

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092822\
 Data File : VY010716.D
 Acq On : 28 Sep 2022 15:06
 Operator : KP/MD
 Sample : N4855-01
 Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 OE-02-092822

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\82Y091922S.M
 Title : SW846 8260

Signal : TIC: VY010716.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.131	43	51	53	rBV4	2048	5386	1.56%	0.134%
2	2.516	106	114	118	rBV7	3751	9513	2.76%	0.237%
3	2.692	138	143	148	rVB2	2868	5824	1.69%	0.145%
4	4.710	462	474	486	rBV2	19871	60431	17.53%	1.508%
5	5.734	636	642	651	rBV6	2366	7035	2.04%	0.176%
6	6.972	835	845	860	rBV	5761	17859	5.18%	0.446%
7	7.716	957	967	973	rBV	70051	163797	47.51%	4.088%
8	7.783	973	978	990	rVB	80275	180940	52.49%	4.516%
9	8.136	1026	1036	1046	rBV	67101	148822	43.17%	3.714%
10	8.685	1117	1126	1137	rBV	141290	273126	79.23%	6.816%
11	10.179	1361	1371	1378	rBV	193269	329019	95.44%	8.211%
12	10.941	1490	1496	1504	rBV2	6761	11651	3.38%	0.291%
13	11.490	1579	1586	1595	rBV	201011	332364	96.41%	8.295%
14	11.703	1612	1621	1629	rBV5	4379	10733	3.11%	0.268%
15	12.026	1664	1674	1683	rBV4	6488	14403	4.18%	0.359%
16	12.239	1705	1709	1718	rVB6	2161	4616	1.34%	0.115%
17	12.477	1742	1748	1759	rVB	160494	252821	73.34%	6.310%
18	12.672	1772	1780	1783	rVB2	3683	7164	2.08%	0.179%
19	12.733	1783	1790	1793	rVV	15140	24637	7.15%	0.615%
20	12.764	1793	1795	1798	rVV2	8135	11959	3.47%	0.298%
21	12.812	1798	1803	1808	rVV2	20675	37257	10.81%	0.930%
22	12.855	1808	1810	1815	rVB5	2862	4091	1.19%	0.102%
23	12.965	1820	1828	1833	rBV	12119	23245	6.74%	0.580%
24	13.032	1836	1839	1846	rVB3	7546	11878	3.45%	0.296%
25	13.117	1846	1853	1859	rBV	32736	49309	14.30%	1.231%
26	13.251	1866	1875	1878	rVB7	5165	11026	3.20%	0.275%
27	13.312	1878	1885	1889	rBV3	12526	22352	6.48%	0.558%
28	13.367	1889	1894	1897	rVV7	4490	7250	2.10%	0.181%
29	13.422	1897	1903	1913	rVV2	204601	344734	100.00%	8.603%
30	13.587	1927	1930	1931	rBV2	5440	5503	1.60%	0.137%
31	13.623	1931	1936	1939	rBV2	28457	38421	11.15%	0.959%
32	13.660	1939	1942	1946	rVV3	15918	21179	6.14%	0.529%
33	13.745	1951	1956	1961	rVV3	13564	24008	6.96%	0.599%
34	13.818	1961	1968	1973	rVV	13750	28287	8.21%	0.706%
35	13.879	1974	1978	1981	rVV	15406	26425	7.67%	0.659%
36	13.910	1981	1983	1988	rVV	20148	27431	7.96%	0.685%

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Sampling : 1

Min Area: 3 % of largest Peak

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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\82Y091922S.M
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37	13.965	1988	1992	1997	rVV	28678	43981	12.76%	1.098%
38	14.026	1997	2002	2005	rVV4	6197	9734	2.82%	0.243%
39	14.074	2005	2010	2015	rVB	33341	51558	14.96%	1.287%
40	14.141	2015	2021	2022	rBV4	5647	9640	2.80%	0.241%
41	14.209	2026	2032	2035	rVV4	10539	19504	5.66%	0.487%
42	14.251	2035	2039	2043	rVV6	13907	26505	7.69%	0.661%
43	14.294	2043	2046	2049	rVV	15945	23381	6.78%	0.584%
44	14.330	2049	2052	2056	rVV	26726	35861	10.40%	0.895%
45	14.379	2056	2060	2063	rVV	14129	20416	5.92%	0.510%
46	14.416	2063	2066	2070	rVV3	14975	21111	6.12%	0.527%
47	14.477	2073	2076	2078	rVV3	6222	8728	2.53%	0.218%
48	14.513	2078	2082	2086	rVV3	13526	23135	6.71%	0.577%
49	14.562	2086	2090	2095	rVV2	32702	61488	17.84%	1.535%
50	14.611	2095	2098	2102	rVV	18393	25120	7.29%	0.627%
51	14.678	2102	2109	2115	rVV4	71686	150984	43.80%	3.768%
52	14.733	2115	2118	2125	rVV2	15639	26890	7.80%	0.671%
53	14.830	2128	2134	2142	rVV	59296	98686	28.63%	2.463%
54	14.904	2142	2146	2147	rVV2	8603	11307	3.28%	0.282%
55	14.940	2147	2152	2155	rVV4	29866	57072	16.56%	1.424%
56	14.977	2155	2158	2164	rVV2	16146	32929	9.55%	0.822%
57	15.050	2164	2170	2177	rVB	29070	65191	18.91%	1.627%
58	15.233	2190	2200	2204	rBV6	9712	28965	8.40%	0.723%
59	15.288	2204	2209	2214	rBV2	27848	46012	13.35%	1.148%
60	15.342	2214	2218	2219	rBV2	14264	17155	4.98%	0.428%
61	15.434	2230	2233	2240	rVB8	9117	16877	4.90%	0.421%
62	15.519	2240	2247	2256	rBV3	22984	49620	14.39%	1.238%
63	15.599	2258	2260	2265	rVB6	4704	5726	1.66%	0.143%
64	15.690	2266	2275	2276	rBV5	17766	36023	10.45%	0.899%
65	15.727	2276	2281	2289	rVB	63190	117829	34.18%	2.941%
66	15.812	2291	2295	2300	rVB7	6988	12565	3.64%	0.314%
67	15.879	2300	2306	2313	rBV3	13460	32200	9.34%	0.804%
68	16.031	2324	2331	2337	rBV2	37732	75134	21.79%	1.875%
69	16.226	2353	2363	2368	rBV2	29775	69089	20.04%	1.724%
70	16.287	2368	2373	2386	rVB5	18701	50852	14.75%	1.269%
71	16.489	2399	2406	2411	rBV5	14236	36449	10.57%	0.910%
72	16.537	2412	2414	2419	rVV4	7070	11435	3.32%	0.285%
73	16.623	2422	2428	2438	rVB4	9443	23261	6.75%	0.581%

