

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101424\
 Data File : VY019893.D
 Acq On : 14 Oct 2024 16:27
 Operator : SY/MD
 Sample : P4395-01
 Misc : 5.24g/5.0mL/MSVOA_Y/SOIL/A
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 F05308-SOLID

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\82Y1009245.M
 Title : SW846 8260

Signal : TIC: VY019893.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.062	55	56	57	rBV	2459	1538	3.20%	0.302%
2	2.092	57	61	63	rBV4	2668	3869	8.04%	0.761%
3	2.220	78	82	86	rVV3	1296	2034	4.23%	0.400%
4	2.251	86	87	90	rVB	1668	1389	2.89%	0.273%
5	2.440	116	118	123	rBV2	1259	1838	3.82%	0.361%
6	2.629	148	149	156	rVB4	1708	3048	6.34%	0.599%
7	2.745	166	168	170	rBV2	1372	1078	2.24%	0.212%
8	3.367	266	270	273	rVB3	974	1361	2.83%	0.268%
9	4.049	378	382	386	rVB3	961	1289	2.68%	0.253%
10	4.110	390	392	394	rBV2	1115	1234	2.57%	0.243%
11	4.287	416	421	423	rVB3	1144	1421	2.95%	0.279%
12	4.549	462	464	467	rVB	920	1068	2.22%	0.210%
13	6.366	757	762	764	rBV2	834	1370	2.85%	0.269%
14	6.451	773	776	779	rVB2	1207	1439	2.99%	0.283%
15	6.543	789	791	794	rBV	900	1136	2.36%	0.223%
16	6.573	794	796	799	rVB2	995	1135	2.36%	0.223%
17	6.860	837	843	844	rVV2	1567	2398	4.99%	0.472%
18	6.878	844	846	851	rVV3	2458	4382	9.11%	0.862%
19	7.640	963	971	976	rBV4	6035	14456	30.06%	2.843%
20	7.707	976	982	989	rVV4	16110	38957	81.01%	7.660%
21	7.768	989	992	994	rVV2	839	1113	2.31%	0.219%
22	8.067	1035	1041	1049	rVB6	4956	12592	26.18%	2.476%
23	8.335	1083	1085	1088	rBV2	1106	1532	3.19%	0.301%
24	8.372	1088	1091	1094	rVV2	1025	1238	2.57%	0.243%
25	8.616	1124	1131	1138	rBV	23138	44556	92.65%	8.761%
26	9.146	1215	1218	1222	rBV	1070	2120	4.41%	0.417%
27	9.487	1270	1274	1280	rBV2	878	1576	3.28%	0.310%
28	9.664	1298	1303	1305	rBV2	944	1211	2.52%	0.238%
29	9.701	1307	1309	1312	rVB	1441	1119	2.33%	0.220%
30	9.878	1336	1338	1343	rVB	1128	1200	2.50%	0.236%
31	9.993	1354	1357	1360	rBV2	997	1063	2.21%	0.209%
32	10.109	1369	1376	1384	rBV	27224	48092	100.00%	9.457%
33	10.274	1400	1403	1407	rBV	1289	1458	3.03%	0.287%
34	10.414	1423	1426	1431	rVB3	667	1148	2.39%	0.226%
35	11.213	1555	1557	1564	rVB6	1380	2291	4.76%	0.451%
36	11.414	1584	1590	1596	rBV3	12433	22998	47.82%	4.522%

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Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
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37	11.487	1599	1602	1607	rVB4	753	1298	2.70%	0.255%
38	11.554	1609	1613	1616	rBV6	1139	1772	3.68%	0.348%
39	11.609	1619	1622	1625	rBV4	1294	1880	3.91%	0.370%
40	11.670	1627	1632	1635	rBV4	1668	2383	4.96%	0.469%
41	11.798	1650	1653	1654	rVV3	1356	1124	2.34%	0.221%
42	11.822	1654	1657	1659	rVB2	877	1132	2.35%	0.223%
43	11.920	1667	1673	1674	rBV5	3326	5572	11.59%	1.096%
44	12.048	1693	1694	1697	rVB2	2367	1717	3.57%	0.338%
45	12.091	1697	1701	1702	rBV4	1520	1098	2.28%	0.216%
46	12.127	1702	1707	1709	rBV4	4273	6702	13.94%	1.318%
47	12.231	1718	1724	1725	rBV4	2369	3630	7.55%	0.714%
48	12.298	1732	1735	1736	rBV3	1774	1692	3.52%	0.333%
49	12.316	1736	1738	1740	rVV3	2102	1807	3.76%	0.355%
50	12.347	1740	1743	1747	rVB4	1940	2480	5.16%	0.488%
51	12.408	1749	1753	1757	rBV5	5236	9034	18.78%	1.776%
52	12.511	1759	1770	1771	rBV9	6148	17264	35.90%	3.395%
53	12.536	1773	1774	1776	rVV2	5512	4605	9.58%	0.906%
54	12.566	1776	1779	1786	rVB7	8705	17946	37.32%	3.529%
55	12.658	1786	1794	1800	rBV10	6543	20089	41.77%	3.950%
56	12.719	1800	1804	1805	rBV4	4326	5394	11.22%	1.061%
57	12.737	1805	1807	1811	rVB5	3499	4065	8.45%	0.799%
58	12.871	1822	1829	1833	rVB8	4882	8507	17.69%	1.673%
59	12.914	1833	1836	1837	rBV3	1749	1682	3.50%	0.331%
60	12.968	1841	1845	1846	rBV4	2316	3190	6.63%	0.627%
61	12.993	1847	1849	1852	rVB4	2556	2383	4.96%	0.469%
62	13.054	1855	1859	1862	rBV5	2673	4128	8.58%	0.812%
63	13.084	1862	1864	1867	rBV4	3257	4093	8.51%	0.805%
64	13.115	1867	1869	1873	rVB5	3722	4408	9.17%	0.867%
65	13.176	1873	1879	1882	rBV7	3306	6344	13.19%	1.247%
66	13.212	1882	1885	1886	rBV3	2561	2541	5.28%	0.500%
67	13.279	1894	1896	1899	rVB4	2289	2024	4.21%	0.398%
68	13.346	1903	1907	1915	rVB7	5448	11441	23.79%	2.250%
69	13.450	1922	1924	1926	rBV3	998	1272	2.64%	0.250%
70	13.474	1926	1928	1932	rVB5	1787	2465	5.13%	0.485%
71	13.529	1935	1937	1938	rBV	1729	1295	2.69%	0.255%
72	13.621	1947	1952	1954	rBV6	1654	2907	6.04%	0.572%
73	13.670	1954	1960	1968	rVV10	6016	13968	29.04%	2.747%
74	13.773	1972	1977	1984	rVV10	3220	9387	19.52%	1.846%
75	13.834	1984	1987	1989	rVV3	1425	1833	3.81%	0.360%
76	13.859	1989	1991	1992	rVV2	1771	1253	2.61%	0.246%

