

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041124\
 Data File : VY017913.D
 Acq On : 11 Apr 2024 16:01
 Operator : SY/MD
 Sample : P2072-03
 Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 HR-02-041024

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\82Y032924S.M
 Title : SW846 8260

Signal : TIC: VY017913.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.497	106	111	122	rBV7	4609	16982	2.10%	0.111%
2	4.710	463	474	487	rVB3	5517	16480	2.03%	0.108%
3	7.716	956	967	972	rBV2	73458	169791	20.96%	1.110%
4	7.789	972	979	992	rVB	107375	245619	30.32%	1.605%
5	8.136	1026	1036	1048	rBV	61897	136026	16.79%	0.889%
6	8.685	1118	1126	1141	rBV	174837	342073	42.22%	2.235%
7	10.179	1363	1371	1380	rBV	271269	463724	57.24%	3.030%
8	10.947	1493	1497	1504	rBV6	3633	9486	1.17%	0.062%
9	11.038	1507	1512	1516	rVV2	7288	12095	1.49%	0.079%
10	11.087	1516	1520	1526	rVB3	7459	13119	1.62%	0.086%
11	11.209	1533	1540	1543	rBV4	4695	9091	1.12%	0.059%
12	11.282	1543	1552	1563	rVB6	16605	45379	5.60%	0.297%
13	11.380	1563	1568	1573	rBV2	15520	24989	3.08%	0.163%
14	11.490	1580	1586	1597	rBV	264380	418055	51.60%	2.732%
15	11.593	1598	1603	1606	rBV2	9538	16873	2.08%	0.110%
16	11.715	1612	1623	1630	rBV3	79220	173776	21.45%	1.136%
17	11.782	1630	1634	1639	rVB	16257	27746	3.42%	0.181%
18	11.935	1654	1659	1663	rVB4	6732	10627	1.31%	0.069%
19	12.026	1663	1674	1682	rBV2	79884	204964	25.30%	1.339%
20	12.124	1686	1690	1694	rVB	32977	43079	5.32%	0.282%
21	12.203	1699	1703	1705	rBV3	22927	33408	4.12%	0.218%
22	12.246	1705	1710	1714	rVB5	31574	60914	7.52%	0.398%
23	12.325	1719	1723	1726	rBV2	9461	14435	1.78%	0.094%
24	12.386	1727	1733	1735	rVV4	15908	34673	4.28%	0.227%
25	12.416	1735	1738	1740	rVV2	36892	52801	6.52%	0.345%
26	12.441	1740	1742	1745	rVV	35831	40186	4.96%	0.263%
27	12.483	1745	1749	1753	rVV	189464	278176	34.33%	1.818%
28	12.526	1753	1756	1760	rVB2	36530	47513	5.86%	0.310%
29	12.569	1760	1763	1767	rVB4	8257	11393	1.41%	0.074%
30	12.648	1771	1776	1783	rVB6	29148	67658	8.35%	0.442%
31	12.733	1783	1790	1793	rBV3	62610	117793	14.54%	0.770%
32	12.770	1793	1796	1798	rVV	78052	111326	13.74%	0.727%
33	12.806	1798	1802	1808	rVV2	406905	625290	77.18%	4.086%
34	12.861	1808	1811	1816	rVB3	40581	60182	7.43%	0.393%
35	12.916	1816	1820	1821	rBV4	7552	9285	1.15%	0.061%
36	12.971	1821	1829	1834	rBV	120256	243157	30.01%	1.589%

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Max Peaks: 100

Stop Thrs : 0

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37	13.038	1837	1840	1847	rVB	98844	153927	19.00%	1.006%
38	13.123	1847	1854	1858	rBV2	39231	80721	9.96%	0.527%
39	13.166	1858	1861	1867	rBV3	19782	33846	4.18%	0.221%
40	13.227	1867	1871	1872	rBV3	14301	20721	2.56%	0.135%
41	13.251	1872	1875	1879	rVB3	28847	38744	4.78%	0.253%
42	13.318	1879	1886	1890	rBV3	121882	232222	28.66%	1.517%
43	13.367	1890	1894	1897	rVV3	70434	106941	13.20%	0.699%
44	13.428	1897	1904	1907	rVV2	278403	512850	63.30%	3.351%
45	13.459	1907	1909	1913	rVV	185396	211872	26.15%	1.385%
46	13.501	1913	1916	1920	rVB	54286	65745	8.11%	0.430%
47	13.623	1928	1936	1940	rBV	206731	338771	41.81%	2.214%
48	13.666	1940	1943	1951	rVB2	162326	278292	34.35%	1.819%
49	13.751	1951	1957	1962	rBV	481144	693913	85.65%	4.534%
50	13.800	1962	1965	1966	rVV2	29593	37829	4.67%	0.247%
51	13.824	1966	1969	1975	rVV	79927	130908	16.16%	0.855%
52	13.885	1976	1979	1981	rVV	39821	60245	7.44%	0.394%
53	13.916	1981	1984	1988	rVV	90659	132729	16.38%	0.867%
54	13.971	1989	1993	1998	rVV	168872	238066	29.38%	1.556%
55	14.020	1998	2001	2006	rVB2	45154	76858	9.49%	0.502%
56	14.081	2006	2011	2016	rVB	216048	328342	40.53%	2.146%
57	14.148	2016	2022	2027	rBV6	41950	112270	13.86%	0.734%
58	14.215	2027	2033	2035	rBV3	68269	127170	15.70%	0.831%
59	14.257	2035	2040	2044	rVV6	126694	303764	37.49%	1.985%
60	14.294	2044	2046	2049	rVV	96370	159408	19.68%	1.042%
61	14.337	2049	2053	2057	rVV	154189	234743	28.97%	1.534%
62	14.379	2057	2060	2064	rVV	114618	147661	18.23%	0.965%
63	14.422	2064	2067	2072	rVB2	98661	124035	15.31%	0.811%
64	14.477	2073	2076	2079	rBV4	32651	46625	5.75%	0.305%
65	14.519	2080	2083	2086	rVV	63273	81538	10.06%	0.533%
66	14.568	2086	2091	2094	rVV	136563	229743	28.36%	1.501%
67	14.611	2094	2098	2103	rVB	392771	514590	63.51%	3.363%
68	14.684	2103	2110	2115	rBV4	386328	810198	100.00%	5.294%
69	14.733	2115	2118	2124	rVB3	142780	237736	29.34%	1.554%
70	14.837	2129	2135	2143	rVB	269745	446237	55.08%	2.916%
71	14.904	2143	2146	2148	rBV3	28331	37614	4.64%	0.246%
72	14.946	2148	2153	2156	rVV3	113248	192066	23.71%	1.255%
73	14.977	2156	2158	2164	rVB	59366	74010	9.13%	0.484%
74	15.056	2164	2171	2176	rBV4	130519	273295	33.73%	1.786%
75	15.135	2182	2184	2190	rVB4	64280	101561	12.54%	0.664%
76	15.208	2191	2196	2205	rVV4	95853	217293	26.82%	1.420%