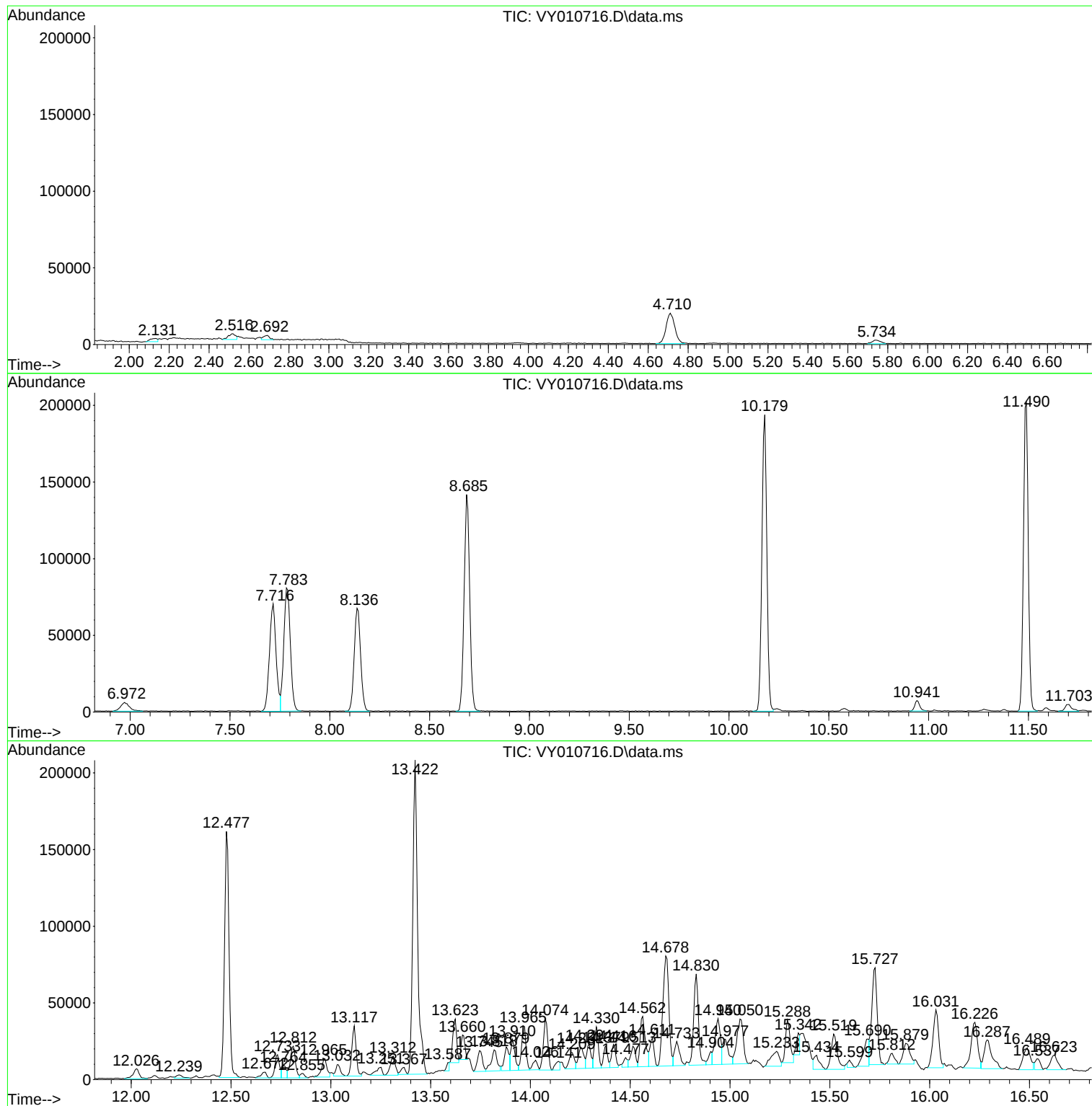
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ClientSampleId :
OE-02-092822

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y091922S.M
Quant Title : SW846 8260

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



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ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
OE-02-092822

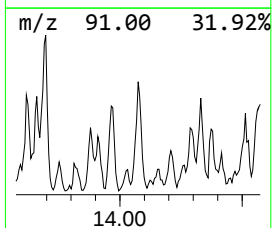
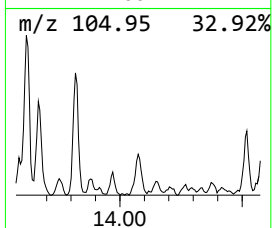
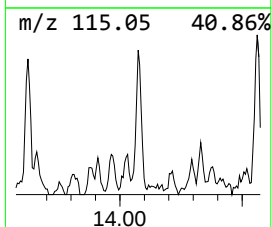
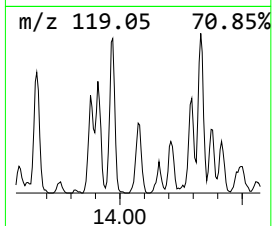
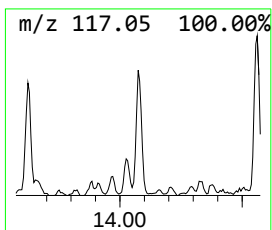
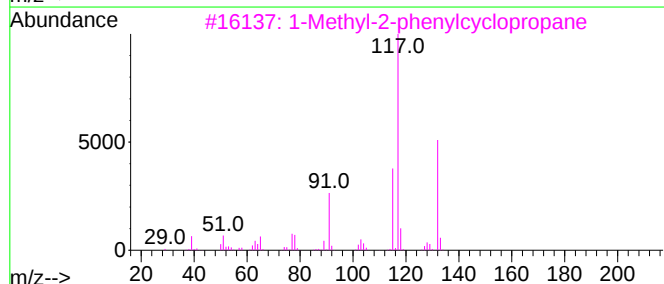
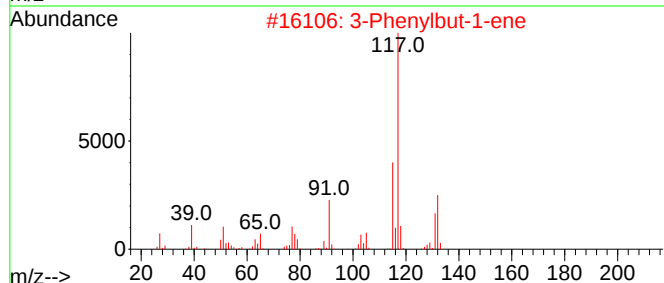
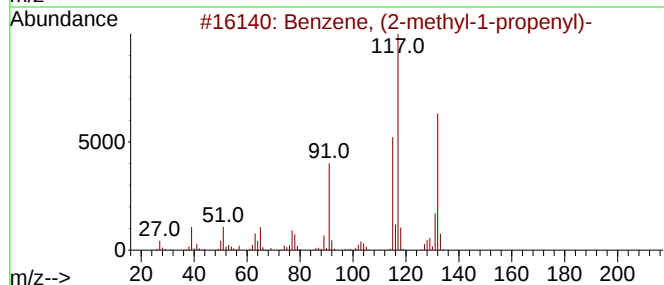
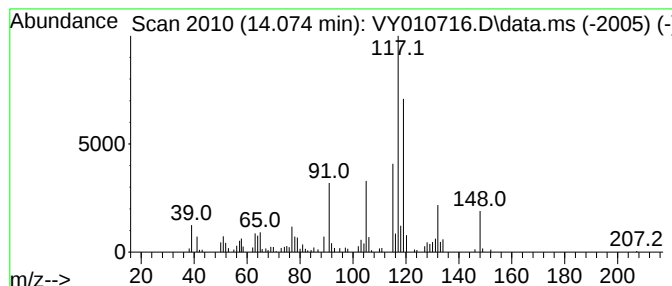
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y091922S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, (2-methyl-1-propen... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.074	7.48 ug/l	51558	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	64
3			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	60
4			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	60
5			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	52