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77	13.883	1992	1995	2000	rVB7	1395	2629	5.47%	0.517%
78	13.974	2009	2010	2013	rBV3	930	1140	2.37%	0.224%
79	14.182	2037	2044	2045	rBV5	3087	4701	9.78%	0.924%
80	14.310	2061	2065	2067	rBV4	1822	3016	6.27%	0.593%
81	14.346	2067	2071	2072	rVV4	2298	3015	6.27%	0.593%
82	14.395	2075	2079	2084	rVB6	1192	2616	5.44%	0.514%
83	14.505	2089	2097	2106	rVV7	9766	32432	67.44%	6.377%
84	14.651	2118	2121	2126	rVB4	1602	2132	4.43%	0.419%
85	14.785	2142	2143	2149	rVV5	1439	1503	3.13%	0.296%
86	14.956	2170	2171	2174	rVB2	1325	1150	2.39%	0.226%
87	15.029	2180	2183	2185	rVV3	1757	1958	4.07%	0.385%
88	15.060	2185	2188	2191	rVB5	733	1139	2.37%	0.224%
89	15.133	2196	2200	2203	rBV6	1763	2433	5.06%	0.478%
90	15.474	2254	2256	2258	rVV2	1455	1313	2.73%	0.258%
91	15.498	2258	2260	2264	rVB5	1342	1721	3.58%	0.338%
92	15.614	2276	2279	2280	rBV2	2568	2074	4.31%	0.408%
93	15.748	2299	2301	2305	rVB4	923	1142	2.37%	0.225%
94	16.047	2347	2350	2353	rBV3	791	1263	2.63%	0.248%
95	16.181	2368	2372	2373	rBV3	1117	1429	2.97%	0.281%
96	16.364	2399	2402	2403	rBV2	2456	2360	4.91%	0.464%
97	16.480	2419	2421	2424	rVB3	947	1186	2.47%	0.233%
98	16.547	2428	2432	2433	rBV4	1527	2237	4.65%	0.440%
99	16.712	2457	2459	2463	rVB4	2194	2185	4.54%	0.430%
100	16.760	2463	2467	2470	rBV6	1219	2219	4.61%	0.436%

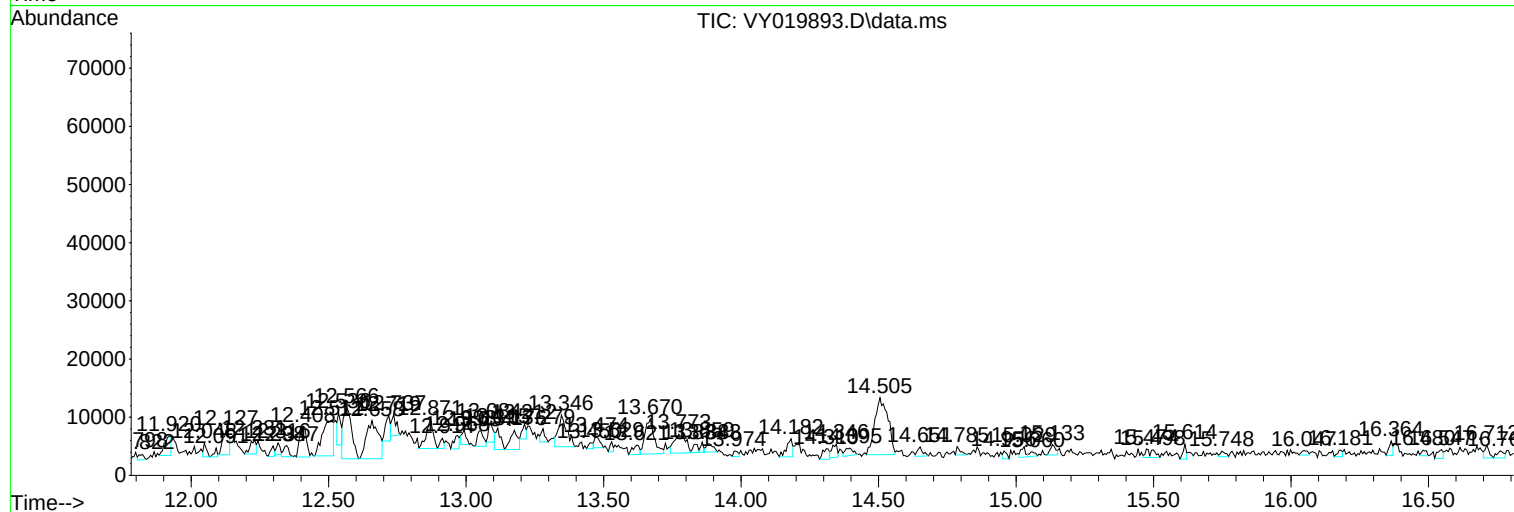
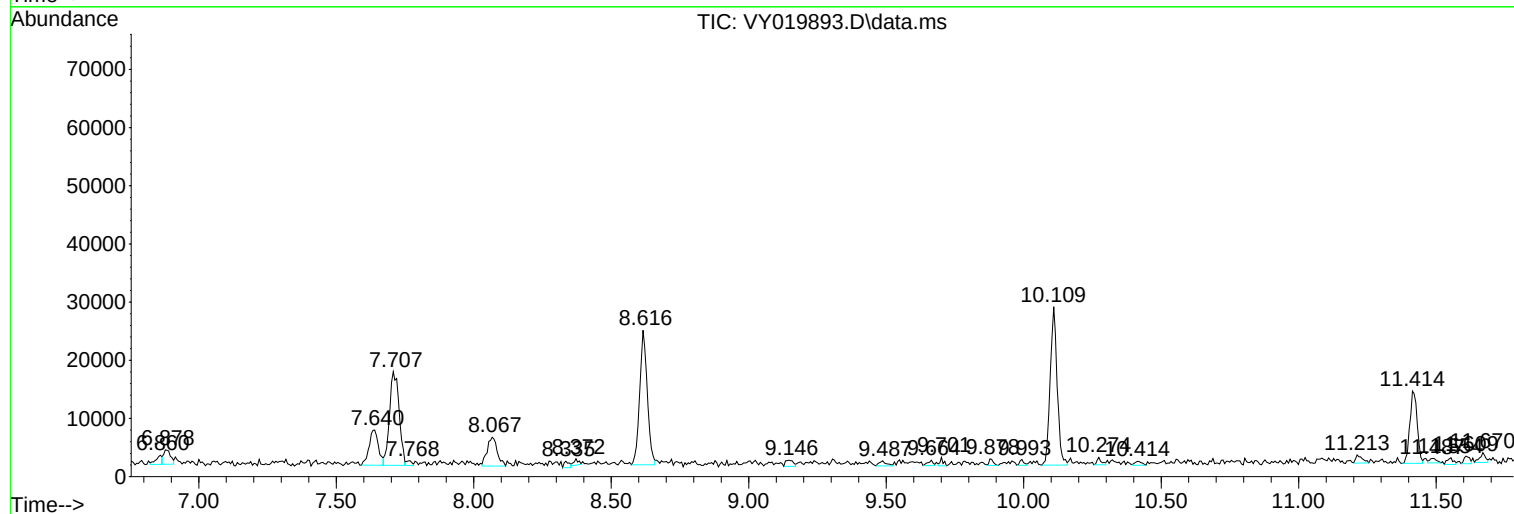
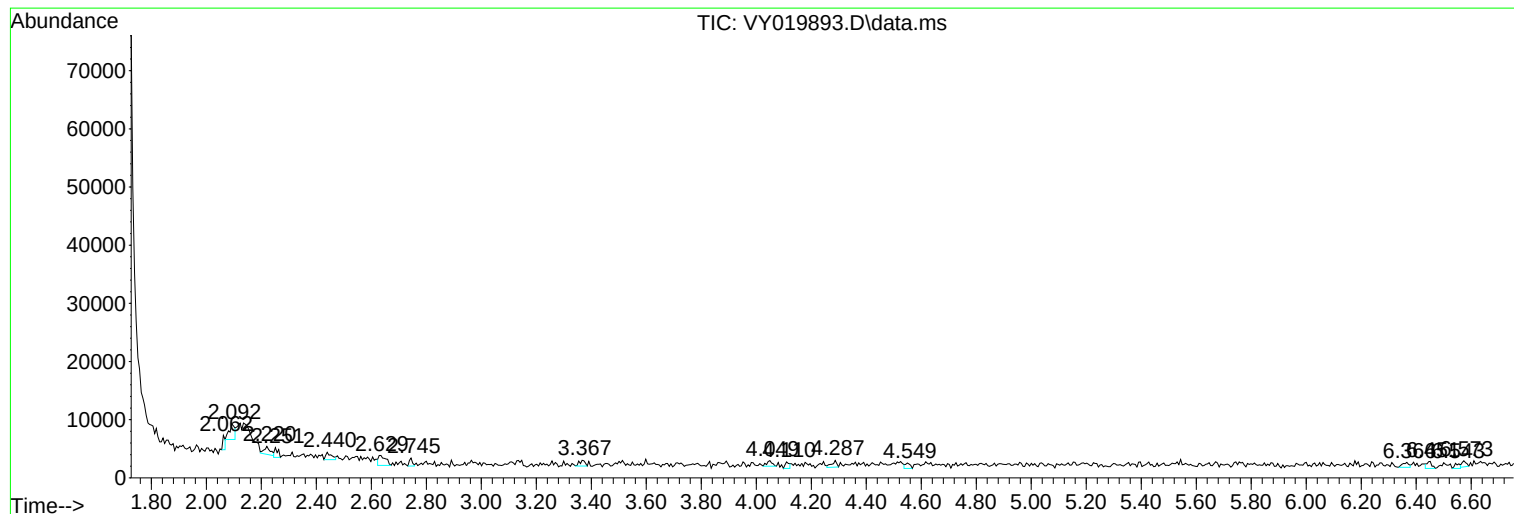
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Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY0101424\
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Acq On : 14 Oct 2024 16:27
Operator : SY/MD
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Misc : 5.24g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
F05308-SOLID

