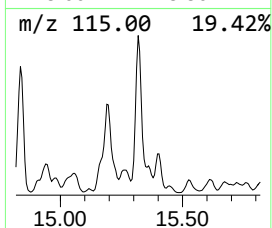
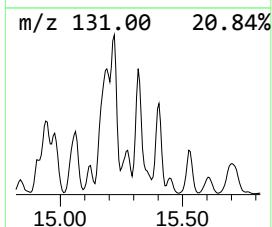
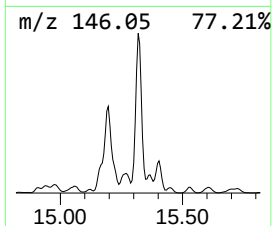
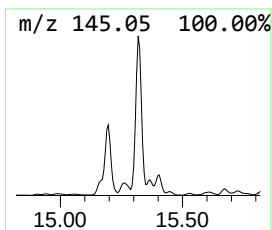
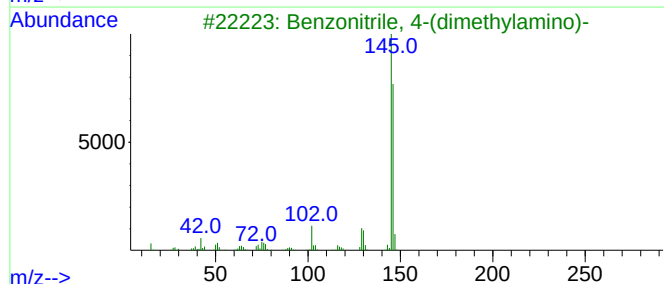
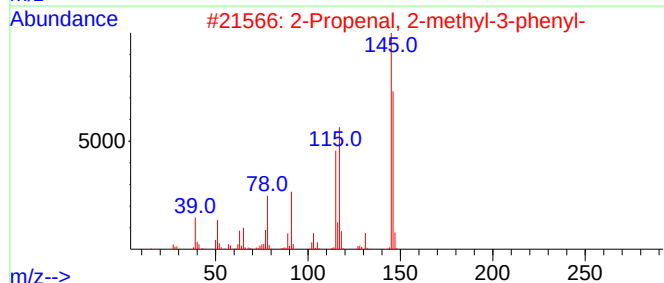
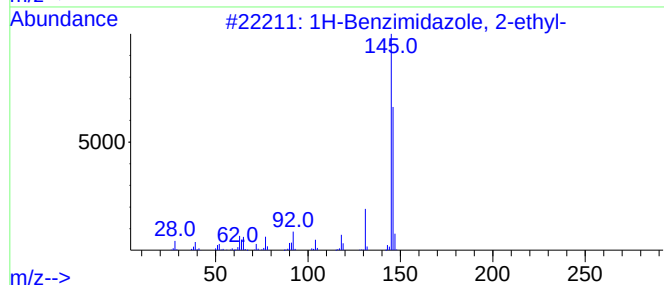
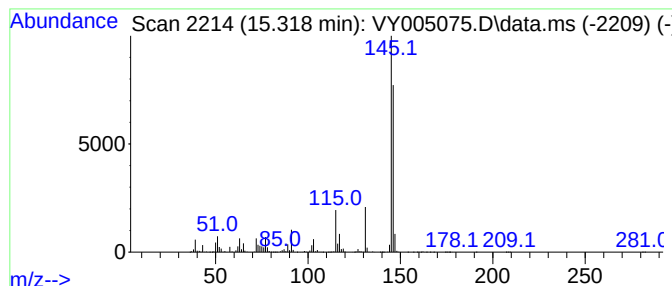
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Quant Title : SW846 8260

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

Peak Number 8 1H-Benzimidazole, 2-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.318	83.37 ug/l	7117210	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Benzimidazole, 2-ethyl-	146	C9H10N2	001848-84-6	72
2			2-Propenal, 2-methyl-3-phenyl-	146	C10H10O	000101-39-3	59
3			Benzonitrile, 4-(dimethylamino)-	146	C9H10N2	001197-19-9	59
4			Benzofuran, 4,7-dimethyl-	146	C10H10O	028715-26-6	58
5			1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY061021\
Data File : VY005075.D
Acq On : 10 Jun 2021 13:21
Operator : SY/MD
Sample : M2674-01
Misc : 5.79G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TMR-DS-01-060921