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Instrument :
MSVOA_Y
ClientSampleId :
OE-02-092822

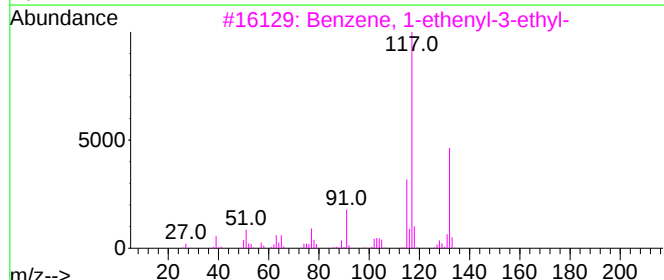
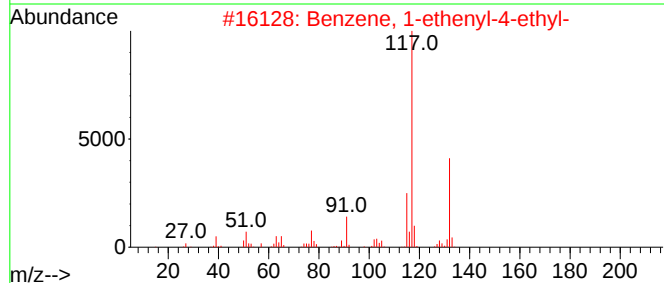
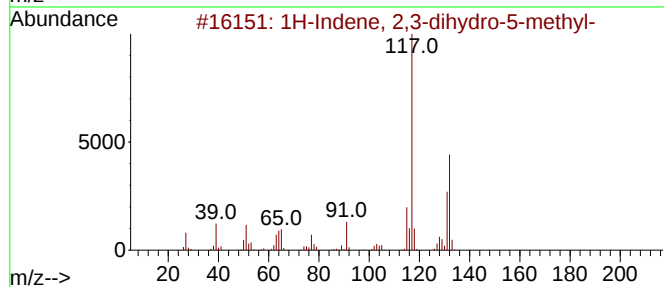
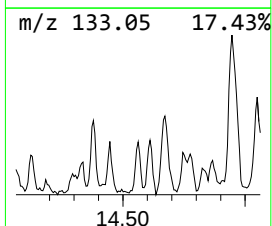
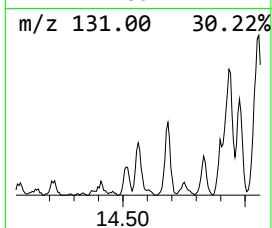
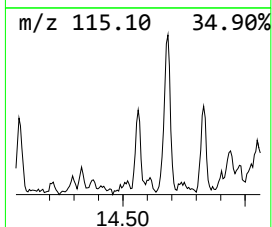
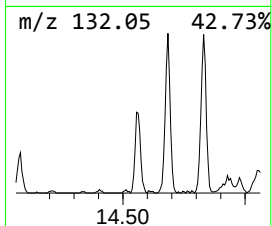
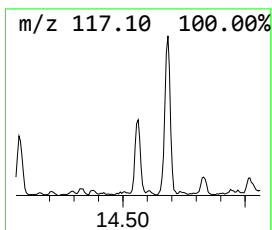
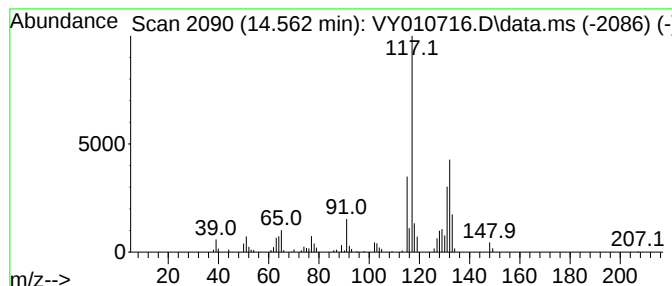
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y091922S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.562	8.92 ug/l	61488	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
4			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	80
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	72



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ClientSampleId :
OE-02-092822