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
Quant Title : SW846 8260

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



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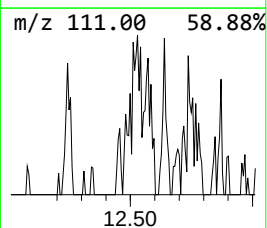
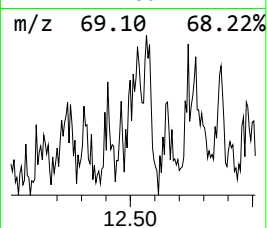
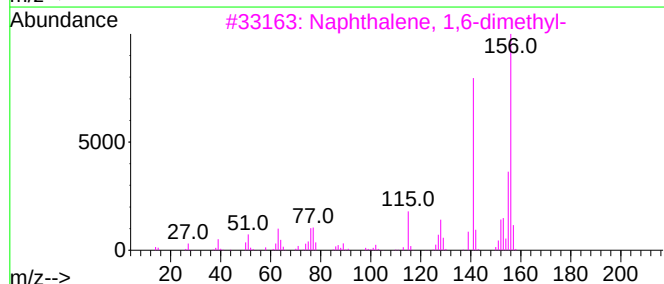
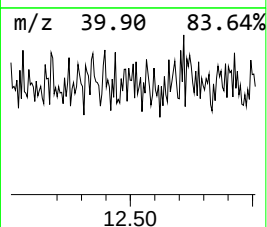
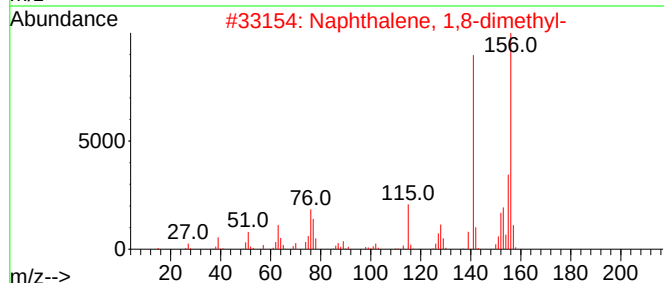
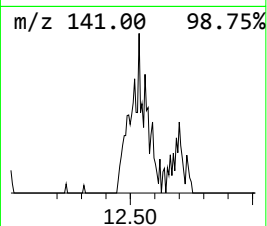
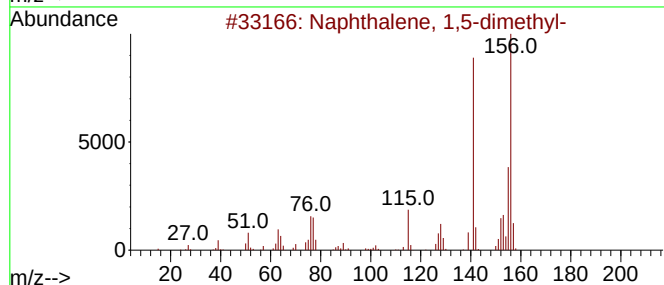
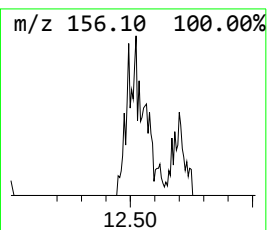
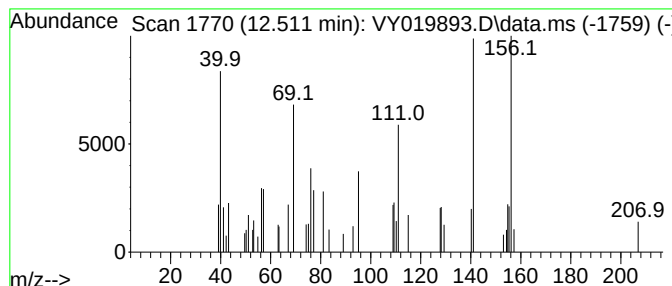
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
Quant Title : SW846 8260

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

Peak Number 1 Naphthalene, 1,5-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.511	75.45 ug/l	17264	1,4-Dichlorobenzene-d4	13.352

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	53
2			Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	49
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	46
4			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	43
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	43