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77	15.294	2205	2210	2214	rVV	129564	213924	26.40%	1.398%
78	15.349	2214	2219	2229	rVV5	97758	355688	43.90%	2.324%
79	15.434	2229	2233	2241	rVB2	228239	361074	44.57%	2.359%
80	15.525	2242	2248	2255	rBV3	89034	218562	26.98%	1.428%
81	15.605	2257	2261	2267	rVB5	39073	68250	8.42%	0.446%
82	15.696	2267	2276	2277	rBV3	64398	137722	17.00%	0.900%
83	15.727	2277	2281	2289	rVB	219549	400019	49.37%	2.614%
84	15.818	2292	2296	2301	rVV3	17673	32201	3.97%	0.210%
85	15.885	2302	2307	2319	rVB4	53817	136449	16.84%	0.892%
86	16.038	2324	2332	2337	rBV3	128114	273749	33.79%	1.789%
87	16.153	2348	2351	2356	rBV3	22867	37942	4.68%	0.248%
88	16.233	2358	2364	2369	rVV3	107323	220654	27.23%	1.442%
89	16.294	2370	2374	2379	rVV3	45531	90645	11.19%	0.592%
90	16.361	2380	2385	2395	rVB	91349	171469	21.16%	1.120%
91	16.489	2400	2406	2407	rBV5	33268	59972	7.40%	0.392%
92	16.623	2423	2428	2437	rVB5	26144	71507	8.83%	0.467%

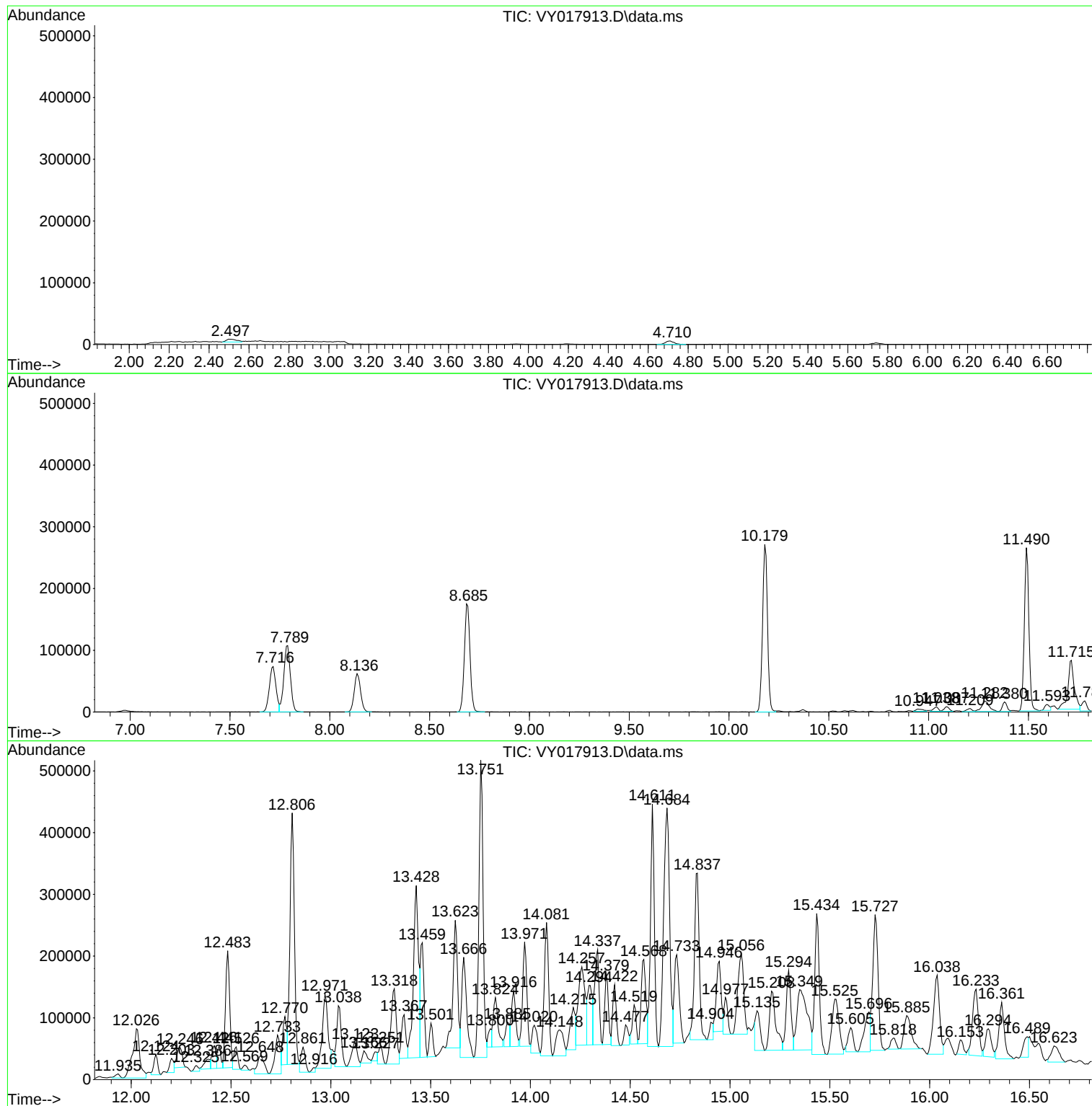
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Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041124\
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ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
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HR-02-041024

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032924S.M
Quant Title : SW846 8260