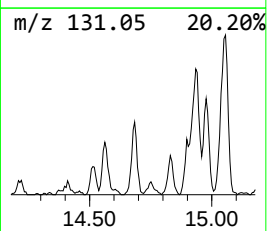
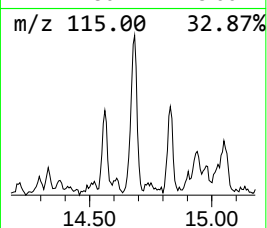
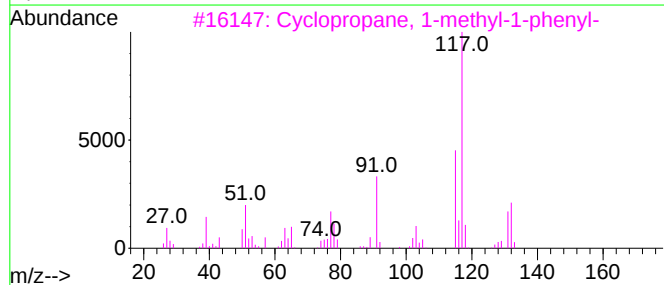
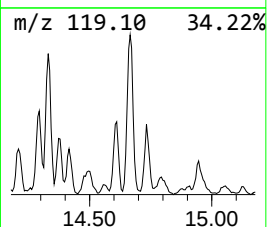
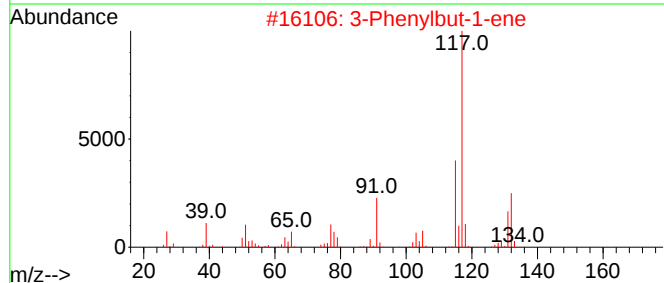
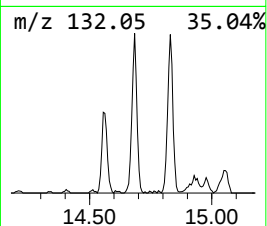
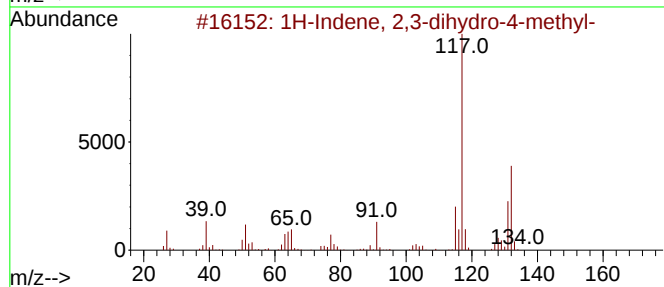
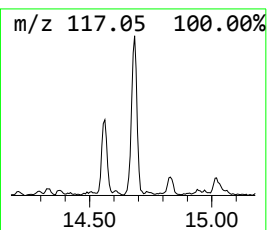
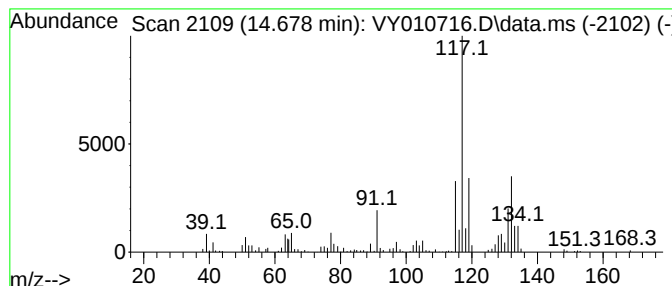
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.678	21.90 ug/l	150984	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	81
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
3			Cyclopropane, 1-methyl-1-phenyl-	132	C10H12	002214-14-4	76
4			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	70
5			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	70



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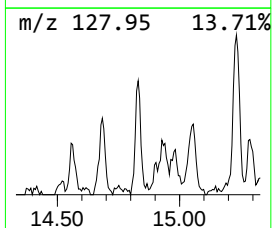
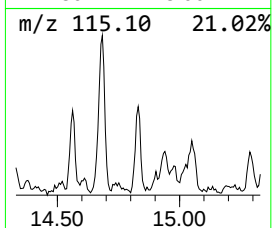
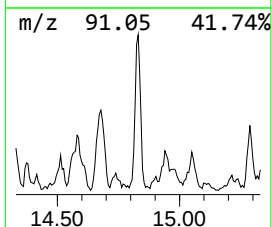
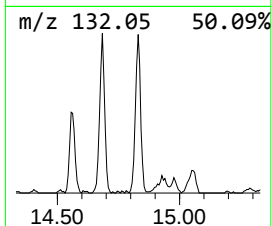
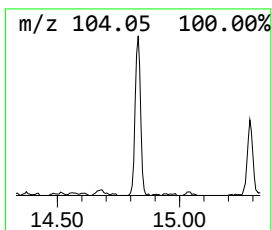
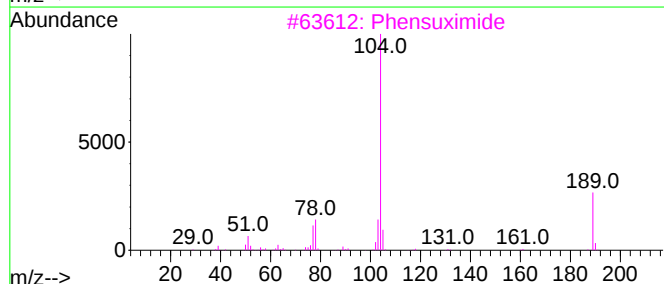
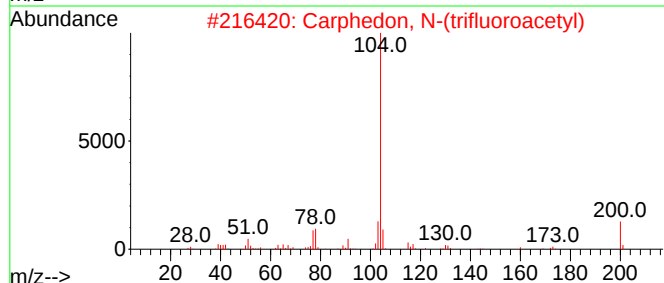
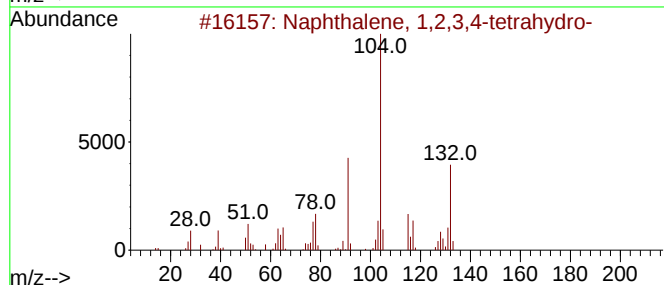
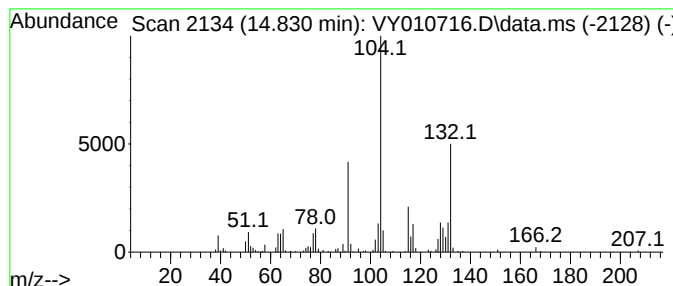
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.830	14.31 ug/l	98686	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	91
2			Carphedon, N-(trifluoroacetyl)	314	C14H13F3N2O3	1000445-42-6	38
3			Phensuximide	189	C11H11NO2	000086-34-0	35
4			3(2H)-Furanone, dihydro-2,2-dime...	190	C12H14O2	063678-00-2	35
5			Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	007694-30-6	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092822\
Data File : VY010716.D
Acq On : 28 Sep 2022 15:06
Operator : KP/MD
Sample : N4855-01
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
OE-02-092822

