

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110824\
 Data File : VY020225.D
 Acq On : 08 Nov 2024 12:05
 Operator : SY/MD
 Sample : P4734-01
 Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 EO-03-11062024

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103024S.M

Title : SW846 8260

Signal : TIC: VY020225.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.616	464	475	488	rBV2	12260	37479	2.65%	0.222%
2	6.884	838	847	858	rBV2	19012	64790	4.58%	0.383%
3	7.634	961	970	976	rBV	194632	463573	32.78%	2.742%
4	7.713	976	983	993	rVB	285449	653833	46.23%	3.867%
5	8.067	1029	1041	1052	rBV	189031	435324	30.78%	2.575%
6	8.616	1121	1131	1144	rBV	521841	1026539	72.58%	6.072%
7	10.103	1368	1375	1385	rBV	802877	1414379	100.00%	8.365%
8	10.506	1435	1441	1450	rBV2	18050	36578	2.59%	0.216%
9	10.877	1496	1502	1508	rBV5	11042	23163	1.64%	0.137%
10	11.012	1520	1524	1530	rBV5	8611	14275	1.01%	0.084%
11	11.213	1548	1557	1567	rVB9	14598	47970	3.39%	0.284%
12	11.304	1567	1572	1579	rBV2	10096	19173	1.36%	0.113%
13	11.414	1583	1590	1600	rBV	796645	1285975	90.92%	7.606%
14	11.639	1616	1627	1630	rBV6	32673	85923	6.07%	0.508%
15	11.707	1635	1638	1643	rVB3	15684	22019	1.56%	0.130%
16	11.926	1668	1674	1675	rBV4	33526	51104	3.61%	0.302%
17	12.048	1690	1694	1698	rVB	26181	34119	2.41%	0.202%
18	12.121	1702	1706	1710	rBV5	19035	36466	2.58%	0.216%
19	12.164	1710	1713	1719	rVB4	26991	46037	3.25%	0.272%
20	12.255	1723	1728	1733	rBV4	58278	111746	7.90%	0.661%
21	12.334	1739	1741	1744	rBV3	13321	14605	1.03%	0.086%
22	12.408	1747	1753	1758	rVV	546232	855511	60.49%	5.060%
23	12.566	1776	1779	1782	rBV3	15914	22929	1.62%	0.136%
24	12.658	1787	1794	1797	rBV2	65010	114860	8.12%	0.679%
25	12.694	1797	1800	1801	rVV2	28769	34125	2.41%	0.202%
26	12.731	1801	1806	1811	rVV2	235759	378670	26.77%	2.240%
27	12.780	1811	1814	1819	rVB2	33207	48779	3.45%	0.289%
28	12.834	1819	1823	1826	rBV5	11555	20719	1.46%	0.123%
29	12.889	1826	1832	1837	rBV2	49897	116092	8.21%	0.687%
30	12.962	1840	1844	1850	rVB2	84225	126291	8.93%	0.747%
31	13.042	1850	1857	1862	rBV	102889	171826	12.15%	1.016%
32	13.090	1862	1865	1870	rBV4	20144	29523	2.09%	0.175%
33	13.170	1871	1878	1882	rVB4	32002	72839	5.15%	0.431%
34	13.237	1882	1889	1893	rBV4	101922	197473	13.96%	1.168%
35	13.286	1893	1897	1900	rVV4	62172	95244	6.73%	0.563%
36	13.346	1900	1907	1916	rVV2	690342	1256587	88.84%	7.432%

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Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103024S.M
 Title : SW846 8260

37	13.426	1916	1920	1923	rVB3	37686	47803	3.38%	0.283%
38	13.542	1932	1939	1943	rBV	142343	237434	16.79%	1.404%
39	13.584	1943	1946	1955	rVB3	119725	231956	16.40%	1.372%
40	13.676	1955	1961	1965	rBV2	296104	474689	33.56%	2.808%
41	13.712	1965	1967	1969	rBV	22866	27789	1.96%	0.164%
42	13.743	1969	1972	1978	rVB	51584	77001	5.44%	0.455%
43	13.804	1979	1982	1984	rBV	32272	44170	3.12%	0.261%
44	13.834	1984	1987	1992	rVB4	77359	108539	7.67%	0.642%
45	13.889	1992	1996	2001	rVB2	158552	231207	16.35%	1.367%
46	13.938	2001	2004	2009	rVB3	43904	72183	5.10%	0.427%
47	13.999	2009	2014	2019	rVB	172610	261558	18.49%	1.547%
48	14.060	2019	2024	2030	rBV6	39527	100522	7.11%	0.595%
49	14.133	2030	2036	2038	rBV4	52720	102397	7.24%	0.606%
50	14.176	2038	2043	2047	rBV7	82648	144775	10.24%	0.856%
51	14.255	2052	2056	2060	rVV	141468	215615	15.24%	1.275%
52	14.298	2060	2063	2067	rVV	90547	131442	9.29%	0.777%
53	14.340	2067	2070	2075	rVV	93802	123400	8.72%	0.730%
54	14.401	2076	2080	2082	rVV4	31649	49563	3.50%	0.293%
55	14.438	2083	2086	2089	rVV	60482	80367	5.68%	0.475%
56	14.480	2089	2093	2098	rVV2	110618	206712	14.62%	1.223%
57	14.529	2098	2101	2106	rVB	193272	260881	18.44%	1.543%
58	14.602	2106	2113	2118	rBV4	337946	707052	49.99%	4.182%
59	14.651	2118	2121	2127	rVB3	114346	188641	13.34%	1.116%
60	14.749	2127	2137	2145	rBV	215440	394296	27.88%	2.332%
61	14.822	2146	2149	2150	rVV2	26716	29394	2.08%	0.174%
62	14.858	2150	2155	2158	rVV2	112637	202328	14.31%	1.197%
63	14.895	2158	2161	2167	rVV4	68331	118146	8.35%	0.699%
64	14.968	2167	2173	2179	rVB2	113936	239777	16.95%	1.418%
65	15.121	2192	2198	2201	rBV5	50727	98554	6.97%	0.583%
66	15.200	2207	2211	2216	rBV	106259	168686	11.93%	0.998%
67	15.255	2216	2220	2231	rVV4	79195	274540	19.41%	1.624%
68	15.340	2231	2234	2242	rVB4	71064	129507	9.16%	0.766%
69	15.431	2242	2249	2255	rBV3	81618	177067	12.52%	1.047%
70	15.505	2257	2261	2266	rVB6	20212	34901	2.47%	0.206%
71	15.596	2267	2276	2277	rBV3	58851	118007	8.34%	0.698%
72	15.627	2277	2281	2290	rVB	196944	344555	24.36%	2.038%
73	15.712	2291	2295	2301	rBV6	21131	37069	2.62%	0.219%
74	15.785	2301	2307	2322	rVB2	56471	178982	12.65%	1.059%
75	15.931	2324	2331	2338	rBV2	117139	237851	16.82%	1.407%
76	16.120	2353	2362	2368	rVV3	92825	211718	14.97%	1.252%