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



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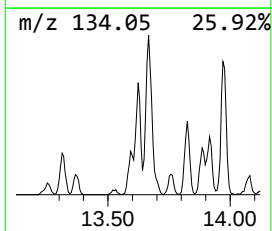
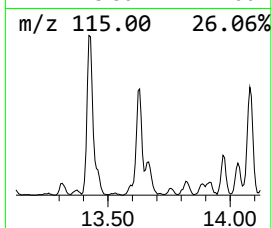
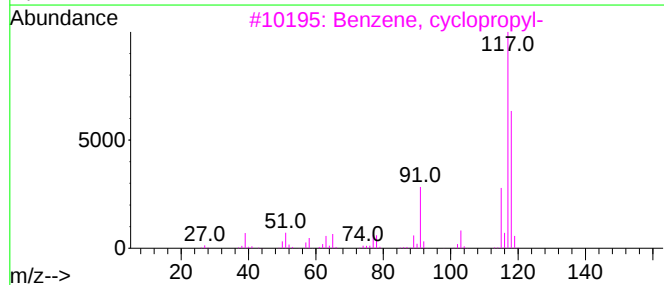
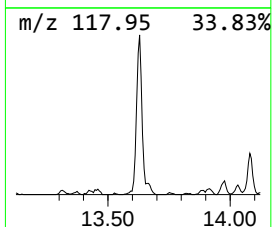
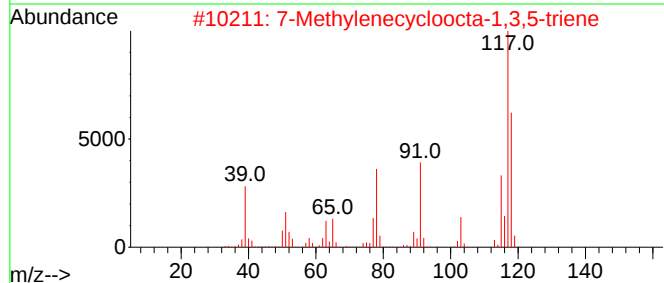
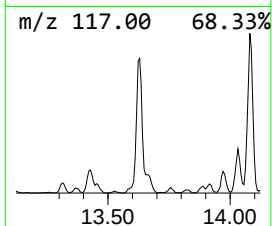
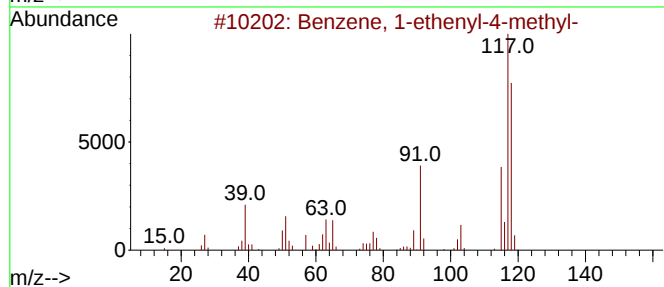
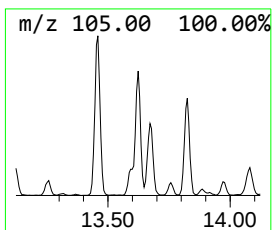
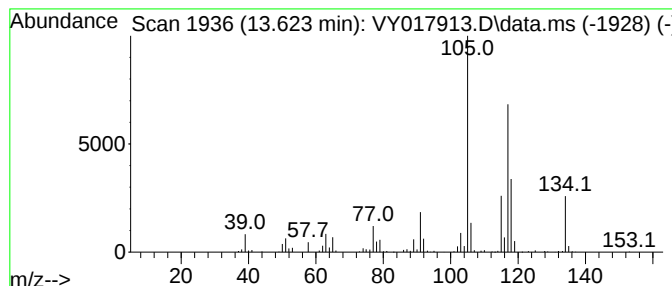
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032924S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethenyl-4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.623	33.03 ug/l	338771	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	50
2			7-Methylenecycloocta-1,3,5-triene	118	C9H10	002570-13-0	50
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	50
4			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	46
5			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	46



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ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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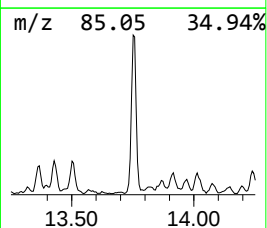
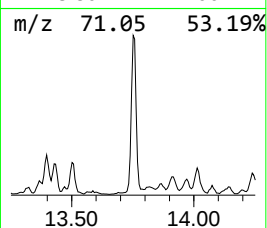
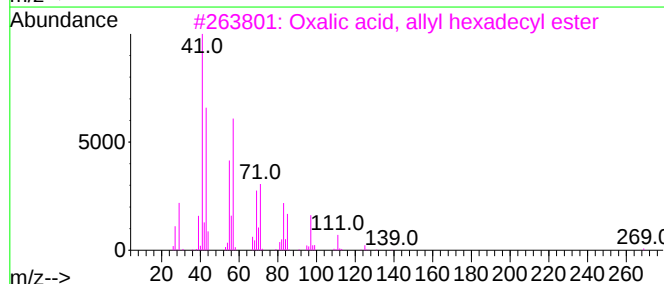
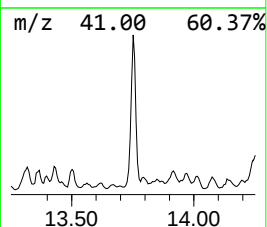
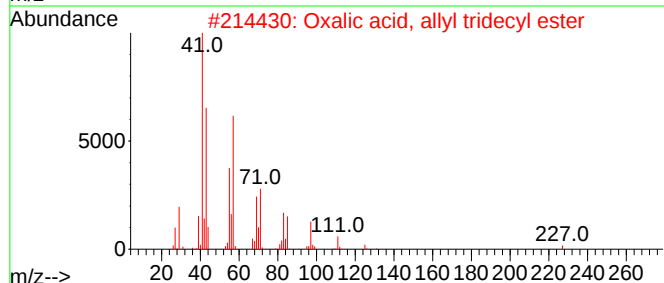
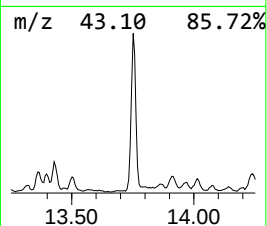
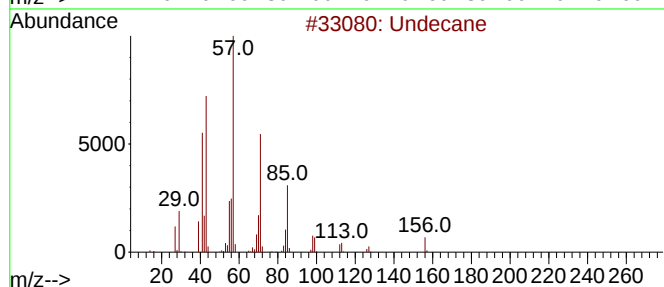
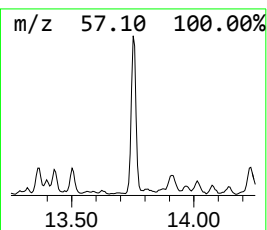
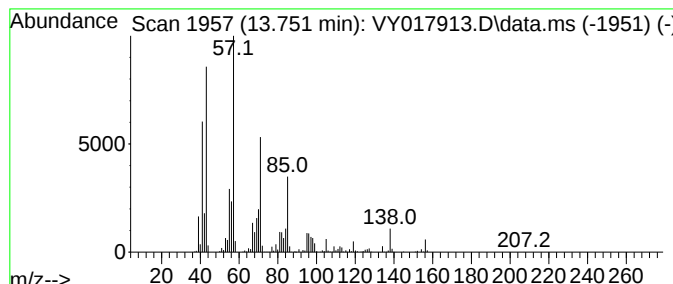
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032924S.M
Quant Title : SW846 8260

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

Peak Number 2 Undecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.751	67.65 ug/l	693913	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	83
2			Oxalic acid, allyl tridecyl ester	312	C18H32O4	1000309-24-1	72
3			Oxalic acid, allyl hexadecyl ester	354	C21H38O4	1000309-24-4	72
4			Decane	142	C10H22	000124-18-5	64
5			Carbonic acid, nonyl vinyl ester	214	C12H22O3	1000383-25-6	64