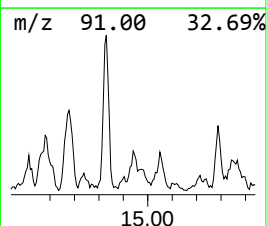
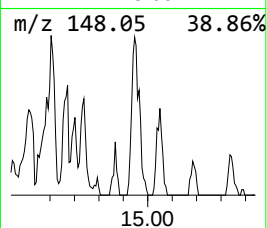
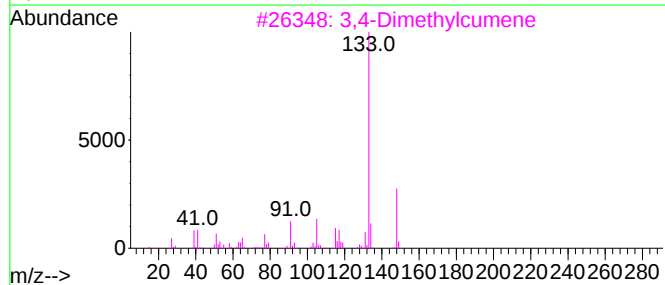
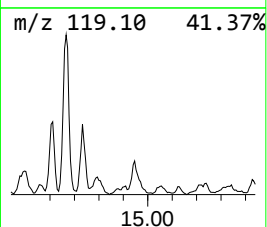
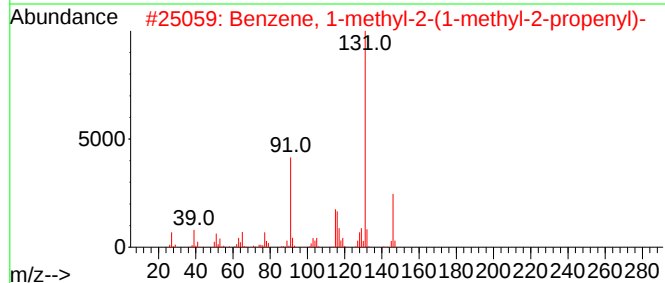
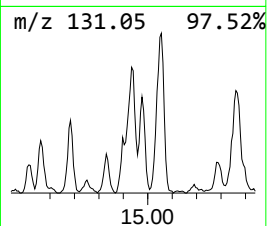
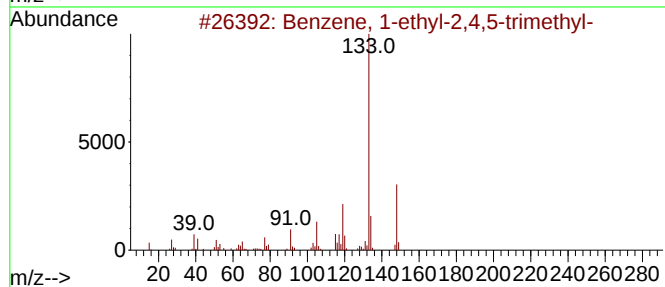
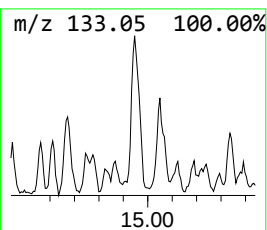
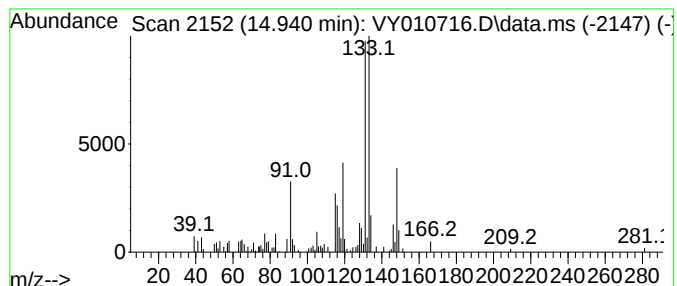
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Quant Title : SW846 8260

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

Peak Number 5 unknown14.940 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.940	8.28 ug/l	57072	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,4,5-trimethyl-	148	C11H16	017851-27-3	46
2			Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	42
3			3,4-Dimethylcumene	148	C11H16	1000370-34-1	38
4			Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	35
5			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092822\
Data File : VY010716.D
Acq On : 28 Sep 2022 15:06
Operator : KP/MD
Sample : N4855-01
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
OE-02-092822