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ClientSampleId :
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Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

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Title : SW846 8260

77	16.181	2368	2372	2386	rVB5	43891	117961	8.34%	0.698%
78	16.376	2398	2404	2410	rBV7	33763	96279	6.81%	0.569%
79	16.431	2410	2413	2418	rVV4	23120	38814	2.74%	0.230%
80	16.511	2420	2426	2433	rVB4	26561	66696	4.72%	0.394%

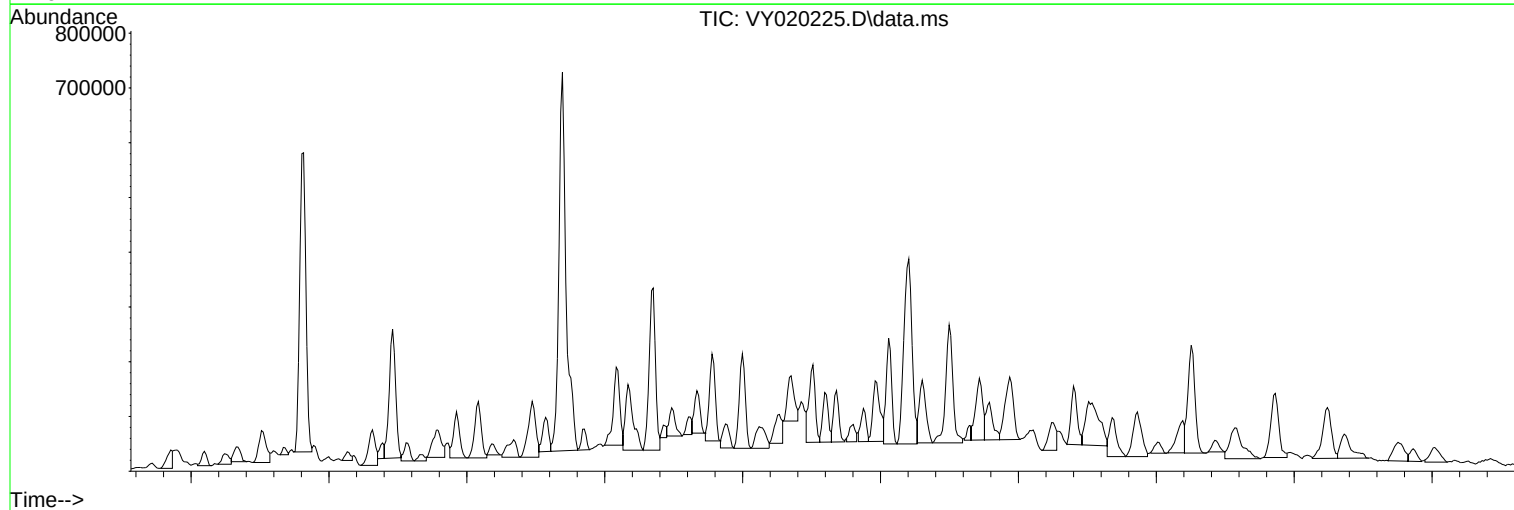
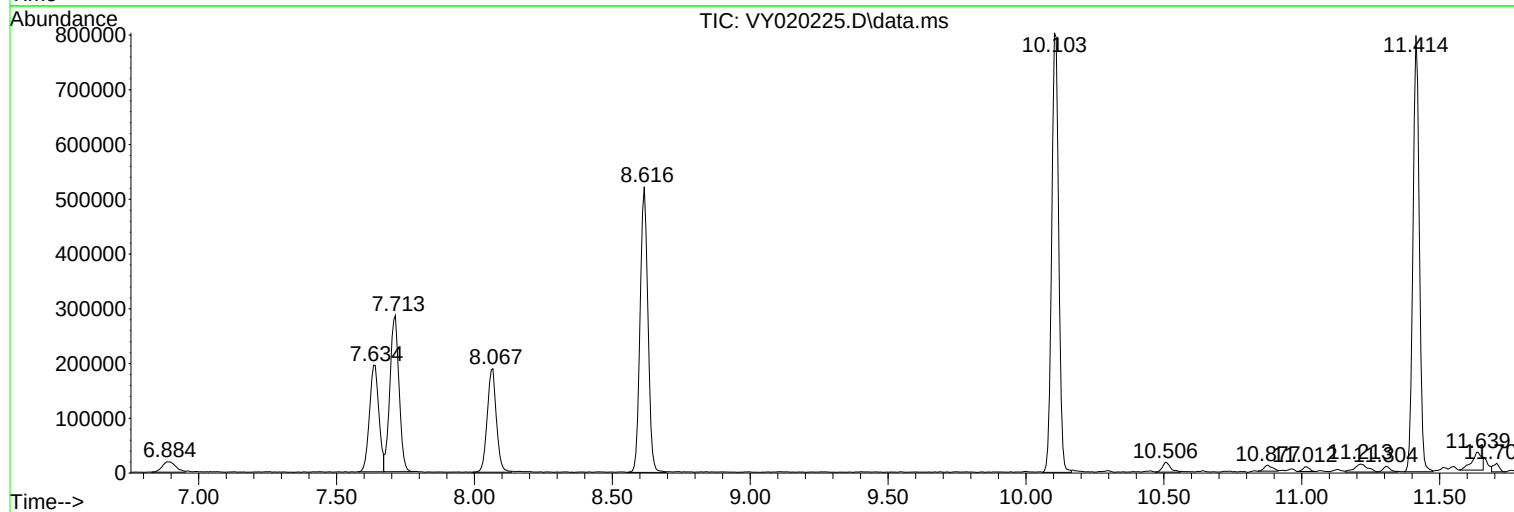
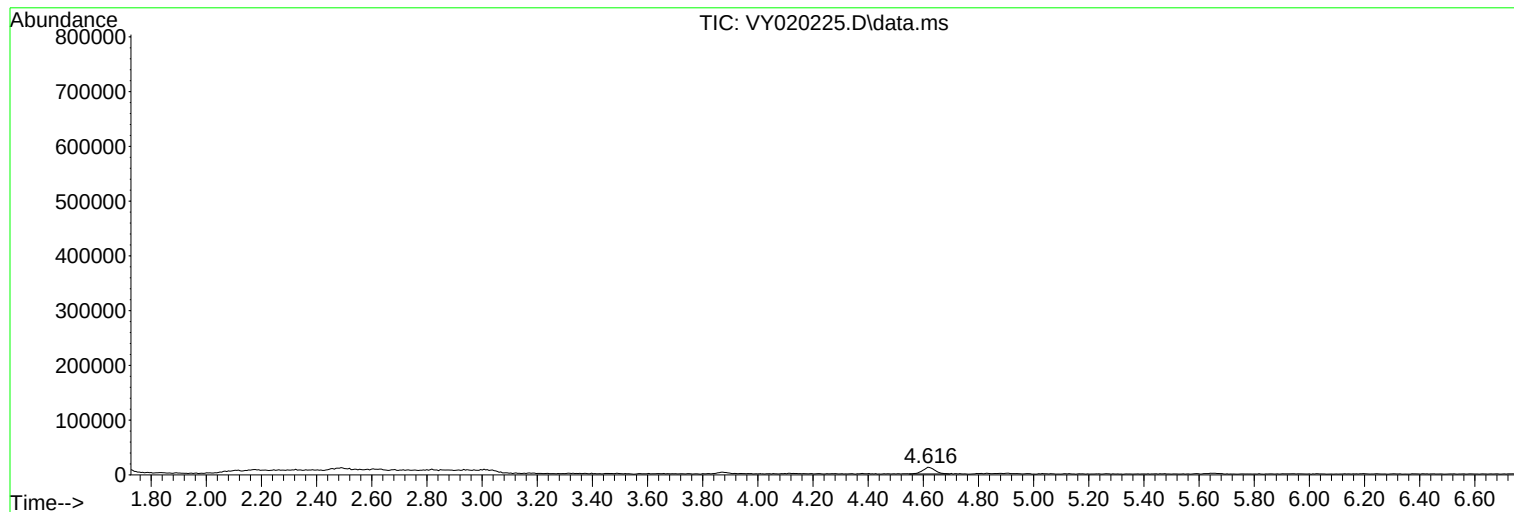
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103024S.M
Quant Title : SW846 8260

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



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ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-03-11062024