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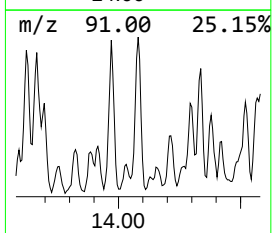
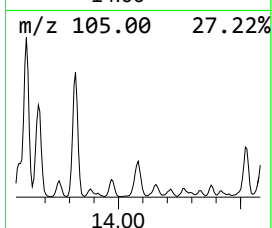
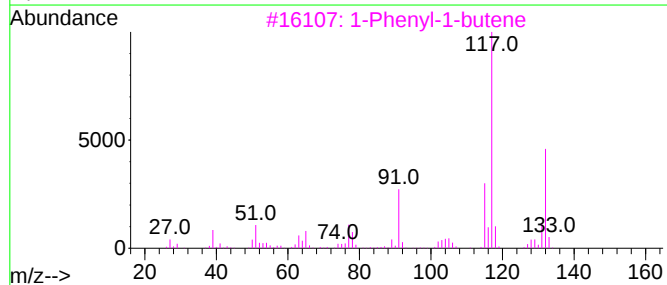
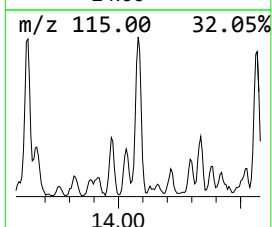
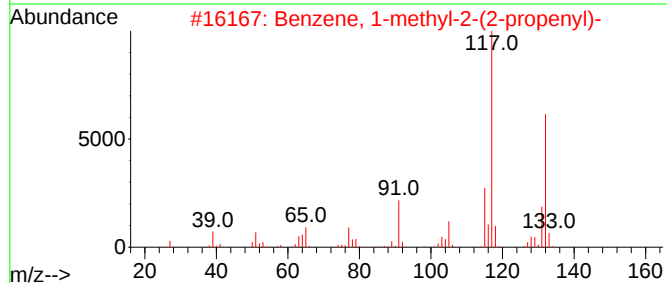
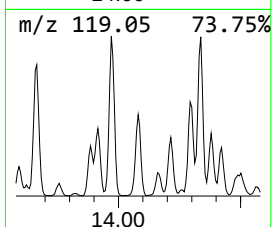
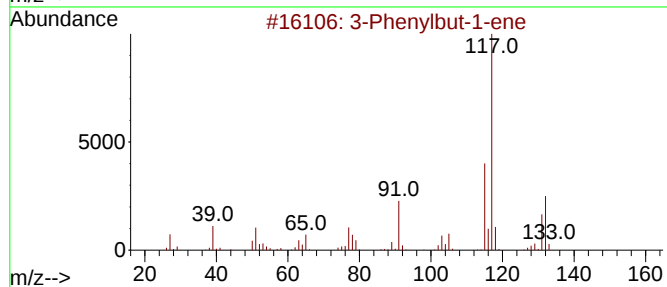
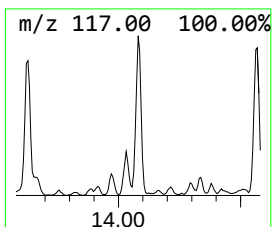
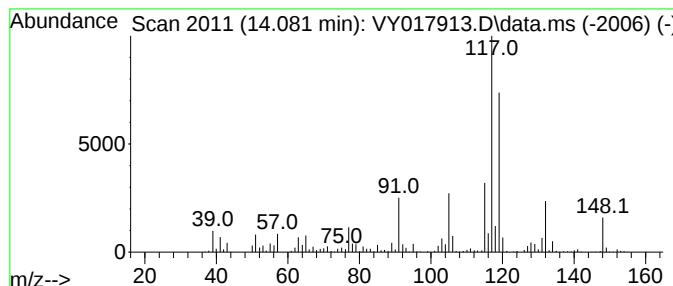
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032924S.M
Quant Title : SW846 8260

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

Peak Number 3 3-Phenylbut-1-ene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.081	32.01 ug/l	328342	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenylbut-1-ene	132	C10H12	000934-10-1	55
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	55
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	55
4			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	50
5			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041124\
Data File : VY017913.D
Acq On : 11 Apr 2024 16:01
Operator : SY/MD
Sample : P2072-03
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-02-041024

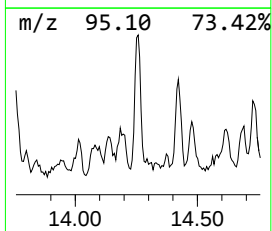
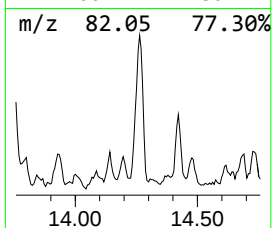
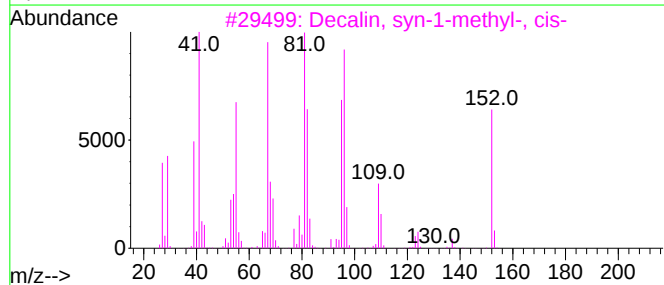
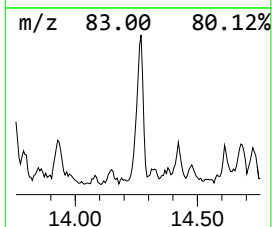
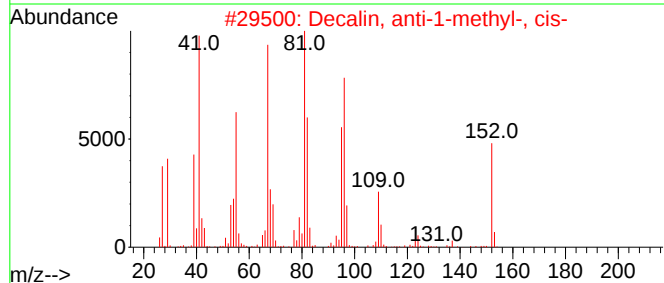
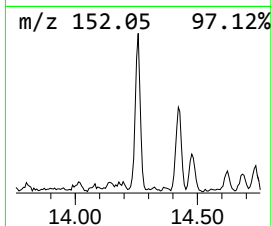
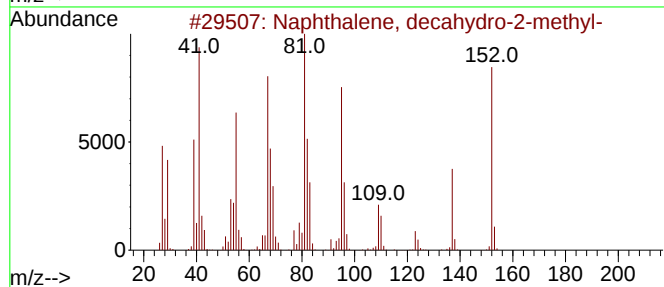
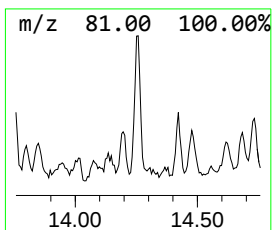
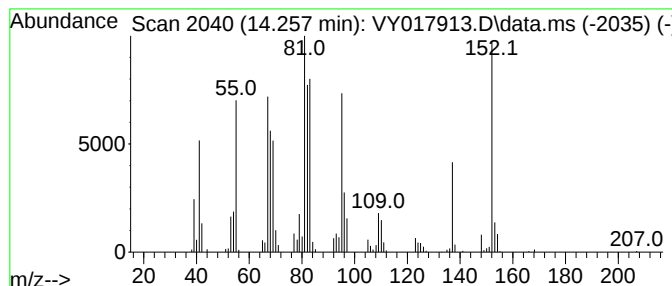
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Quant Title : SW846 8260