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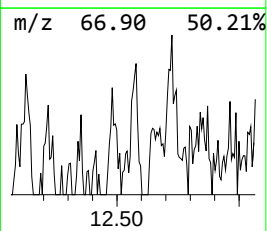
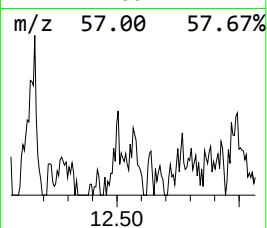
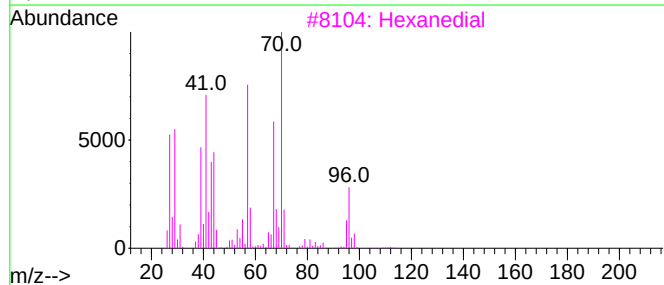
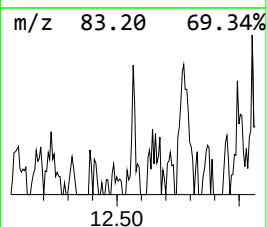
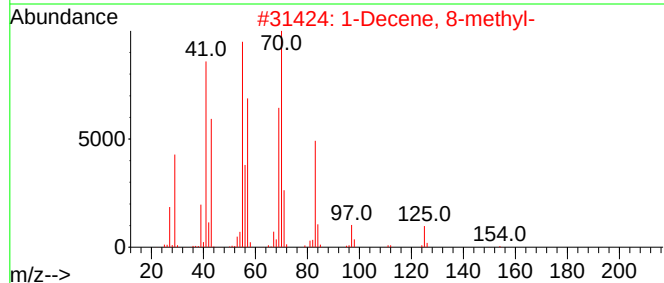
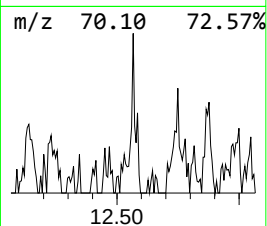
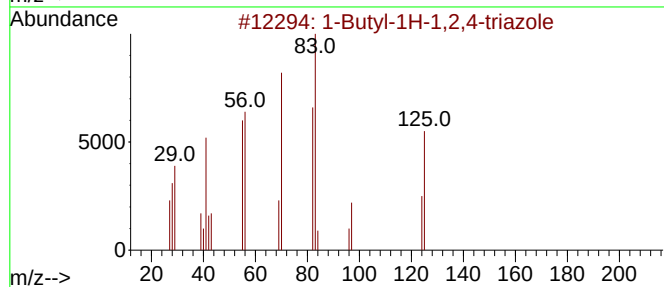
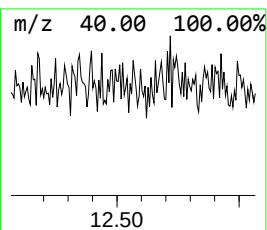
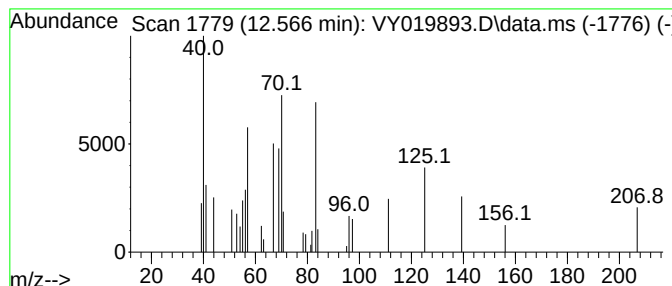
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
Quant Title : SW846 8260

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

Peak Number 2 unknown12.566 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.566	78.43 ug/l	17946	1,4-Dichlorobenzene-d4	13.352

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Butyl-1H-1,2,4-triazole	125	C6H11N3	006086-22-2	38
2			1-Decene, 8-methyl-	154	C11H22	061142-79-8	16
3			Hexanedial	114	C6H10O2	001072-21-5	14
4			3-Undecene, 9-methyl-, (Z)-	168	C12H24	074630-50-5	12
5			4-Undecene, 9-methyl-, (Z)-	168	C12H24	074630-56-1	10



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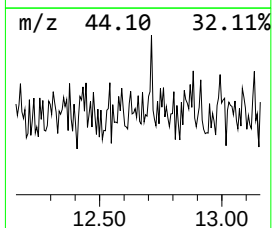
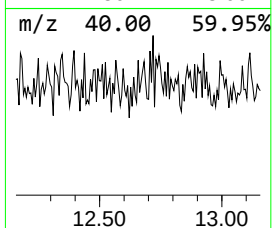
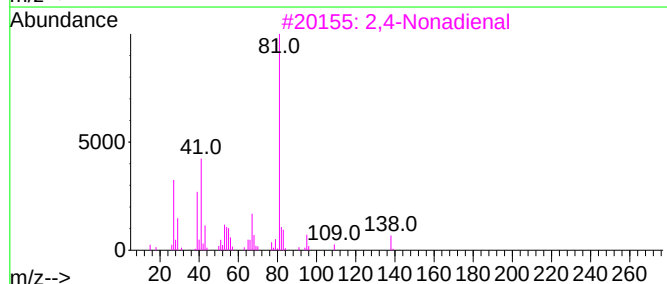
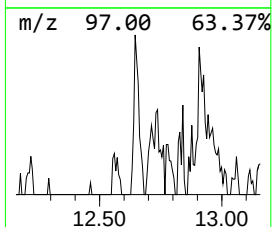
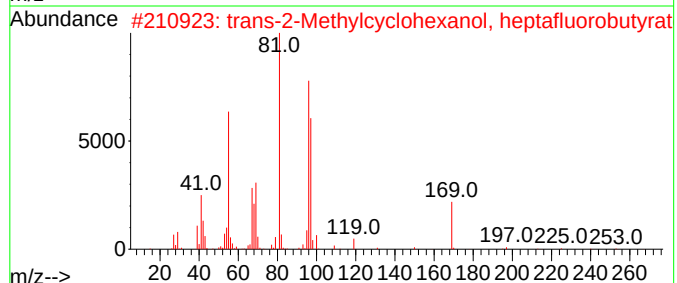
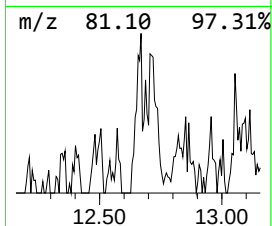
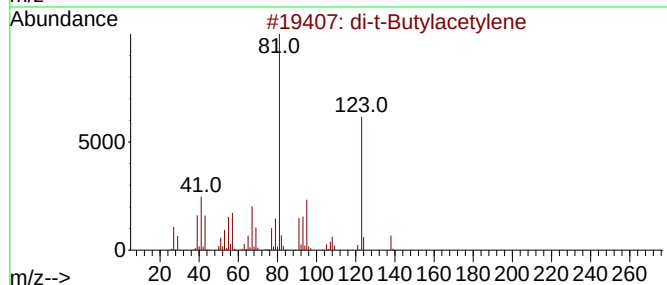
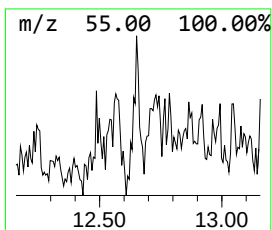
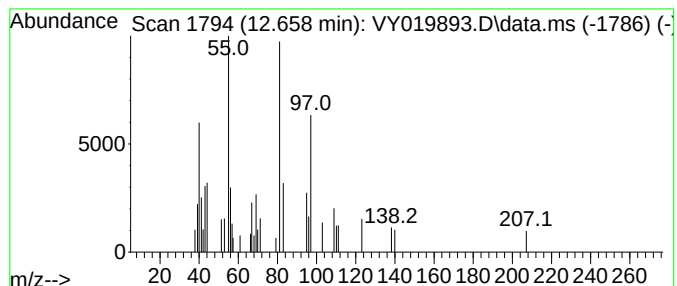
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
Quant Title : SW846 8260