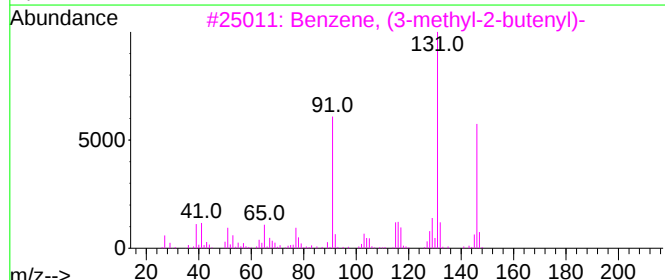
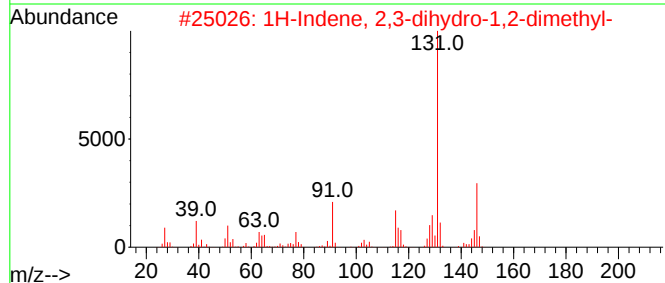
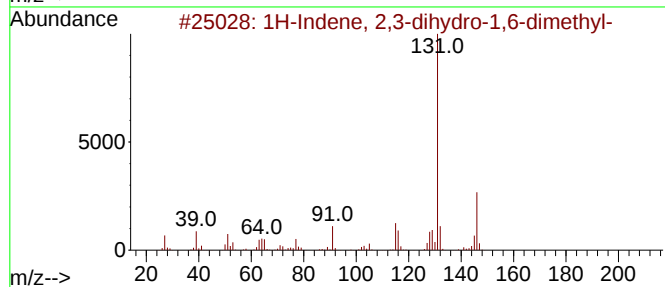
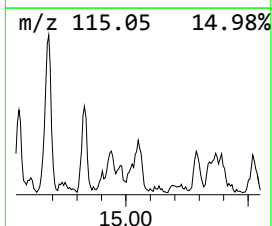
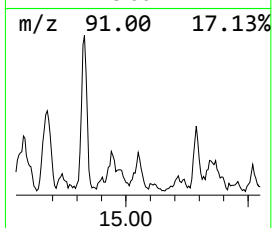
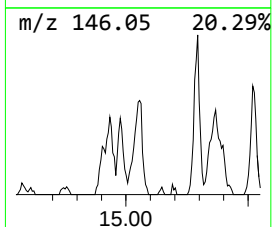
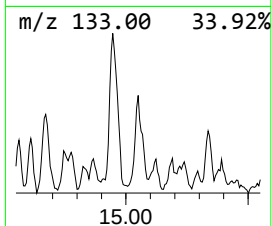
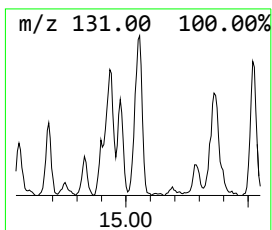
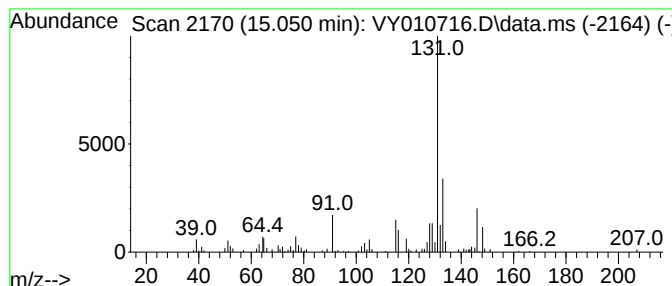
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.050	9.46 ug/l	65191	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	81
3			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	81
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81
5			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092822\
Data File : VY010716.D
Acq On : 28 Sep 2022 15:06
Operator : KP/MD
Sample : N4855-01
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
OE-02-092822

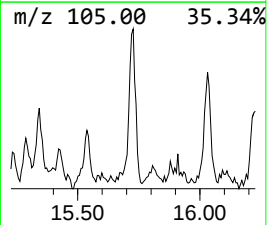
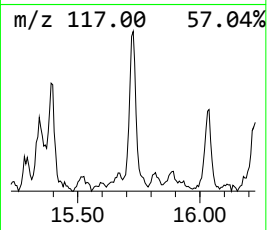
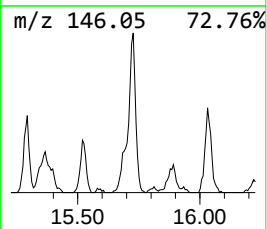
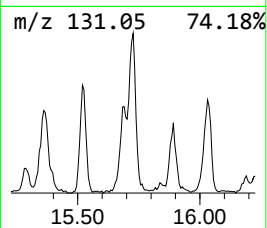
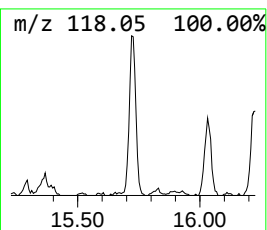
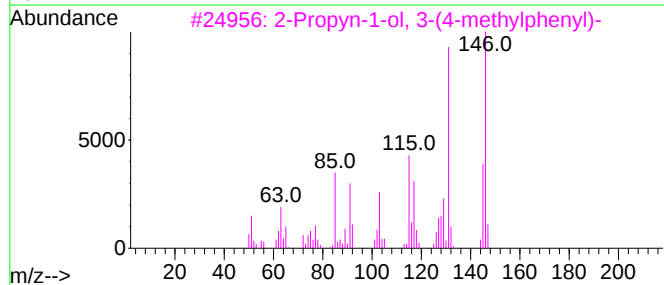
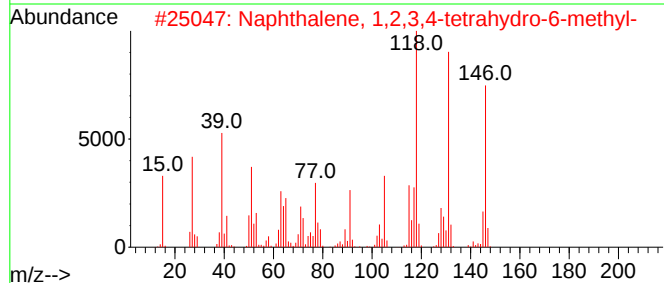
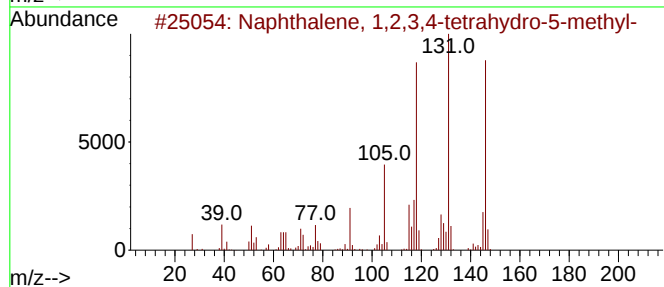
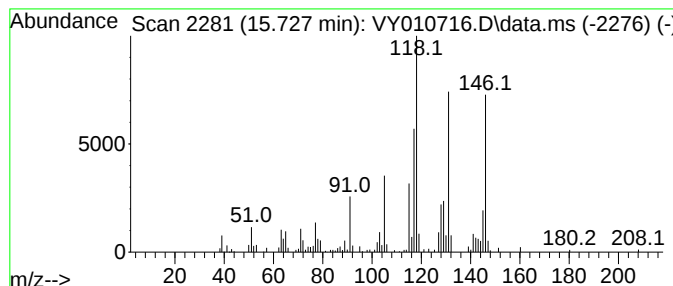
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.727	17.09 ug/l	117829	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	72
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	72
3			2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	55
4			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	55
5			2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1010189-31-0	45



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092822\
Data File : VY010716.D
Acq On : 28 Sep 2022 15:06
Operator : KP/MD
Sample : N4855-01
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
OE-02-092822