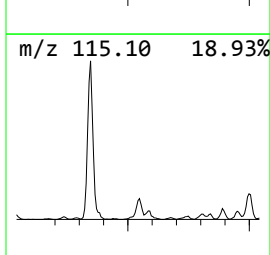
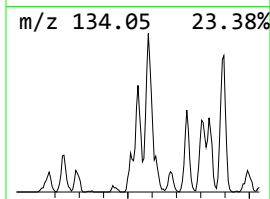
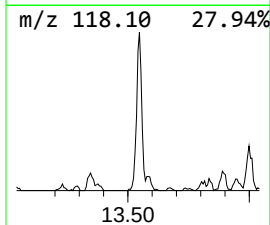
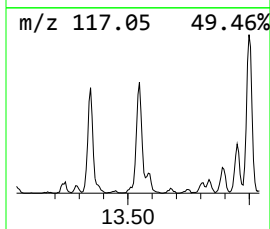
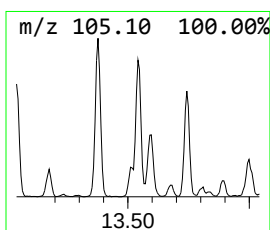
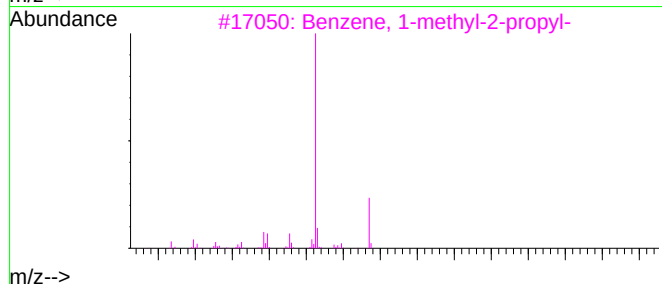
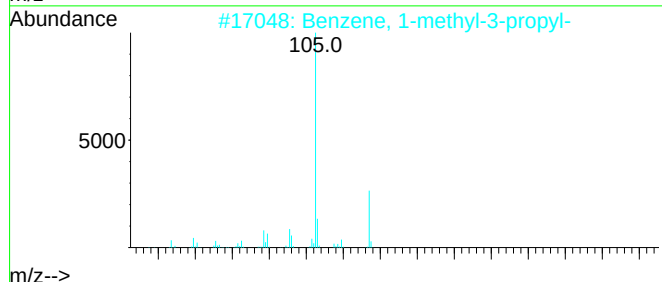
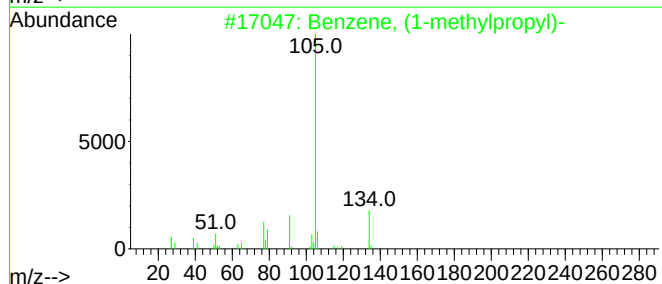
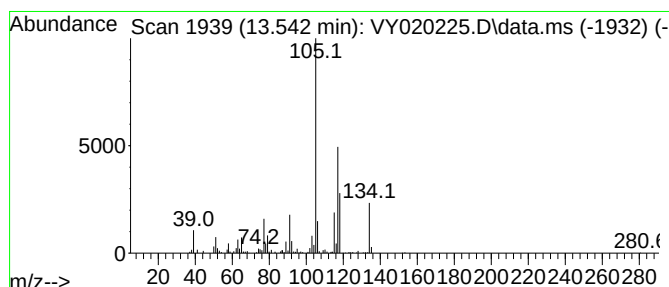
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103024S.M
Quant Title : SW846 8260

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

Peak Number 1 unknown13.542 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.542	9.45 ug/l	237434	1,4-Dichlorobenzene-d4	13.347

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	46
2	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	43
3	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	43
4	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	41
5	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	30



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Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-03-11062024