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

Peak Number 4 Naphthalene, decahydro-2-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.257	29.62 ug/l	303764	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	93
2			Decalin, anti-1-methyl-, cis-	152	C11H20	1000158-89-0	62
3			Decalin, syn-1-methyl-, cis-	152	C11H20	1000158-89-1	60
4			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	58
5			Cyclohexane, cyclopropyl-	124	C9H16	032669-86-6	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041124\
Data File : VY017913.D
Acq On : 11 Apr 2024 16:01
Operator : SY/MD
Sample : P2072-03
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-02-041024

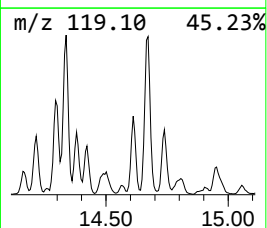
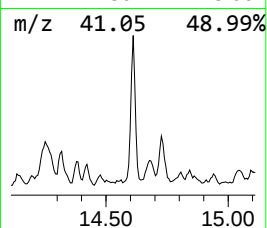
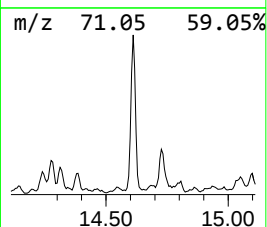
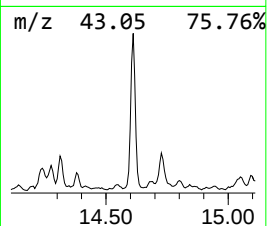
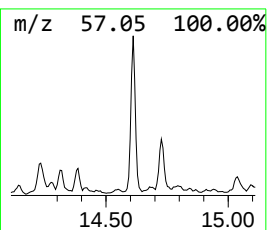
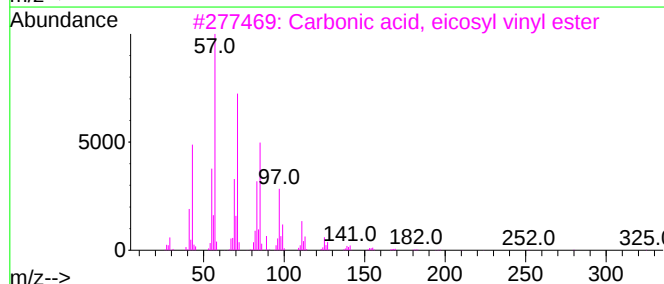
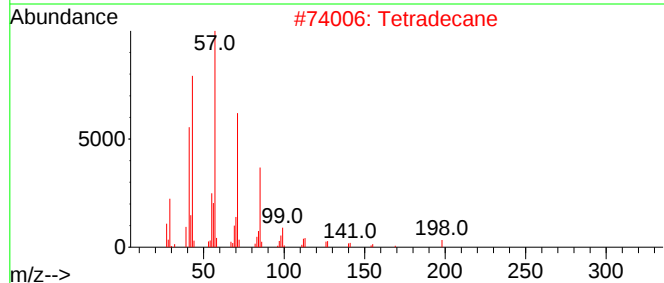
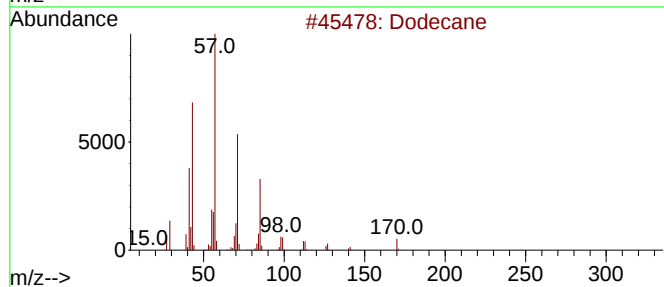
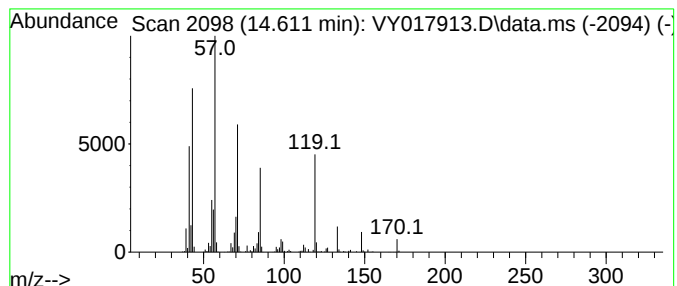
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Quant Title : SW846 8260

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

Peak Number 5 Dodecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.611	50.17 ug/l	514590	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	64
2			Tetradecane	198	C14H30	000629-59-4	50
3			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	50
4			Carbonic acid, dodecyl vinyl ester	256	C15H28O3	1000382-54-8	50
5			Tridecane, 2,5-dimethyl-	212	C15H32	056292-66-1	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041124\
Data File : VY017913.D
Acq On : 11 Apr 2024 16:01
Operator : SY/MD
Sample : P2072-03
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-02-041024