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

Peak Number 3 unknown12.658 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.658	87.79 ug/l	20089	1,4-Dichlorobenzene-d4	13.352

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			di-t-Butylacetylene	138	C10H18	017530-24-4	38
2			trans-2-Methylcyclohexanol, hept...	310	C11H13F7O2	1010364-13-5	37
3			2,4-Nonadienal	138	C9H14O	006750-03-4	35
4			Cyclohexane, ethenyl-	110	C8H14	000695-12-5	35
5			2,4-Heptadienal, (E,E)-	110	C7H10O	004313-03-5	27



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101424\
Data File : VY019893.D
Acq On : 14 Oct 2024 16:27
Operator : SY/MD
Sample : P4395-01
Misc : 5.24g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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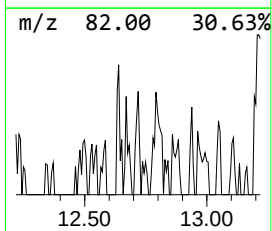
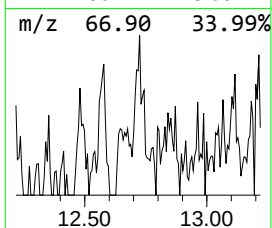
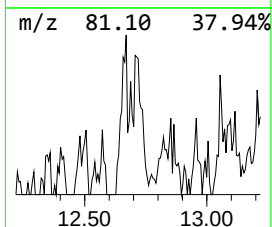
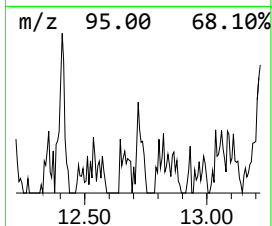
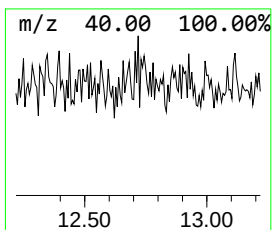
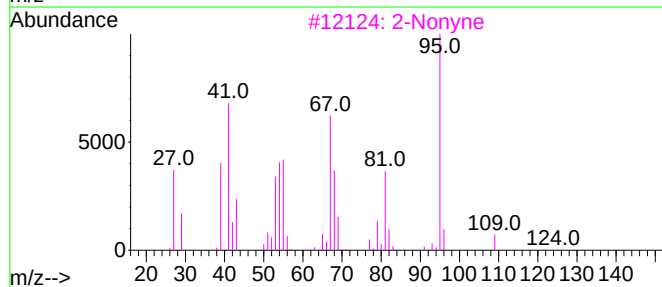
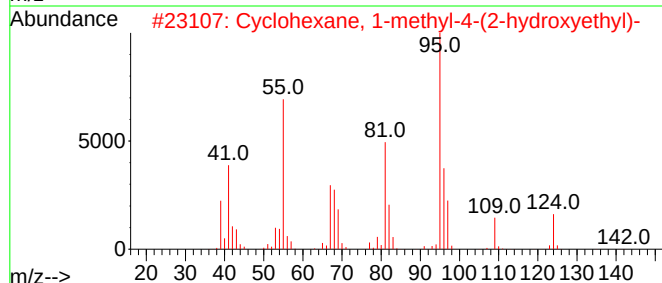
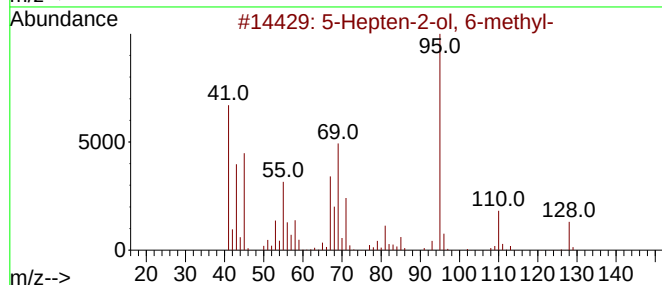
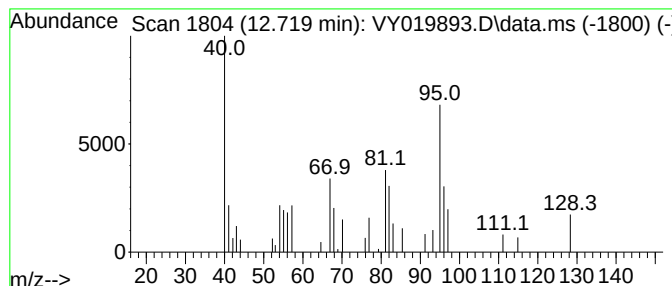
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Quant Title : SW846 8260

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

Peak Number 4 unknown12.719 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.719	23.57 ug/l	5394	1,4-Dichlorobenzene-d4	13.352

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Hepten-2-ol, 6-methyl-	128	C8H16O	001569-60-4	43
2			Cyclohexane, 1-methyl-4-(2-hydro...	142	C9H18O	004916-87-4	40
3			2-Nonyne	124	C9H16	019447-29-1	38
4			1H-Imidazole, 2-ethyl-	96	C5H8N2	001072-62-4	38
5			6-Methylhept-4-en-1-yl 2-methylb...	212	C13H24O2	1215128-05-4	37



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101424\
Data File : VY019893.D
Acq On : 14 Oct 2024 16:27
Operator : SY/MD
Sample : P4395-01
Misc : 5.24g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
F05308-SOLID