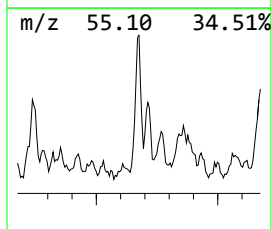
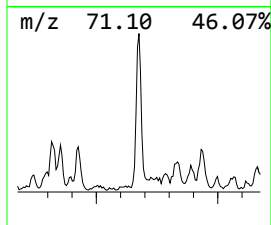
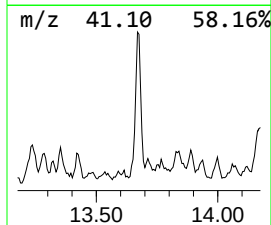
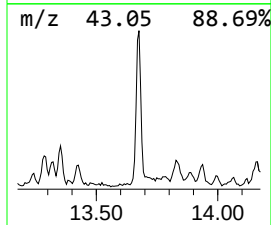
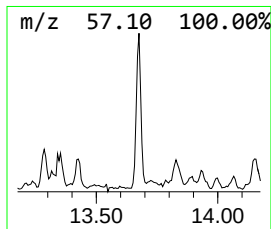
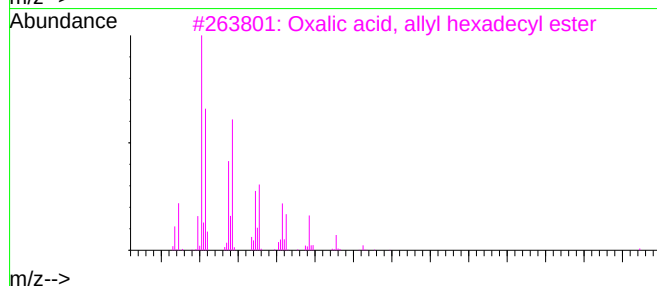
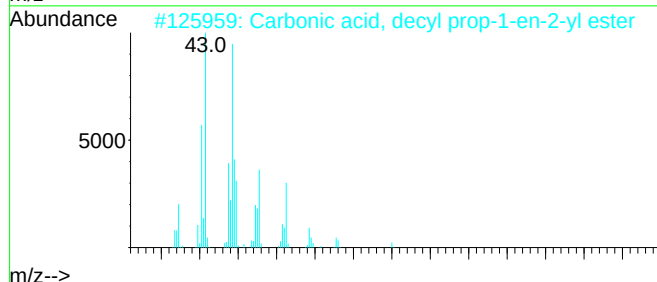
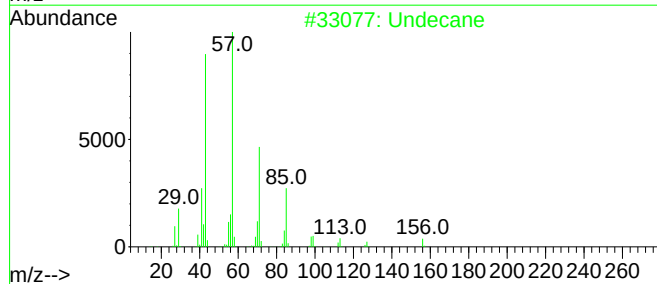
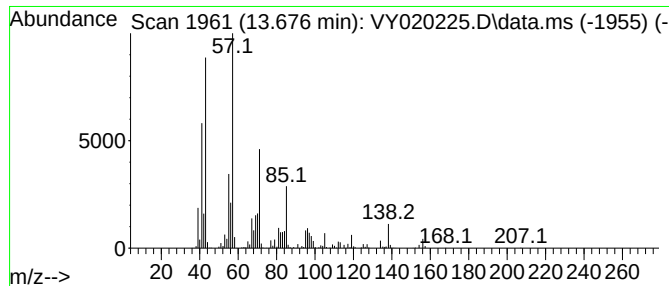
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103024S.M
Quant Title : SW846 8260

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.
13.676	18.89 ug/l	474689	1,4-Dichlorobenzene-d4	13.347

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane	156	C11H24	001120-21-4	68
2	Carbonic acid, decyl prop-1-en-2...	242	C14H26O3	1000382-90-5	64
3	Oxalic acid, allyl hexadecyl ester	354	C21H38O4	1000309-24-4	59
4	Tridecane	184	C13H28	000629-50-5	52
5	Carbonic acid, nonyl vinyl ester	214	C12H22O3	1000383-25-6	52



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ALS Vial : 6 Sample Multiplier: 1

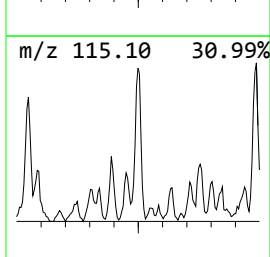
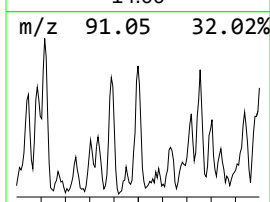
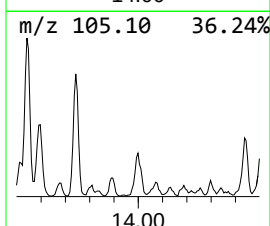
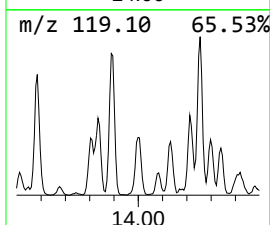
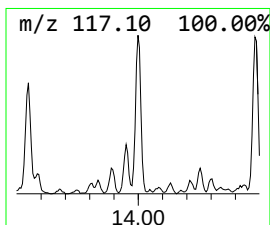
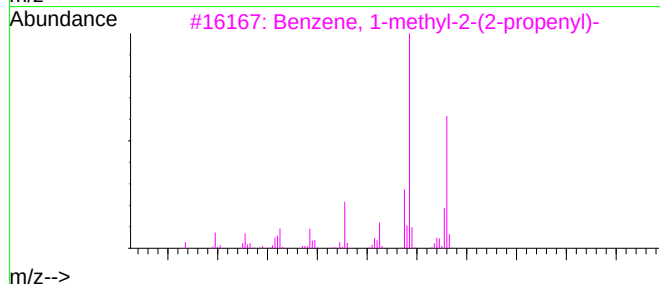
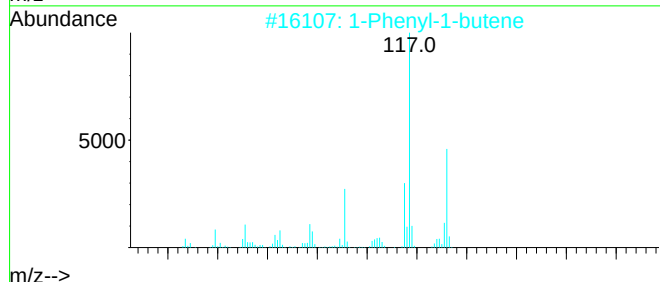
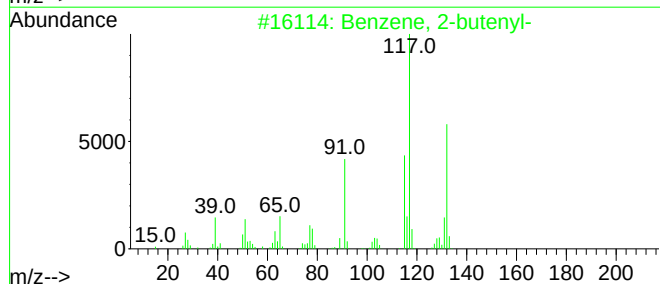
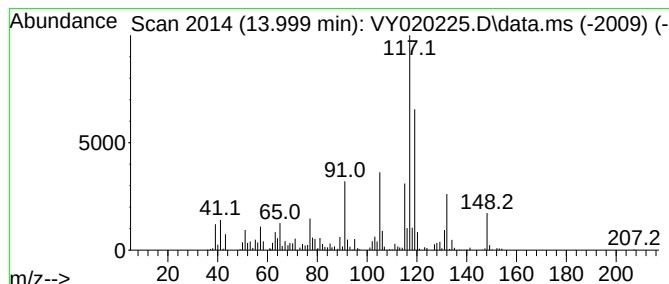
Instrument :
MSVOA_Y
ClientSampleId :
EO-03-11062024