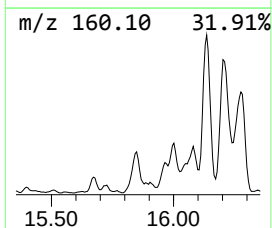
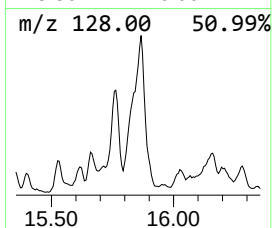
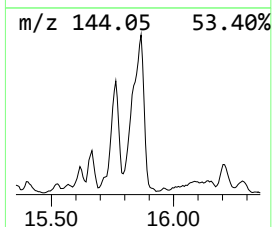
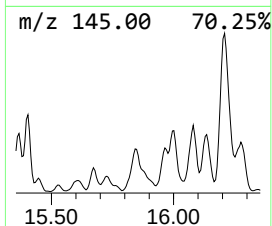
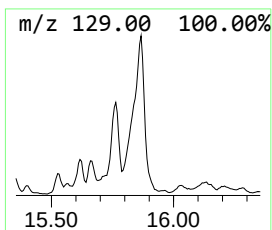
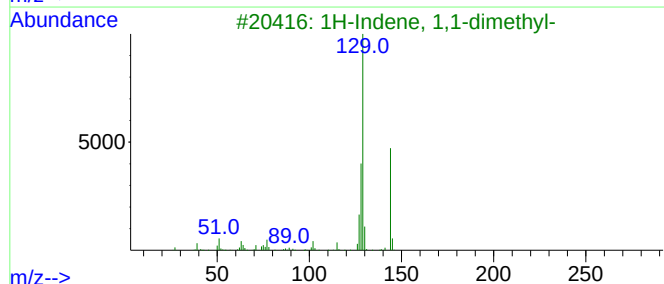
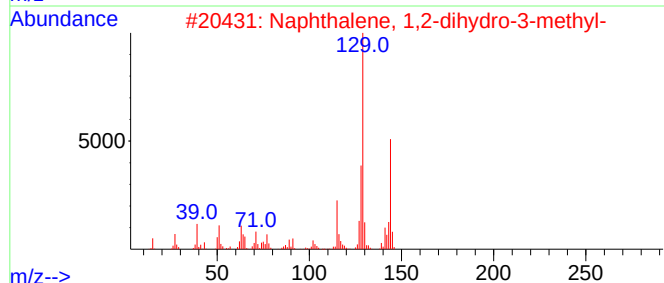
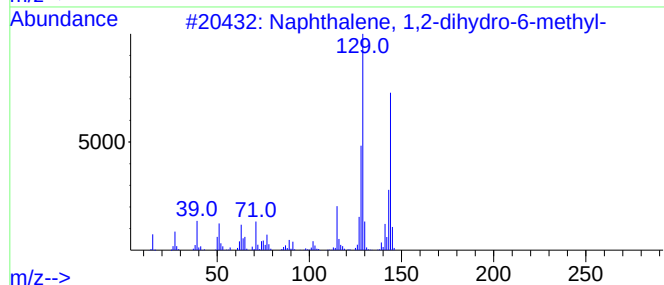
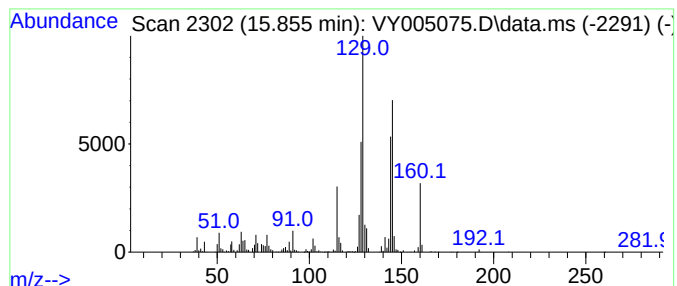
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2-dihydro-6-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.855	27.71 ug/l	2365380	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2-dihydro-6-methyl-	144	C11H12	002717-47-7	76
2			Naphthalene, 1,2-dihydro-3-methyl-	144	C11H12	002717-44-4	76
3			1H-Indene, 1,1-dimethyl-	144	C11H12	018636-55-0	70
4			1H-Indene, 1,3-dimethyl-	144	C11H12	002177-48-2	70
5			1H-Indene, 2,3-dimethyl-	144	C11H12	004773-82-4	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY061021\
Data File : VY005075.D
Acq On : 10 Jun 2021 13:21
Operator : SY/MD
Sample : M2674-01
Misc : 5.79G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TMR-DS-01-060921