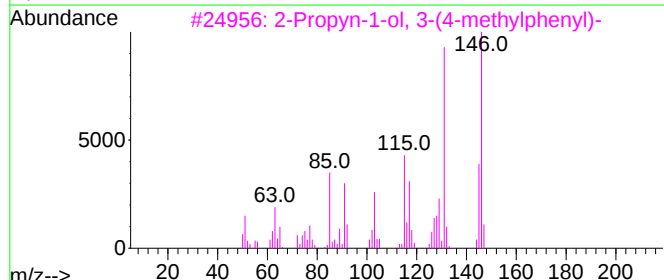
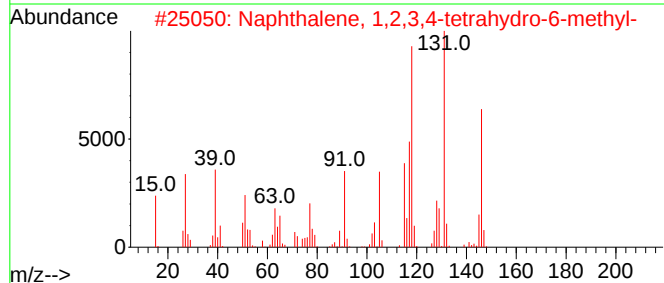
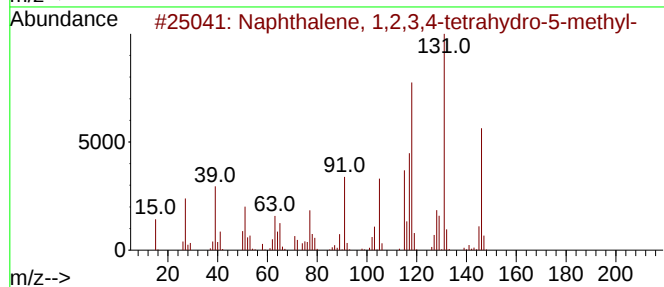
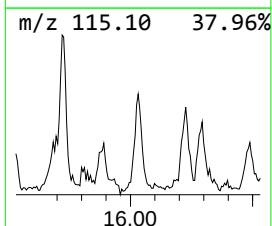
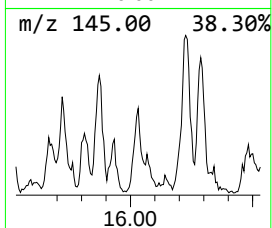
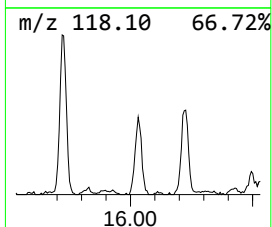
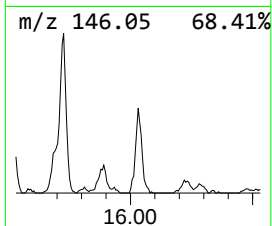
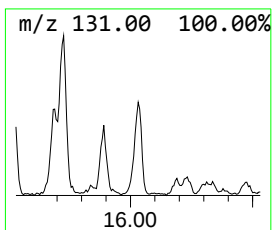
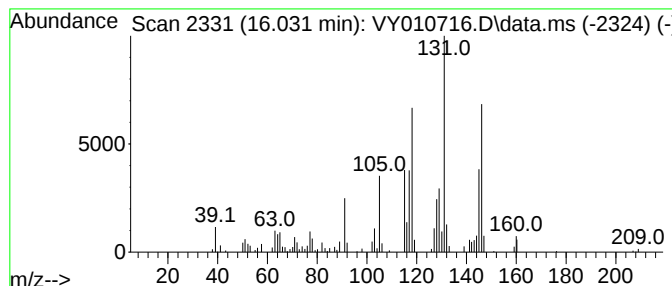
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.031	10.90 ug/l	75134	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	93
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	86
3			2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	68
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	60
5			1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092822\
Data File : VY010716.D
Acq On : 28 Sep 2022 15:06
Operator : KP/MD
Sample : N4855-01
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
OE-02-092822

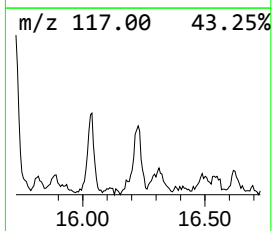
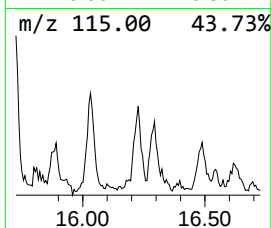
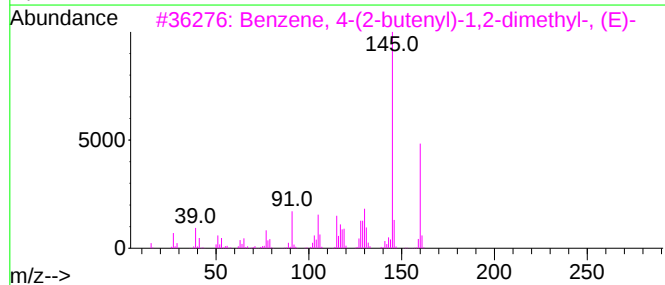
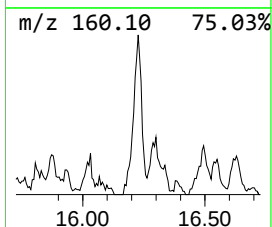
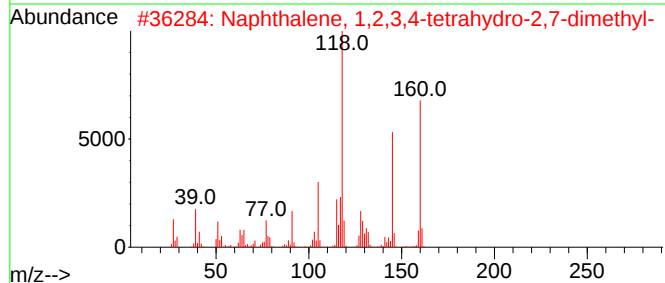
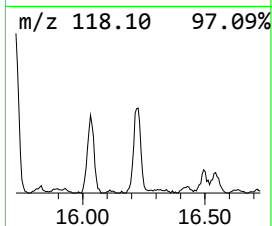
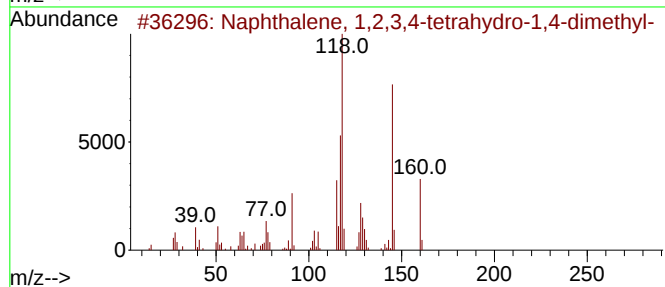
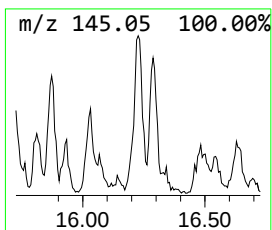
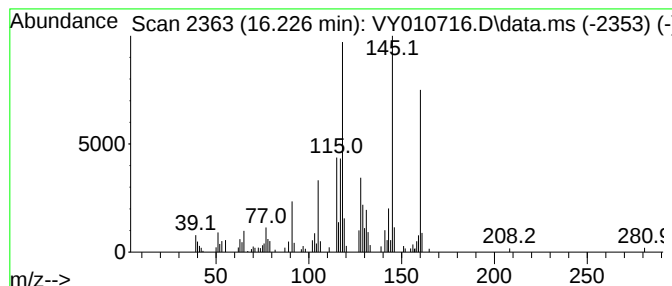
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.227	10.02 ug/l	69089	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	81
2			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	81
3			Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	70
4			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	60
5			Oct-3-ene-1,5-diyne, 3-isopropyl...	160	C12H16	1000211-23-1	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092822\
Data File : VY010716.D
Acq On : 28 Sep 2022 15:06
Operator : KP/MD
Sample : N4855-01
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
OE-02-092822