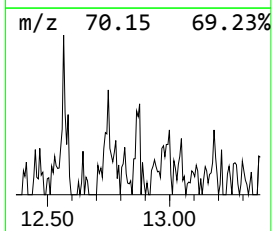
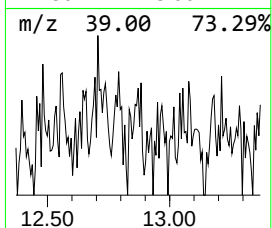
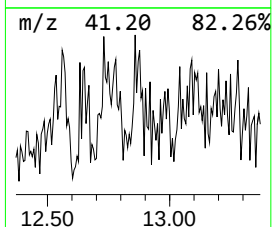
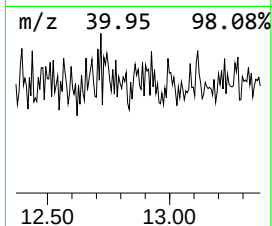
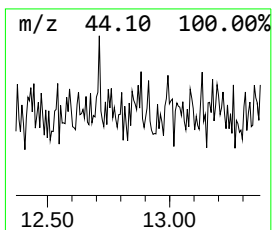
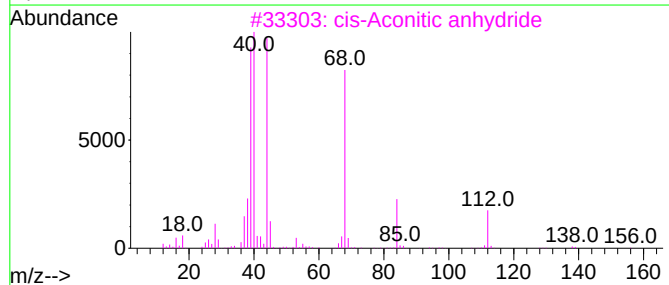
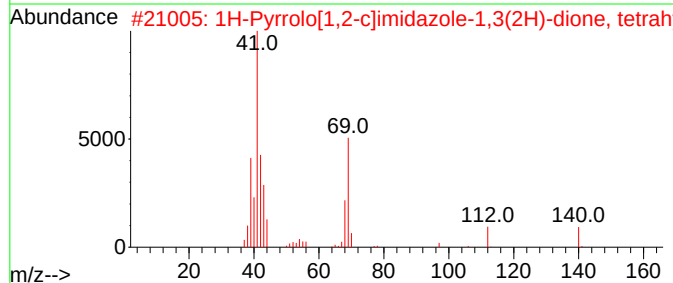
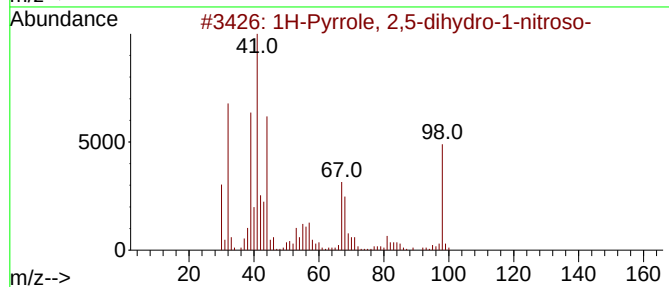
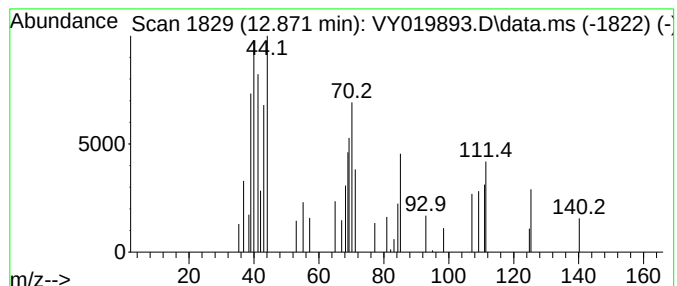
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Quant Title : SW846 8260

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

Peak Number 5 unknown12.871 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.871	37.18 ug/l	8507	1,4-Dichlorobenzene-d4	13.352

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Pyrrole, 2,5-dihydro-1-nitroso-	98	C4H6N2O	010552-94-0	43
2			1H-Pyrrolo[1,2-c]imidazole-1,3(2...	140	C6H8N2O2	1000351-31-8	43
3			cis-Aconitic anhydride	156	C6H4O5	006318-55-4	38
4			2-Butenal, (E)-	70	C4H6O	000123-73-9	38
5			Methacrolein	70	C4H6O	000078-85-3	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101424\
Data File : VY019893.D
Acq On : 14 Oct 2024 16:27
Operator : SY/MD
Sample : P4395-01
Misc : 5.24g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
F05308-SOLID

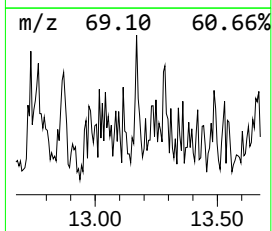
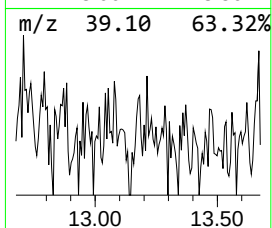
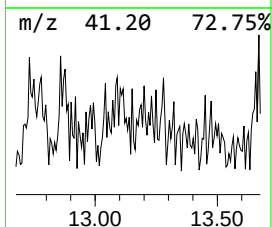
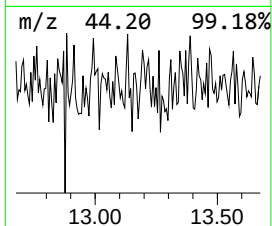
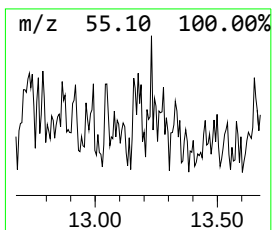
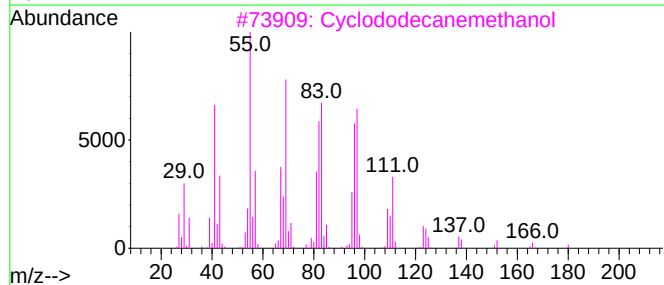
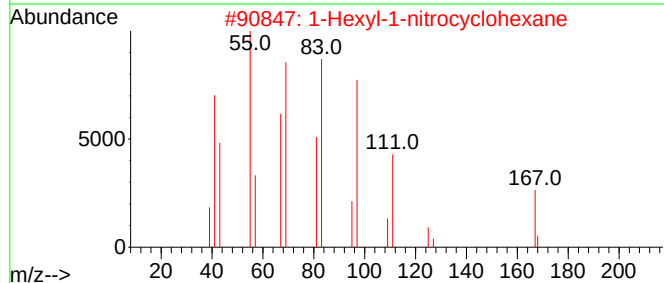
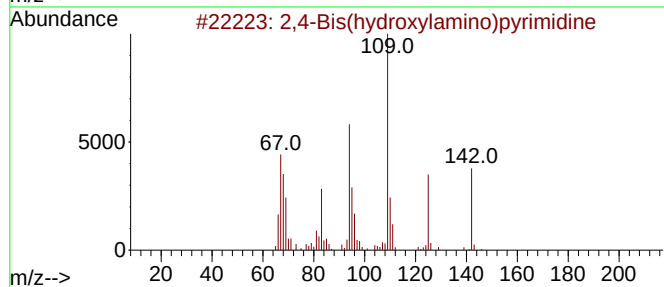
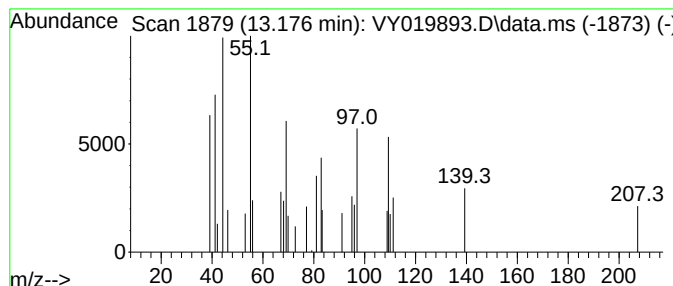
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Quant Title : SW846 8260

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

Peak Number 6 unknown13.176 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.176	27.72 ug/l	6344	1,4-Dichlorobenzene-d4	13.352

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4-Bis(hydroxylamino)pyrimidine	142	C4H6N4O2	154845-38-2	27
2			1-Hexyl-1-nitrocyclohexane	213	C12H23NO2	118252-09-8	22
3			Cyclododecanemethanol	198	C13H26O	001892-12-2	16
4			Cyclopentane, 2-isopropyl-1,3-di...	140	C10H20	032281-85-9	14
5			1-Heneicosanol	312	C21H44O	015594-90-8	12