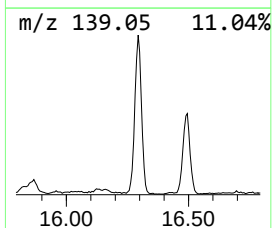
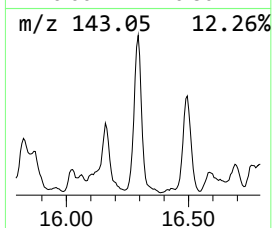
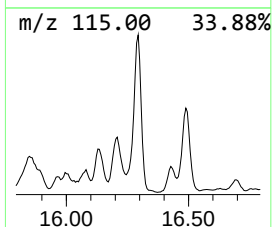
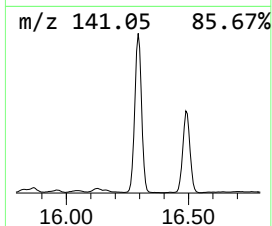
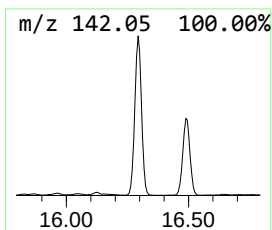
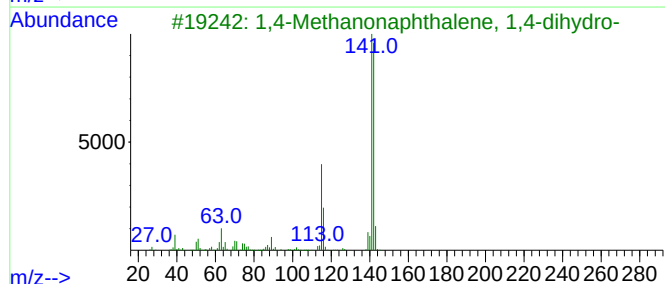
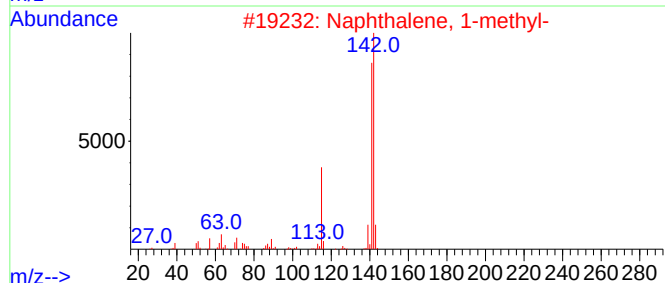
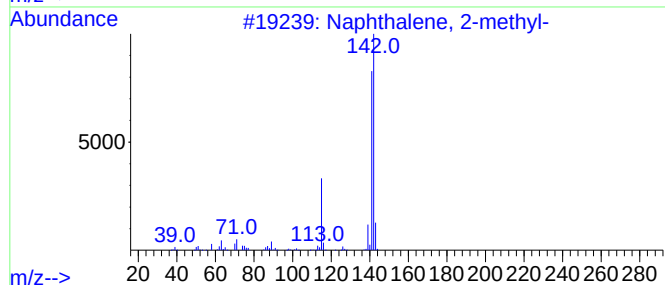
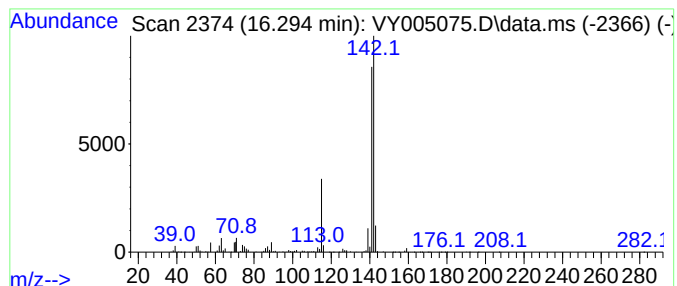
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.294	36.95 ug/l	3154130	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	90
4			Benzocycloheptatriene	142	C11H10	000264-09-5	90
5			3-Indolecarbonitrile	142	C9H6N2	005457-28-3	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY061021\
Data File : VY005075.D
Acq On : 10 Jun 2021 13:21
Operator : SY/MD
Sample : M2674-01
Misc : 5.79G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TMR-DS-01-060921

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y060221S.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...	12.739	54.5	ug/l	4649500	4	13.428	4268620	50.0
Benzene, 1-ethy...	12.971	41.9	ug/l	3578350	4	13.428	4268620	50.0
Indane	13.630	36.5	ug/l	3112840	4	13.428	4268620	50.0
Indan, 1-methyl-	14.081	28.2	ug/l	2409590	4	13.428	4268620	50.0
Benzofuran, 2-m...	14.392	128.9	ug/l	11002700	4	13.428	4268620	50.0
1H-Indene, 2,3-...	14.568	44.6	ug/l	3810240	4	13.428	4268620	50.0
Benzene, 2-ethe...	14.684	47.1	ug/l	4022470	4	13.428	4268620	50.0
1H-Benzimidazol...	15.318	83.4	ug/l	7117210	4	13.428	4268620	50.0
Naphthalene, 1,...	15.855	27.7	ug/l	2365380	4	13.428	4268620	50.0
Naphthalene, 1-...	16.294	37.0	ug/l	3154130	4	13.428	4268620	50.0