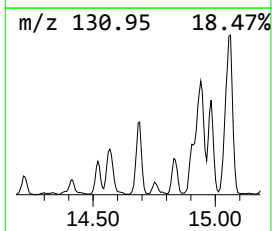
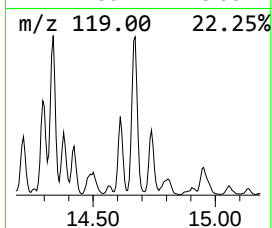
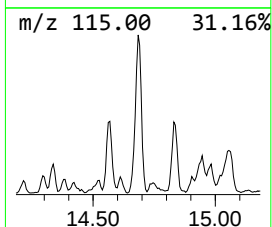
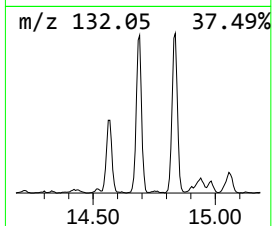
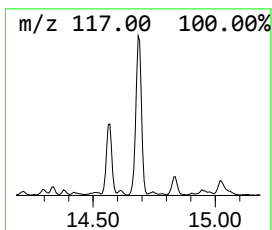
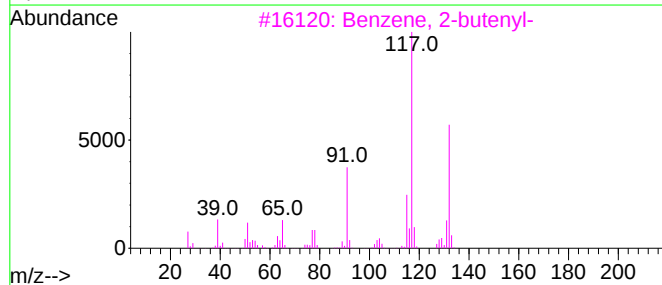
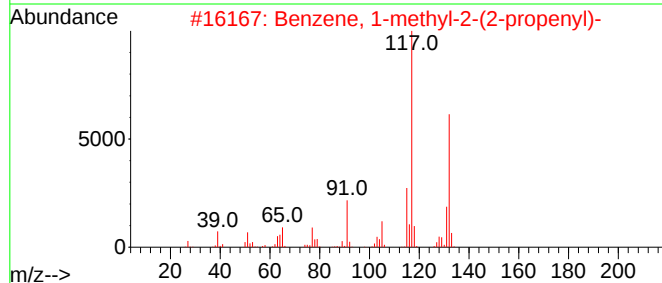
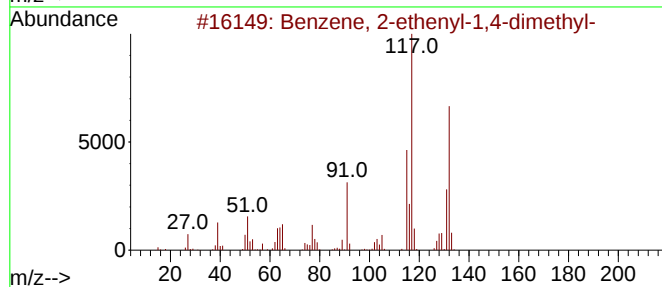
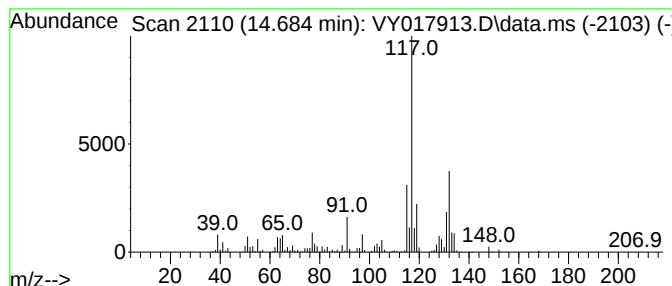
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	78.99 ug/l	810198	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
4			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	83
5			1-Phenyl-1-butene	132	C10H12	000824-90-8	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041124\
Data File : VY017913.D
Acq On : 11 Apr 2024 16:01
Operator : SY/MD
Sample : P2072-03
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-02-041024

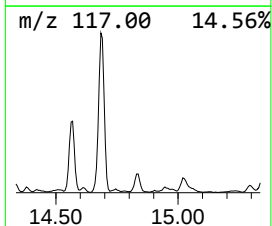
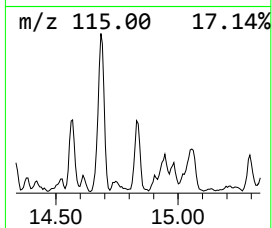
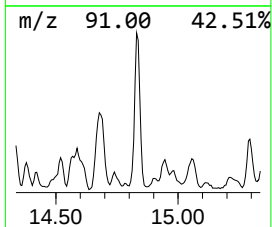
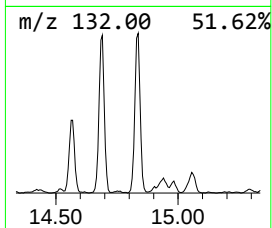
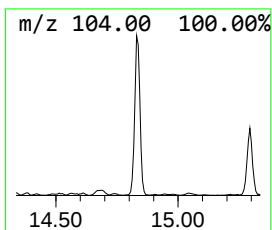
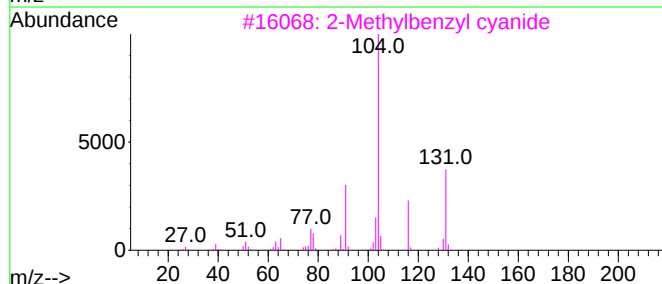
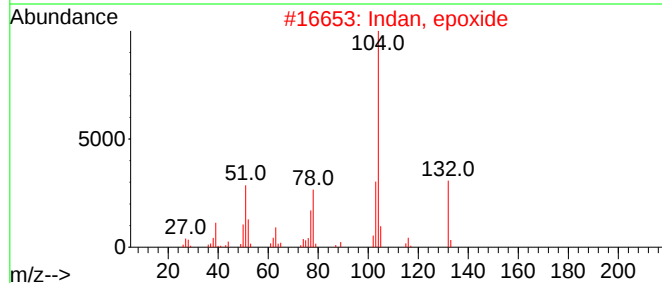
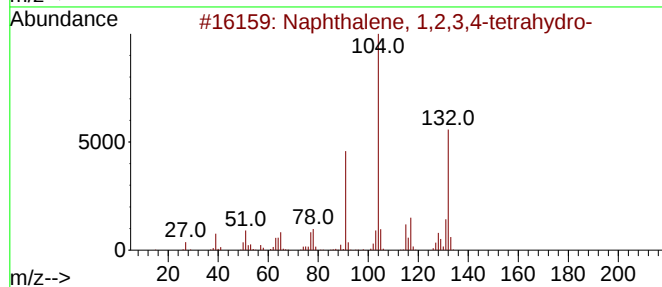
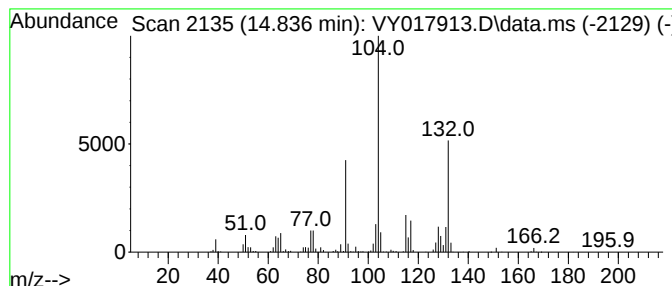
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.836	43.51 ug/l	446237	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2			Indan, epoxide	132	C9H8O	000768-22-9	58
3			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43
4			Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	007694-30-6	38
5			Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041124\
Data File : VY017913.D
Acq On : 11 Apr 2024 16:01
Operator : SY/MD
Sample : P2072-03
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-02-041024