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103024S.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 2-butenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.999	10.41 ug/l	261558	1,4-Dichlorobenzene-d4		13.347	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 2-butenyl-		132	C10H12	001560-06-1	92
2	1-Phenyl-1-butene		132	C10H12	000824-90-8	83
3	Benzene, 1-methyl-2-(2-propenyl)-		132	C10H12	001587-04-8	70
4	Benzene, 1-ethenyl-3-ethyl-		132	C10H12	007525-62-4	64
5	Benzene, (2-methyl-1-propenyl)-		132	C10H12	000768-49-0	64



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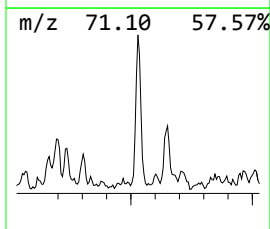
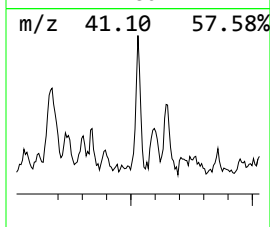
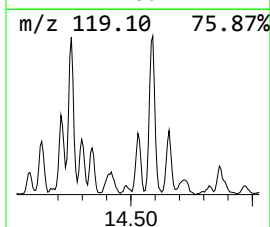
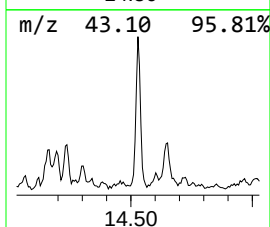
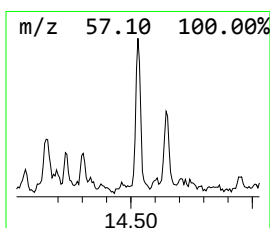
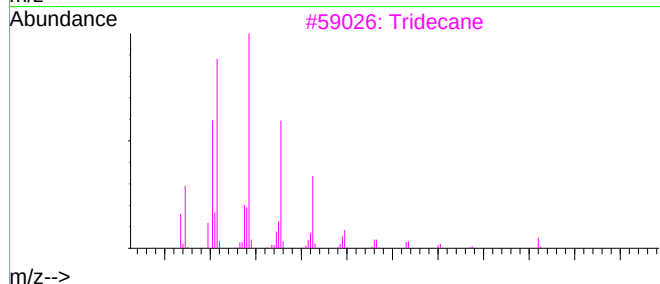
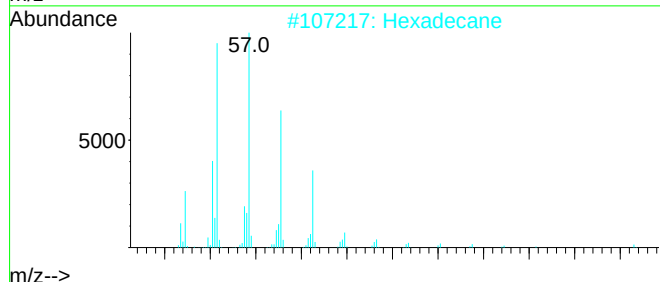
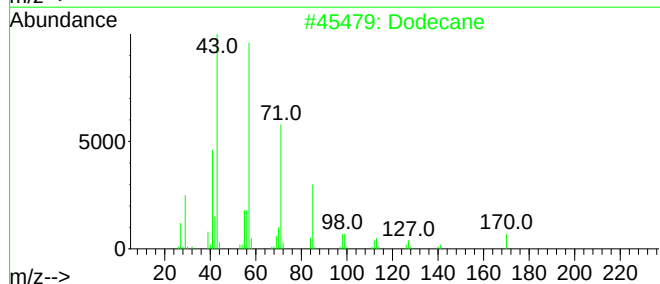
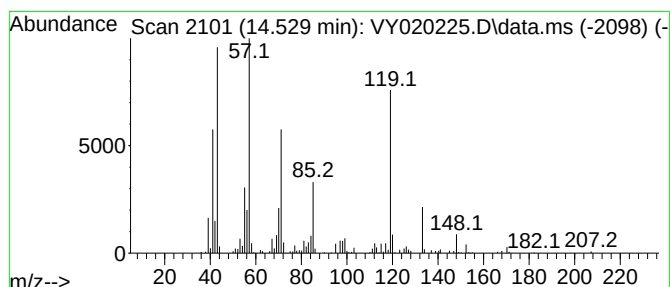
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Quant Title : SW846 8260