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101424\
Data File : VY019893.D
Acq On : 14 Oct 2024 16:27
Operator : SY/MD
Sample : P4395-01
Misc : 5.24g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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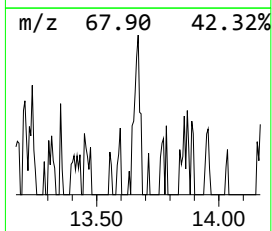
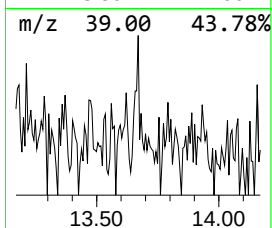
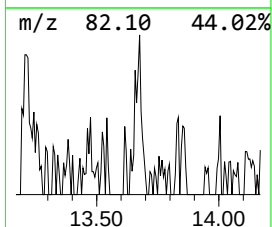
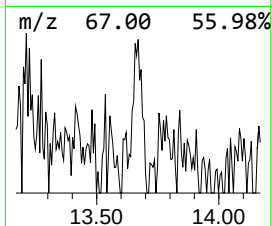
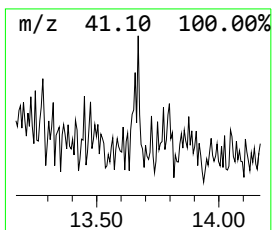
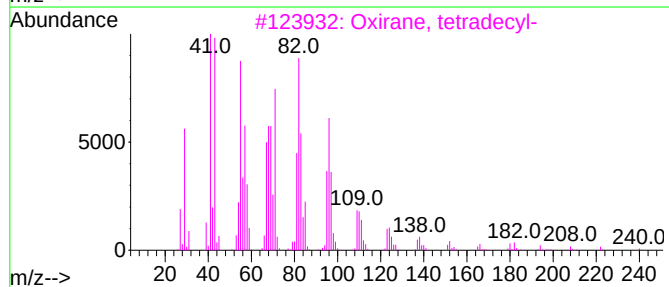
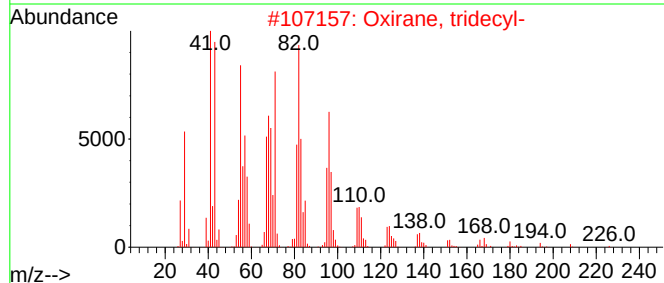
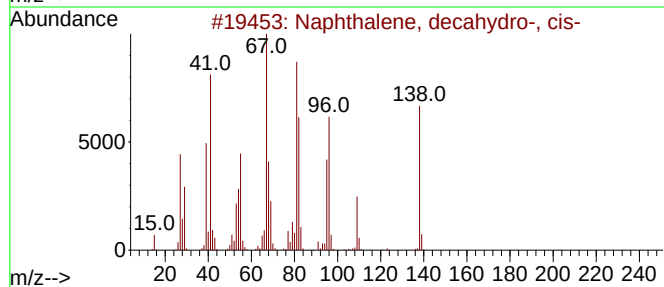
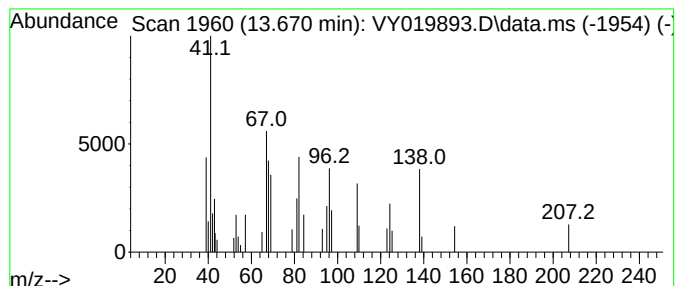
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Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, decahydro-, cis- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.670	61.04 ug/l	13968	1,4-Dichlorobenzene-d4	13.352

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	53
2			Oxirane, tridecyl-	226	C15H30O	018633-25-5	50
3			Oxirane, tetradecyl-	240	C16H32O	007320-37-8	47
4			3-Octyne, 2-methyl-	124	C9H16	055402-15-8	43
5			Naphthalene, decahydro-	138	C10H18	000091-17-8	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101424\
Data File : VY019893.D
Acq On : 14 Oct 2024 16:27
Operator : SY/MD
Sample : P4395-01
Misc : 5.24g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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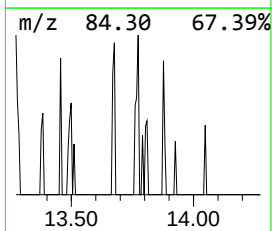
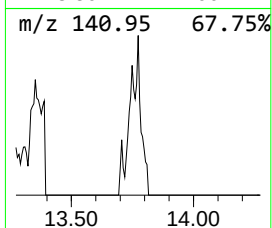
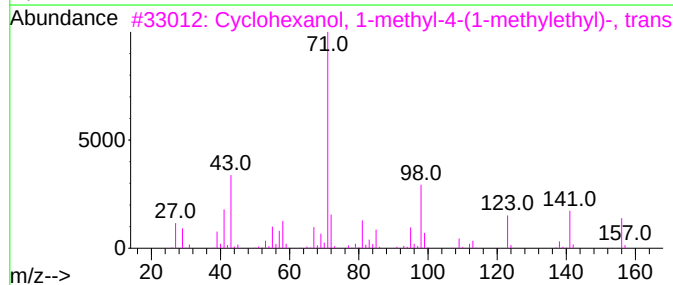
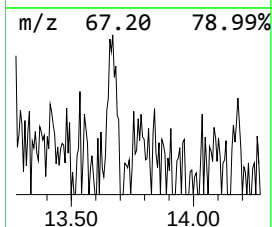
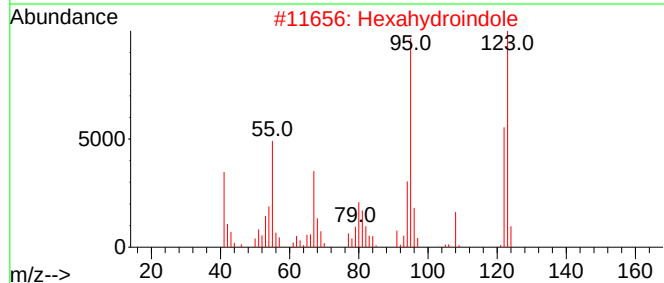
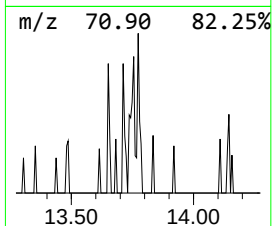
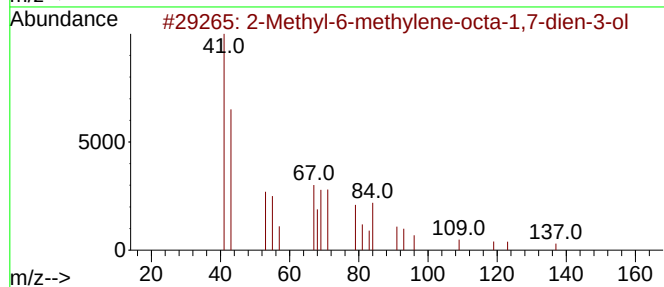
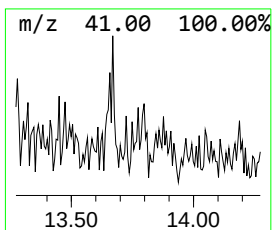
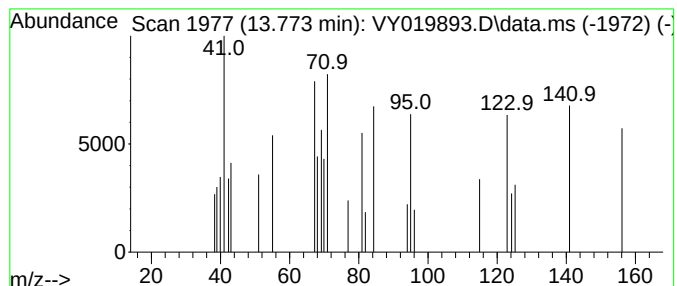
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Quant Title : SW846 8260