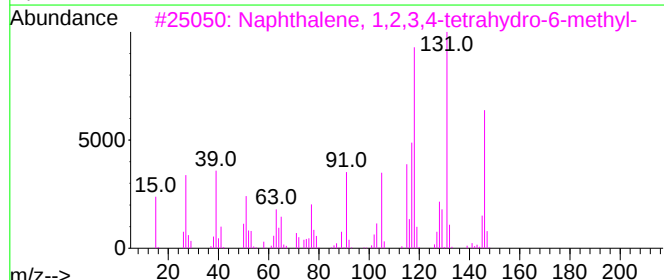
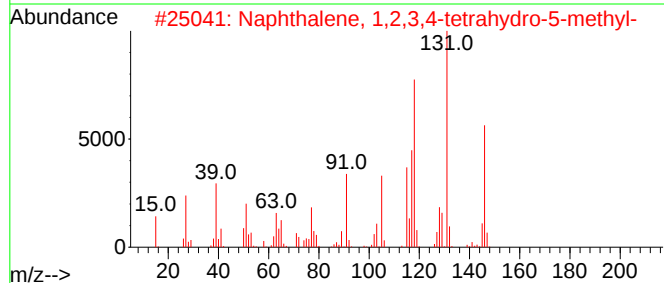
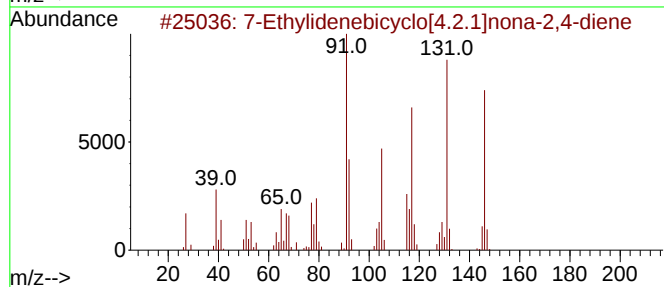
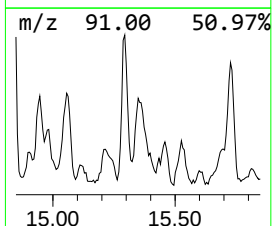
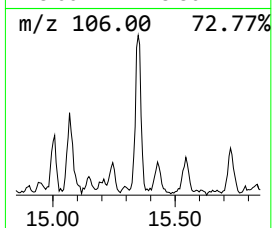
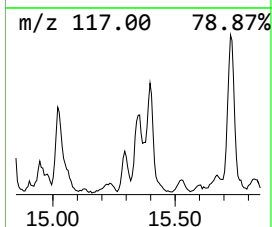
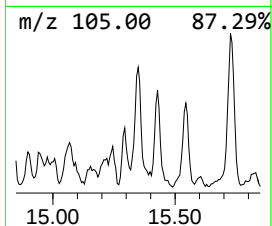
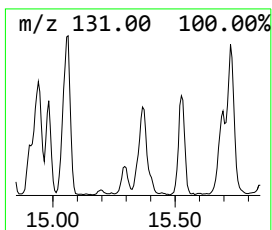
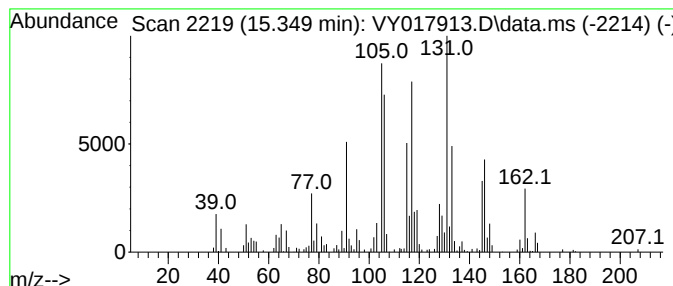
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Quant Title : SW846 8260

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

Peak Number 8 7-Ethylidenebicyclo[4.2.1]n... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.349	34.68 ug/l	355688	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Ethylidenebicyclo[4.2.1]nona-2...	146	C11H14	094400-10-9	55
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	55
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	50
4			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	42
5			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041124\
Data File : VY017913.D
Acq On : 11 Apr 2024 16:01
Operator : SY/MD
Sample : P2072-03
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-02-041024

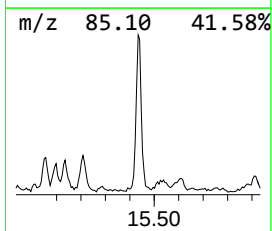
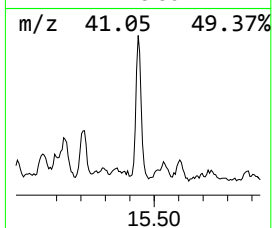
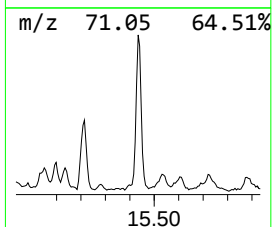
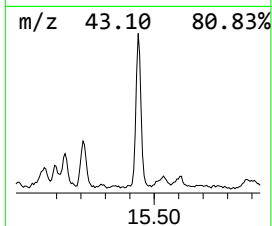
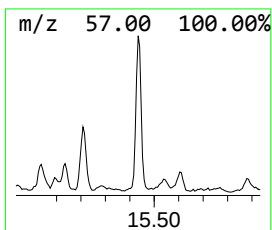
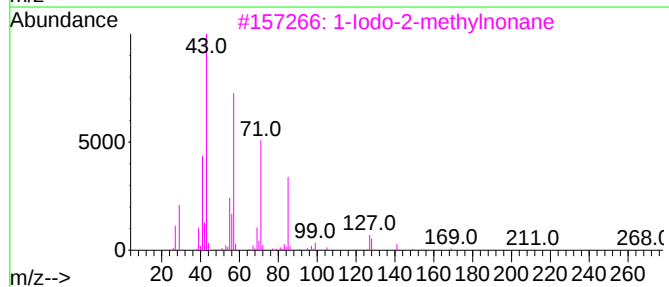
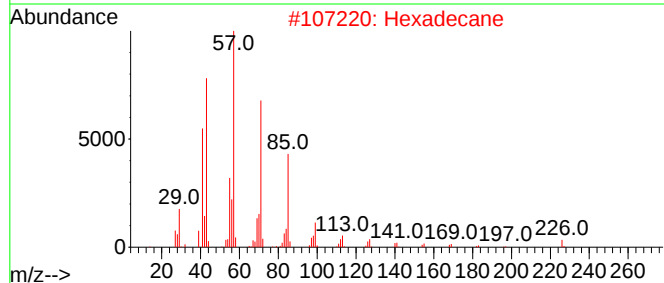
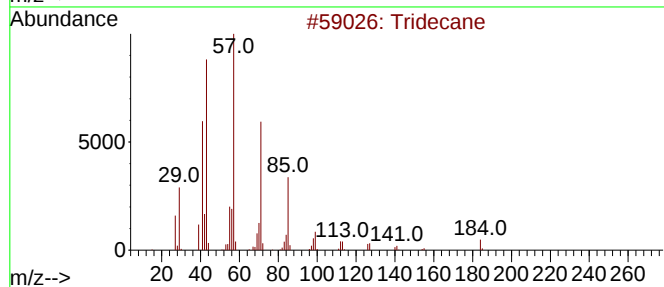
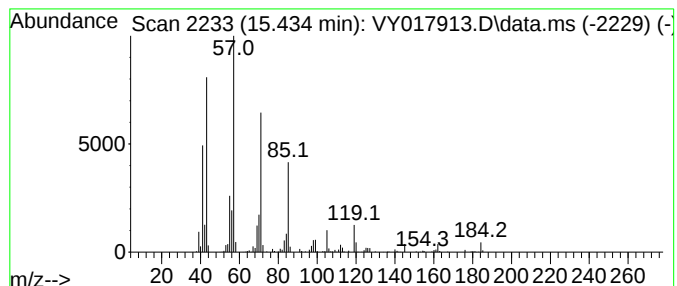
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Quant Title : SW846 8260