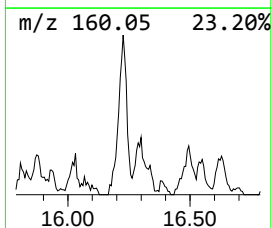
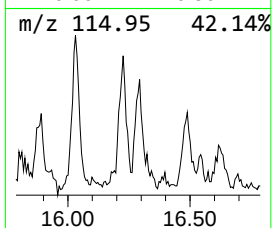
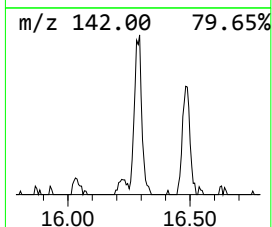
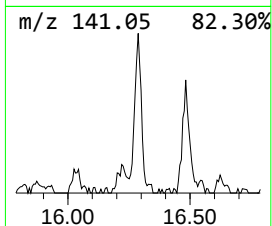
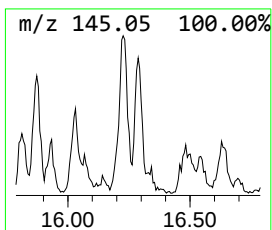
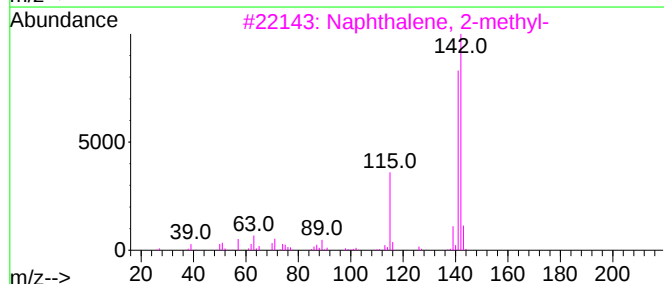
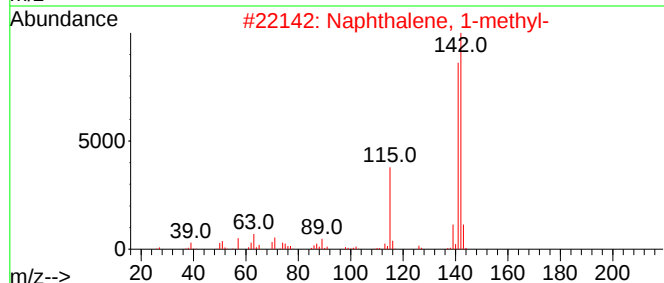
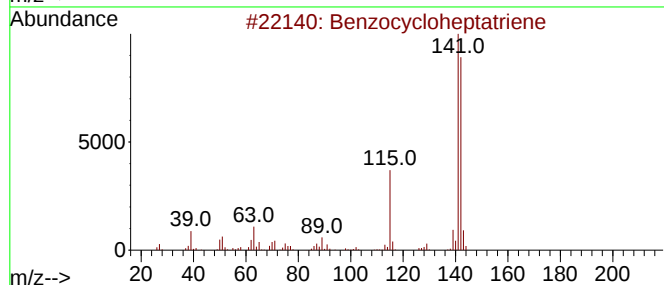
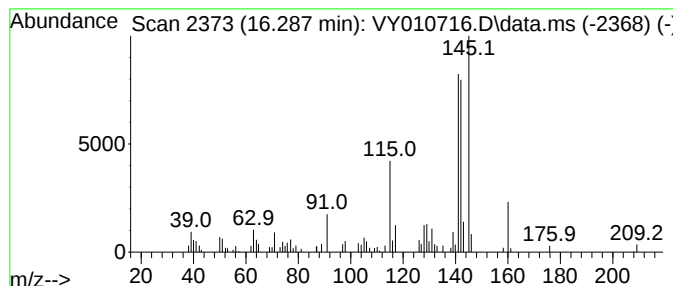
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Quant Title : SW846 8260

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

Peak Number 10 unknown16.287 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.287	7.38 ug/l	50852	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzocycloheptatriene	142	C11H10	000264-09-5	49
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	49
3			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	49
4			Benzeneacetonitrile, 4-cyano-	142	C9H6N2	000876-31-3	30
5			Acetonitrile, 2-(3-cyanophenyl)-	142	C9H6N2	016532-78-8	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092822\
Data File : VY010716.D
Acq On : 28 Sep 2022 15:06
Operator : KP/MD
Sample : N4855-01
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
OE-02-092822

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y091922S.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, (2-met...	14.074	7.5	ug/l	51558	4	13.422	344734	50.0
1H-Indene, 2,3-...	14.562	8.9	ug/l	61488	4	13.422	344734	50.0
1H-Indene, 2,3-...	14.678	21.9	ug/l	150984	4	13.422	344734	50.0
Naphthalene, 1,...	14.830	14.3	ug/l	98686	4	13.422	344734	50.0
unknown14.940	14.940	8.3	ug/l	57072	4	13.422	344734	50.0
1H-Indene, 2,3-...	15.050	9.5	ug/l	65191	4	13.422	344734	50.0
Naphthalene, 1,...	15.727	17.1	ug/l	117829	4	13.422	344734	50.0
Naphthalene, 1,...	16.031	10.9	ug/l	75134	4	13.422	344734	50.0
Naphthalene, 1,...	16.227	10.0	ug/l	69089	4	13.422	344734	50.0
unknown16.287	16.287	7.4	ug/l	50852	4	13.422	344734	50.0