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

Peak Number 4 unknown14.529 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.529	10.38 ug/l	260881	1,4-Dichlorobenzene-d4	13.347

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane	170	C12H26	000112-40-3	46
2	Hexadecane	226	C16H34	000544-76-3	38
3	Tridecane	184	C13H28	000629-50-5	35
4	2,6-Dimethyldecane	170	C12H26	013150-81-7	30
5	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	27



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110824\
Data File : VY020225.D
Acq On : 08 Nov 2024 12:05
Operator : SY/MD
Sample : P4734-01
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-03-11062024

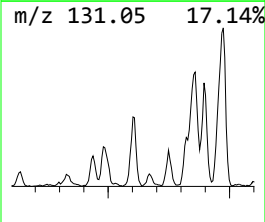
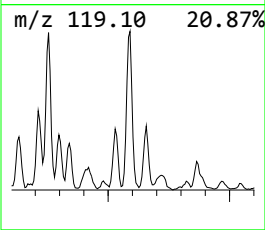
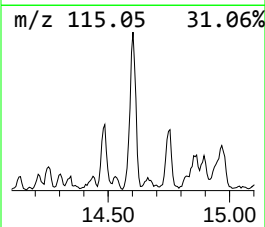
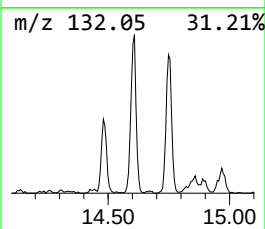
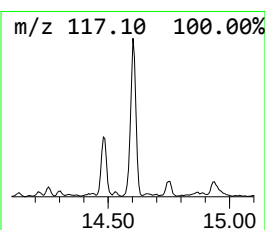
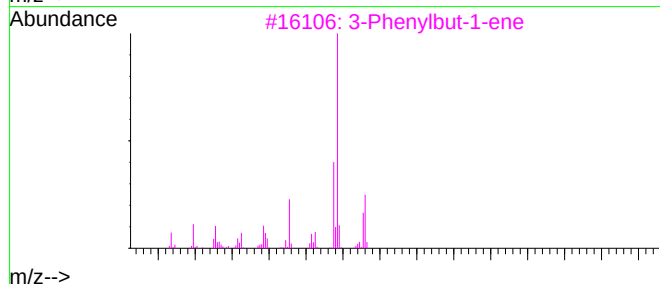
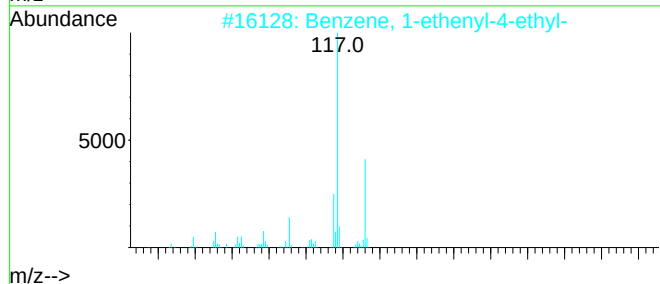
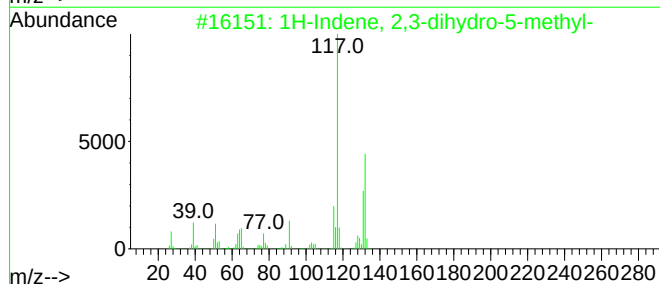
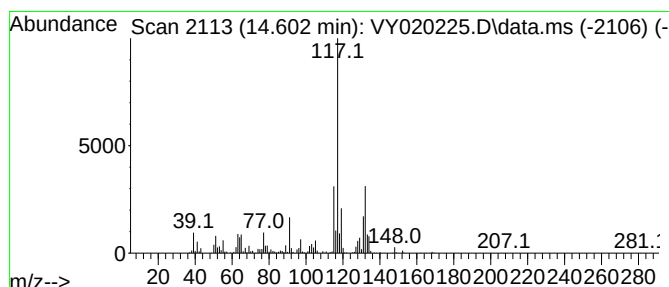
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.602	28.13 ug/l	707052	1,4-Dichlorobenzene-d4	13.347

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76
5	Indan, 1-methyl-	132	C10H12	000767-58-8	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110824\
Data File : VY020225.D
Acq On : 08 Nov 2024 12:05
Operator : SY/MD
Sample : P4734-01
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-03-11062024

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103024S.M
Quant Title : SW846 8260