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

Peak Number 9 Tridecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.434	35.20 ug/l	361074	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	74
2			Hexadecane	226	C16H34	000544-76-3	72
3			1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	59
4			Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	50
5			Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041124\
Data File : VY017913.D
Acq On : 11 Apr 2024 16:01
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Sample : P2072-03
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ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HR-02-041024

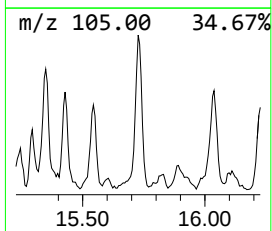
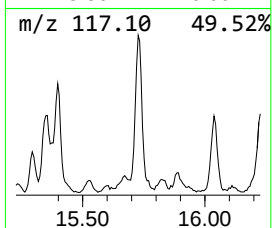
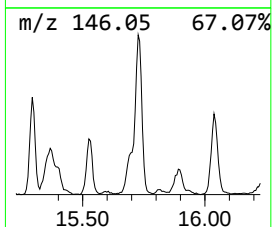
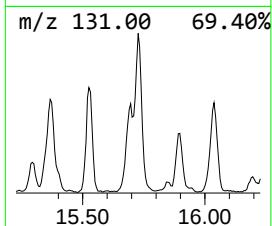
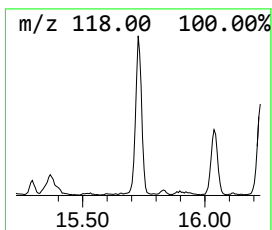
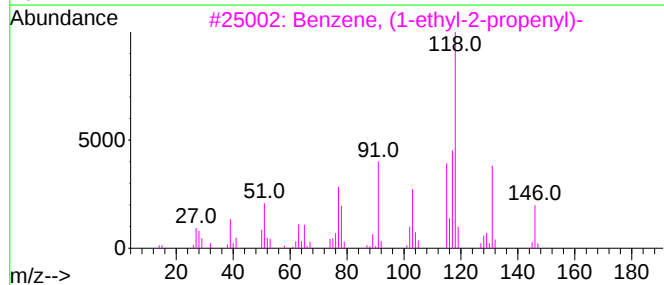
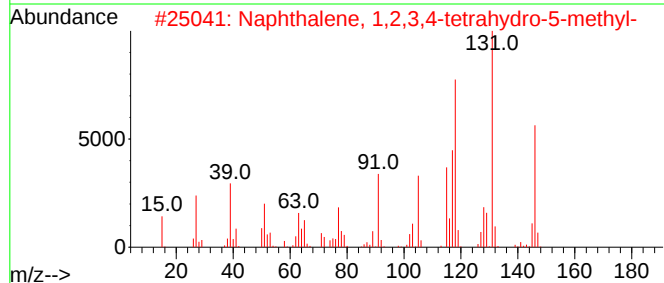
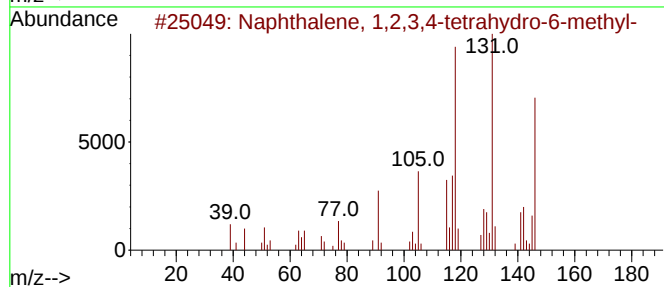
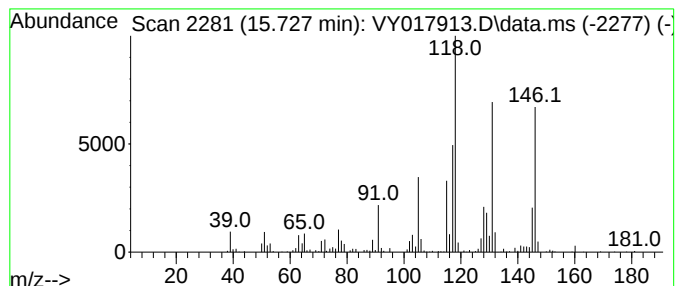
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032924S.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.727	39.00 ug/l	400019	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	83
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	74
3			Benzene, (1-ethyl-2-propenyl)-	146	C11H14	019947-22-9	74
4			Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	58
5			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041124\
Data File : VY017913.D
Acq On : 11 Apr 2024 16:01
Operator : SY/MD
Sample : P2072-03
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-02-041024

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032924S.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, 1-ethe...	13.623	33.0	ug/l	338771	4	13.428	512850	50.0
Undecane	13.751	67.7	ug/l	693913	4	13.428	512850	50.0
3-Phenylbut-1-ene	14.081	32.0	ug/l	328342	4	13.428	512850	50.0
Naphthalene, de...	14.257	29.6	ug/l	303764	4	13.428	512850	50.0
Dodecane	14.611	50.2	ug/l	514590	4	13.428	512850	50.0
Benzene, 2-ethe...	14.684	79.0	ug/l	810198	4	13.428	512850	50.0
Naphthalene, 1,...	14.836	43.5	ug/l	446237	4	13.428	512850	50.0
7-Ethylidenebic...	15.349	34.7	ug/l	355688	4	13.428	512850	50.0
Tridecane	15.434	35.2	ug/l	361074	4	13.428	512850	50.0
Naphthalene, 1,...	15.727	39.0	ug/l	400019	4	13.428	512850	50.0