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

Peak Number 8 unknown13.773 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.773	41.02 ug/l	9387	1,4-Dichlorobenzene-d4	13.352

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methyl-6-methylene-octa-1,7-di...	152	C10H16O	022459-10-5	38
2			Hexahydroindole	123	C8H13N	018159-32-5	27
3			Cyclohexanol, 1-methyl-4-(1-meth...	156	C10H20O	003901-93-7	16
4			Cyclohexanol, 1-methyl-4-(1-meth...	156	C10H20O	003901-95-9	16
5			2-Butenenitrile	67	C4H5N	004786-20-3	14



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101424\
Data File : VY019893.D
Acq On : 14 Oct 2024 16:27
Operator : SY/MD
Sample : P4395-01
Misc : 5.24g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
F05308-SOLID

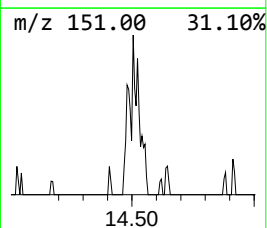
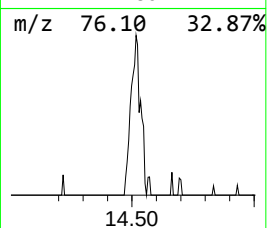
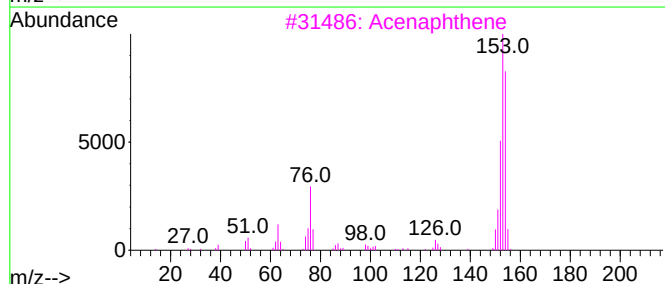
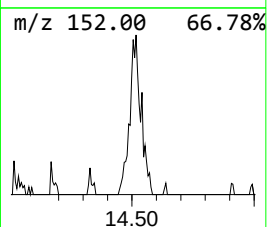
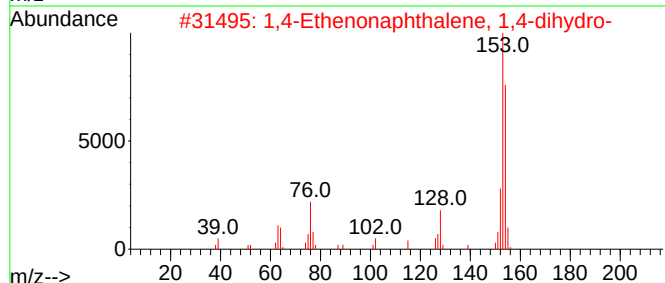
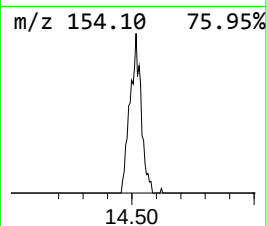
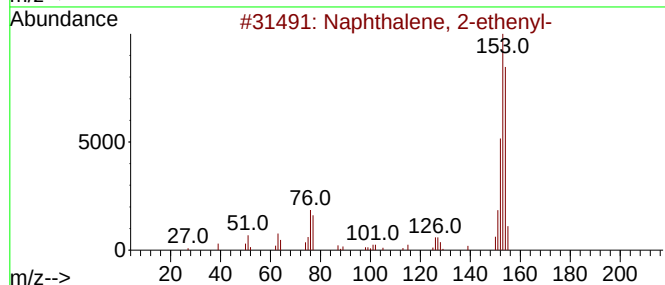
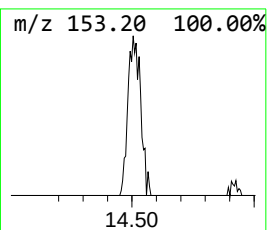
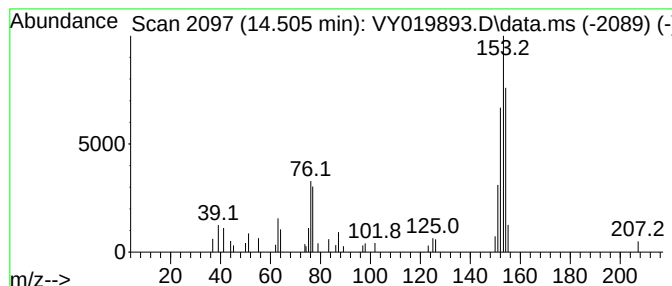
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2-ethenyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.505	141.74 ug/l	32432	1,4-Dichlorobenzene-d4	13.352

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	68
2			1,4-Ethenonaphthalene, 1,4-dihydro-	154	C12H10	007322-47-6	58
3			Acenaphthene	154	C12H10	000083-32-9	52
4			Hexa-2,4-diyn-1-ylbenzene	154	C12H10	000520-74-1	50
5			2,3,4-Trihydroxybenzaldehyde	154	C7H6O4	002144-08-3	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101424\
Data File : VY019893.D
Acq On : 14 Oct 2024 16:27
Operator : SY/MD
Sample : P4395-01
Misc : 5.24g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
F05308-SOLID

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.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
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unknown12.658	12.658	87.8	ug/l	20089	4	13.352	11441	50.0
unknown12.719	12.719	23.6	ug/l	5394	4	13.352	11441	50.0
unknown12.871	12.871	37.2	ug/l	8507	4	13.352	11441	50.0
unknown13.176	13.176	27.7	ug/l	6344	4	13.352	11441	50.0
Naphthalene, de...	13.670	61.0	ug/l	13968	4	13.352	11441	50.0
unknown13.773	13.773	41.0	ug/l	9387	4	13.352	11441	50.0
Naphthalene, 2-...	14.505	141.7	ug/l	32432	4	13.352	11441	50.0