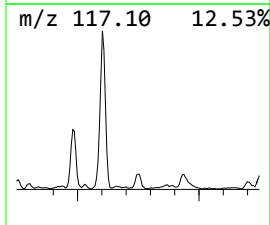
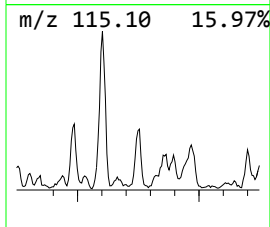
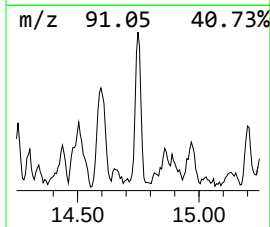
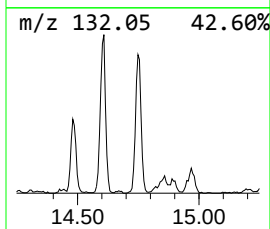
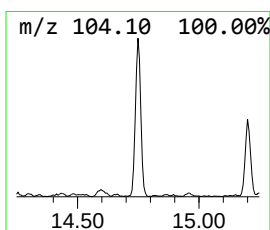
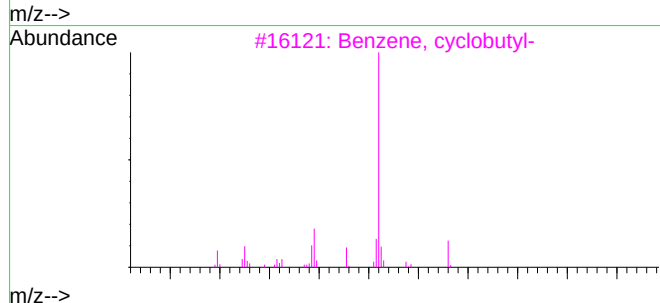
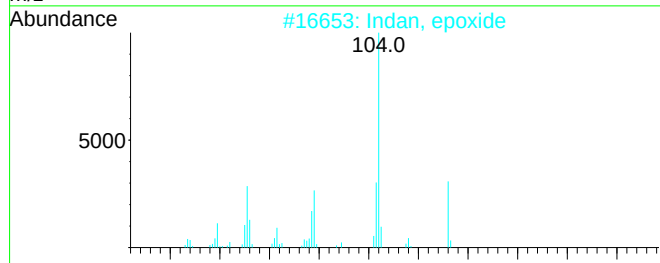
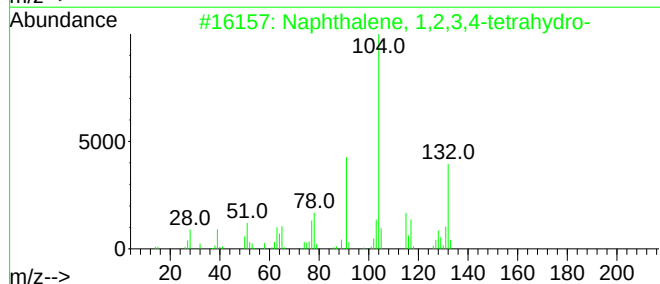
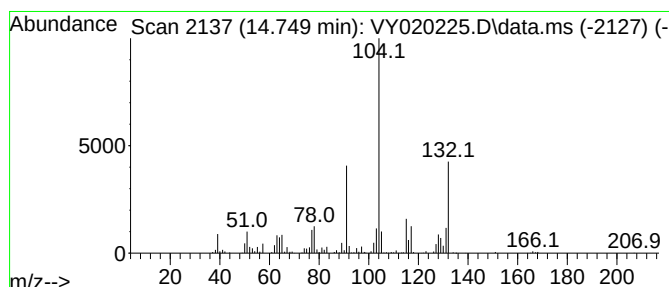
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 6 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.749	15.69 ug/l	394296	1,4-Dichlorobenzene-d4	13.347

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	62
3	Benzene, cyclobutyl-	132	C10H12	004392-30-7	59
4	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	53
5	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110824\
Data File : VY020225.D
Acq On : 08 Nov 2024 12:05
Operator : SY/MD
Sample : P4734-01
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-03-11062024

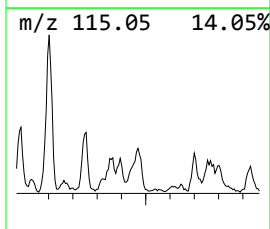
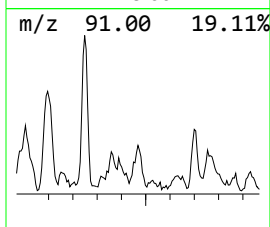
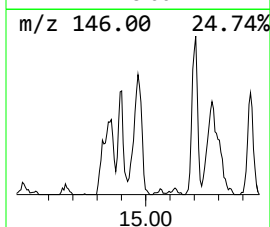
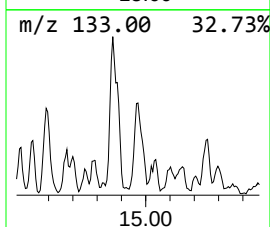
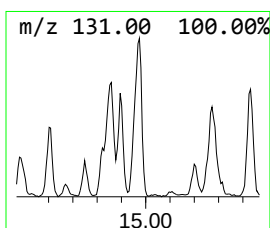
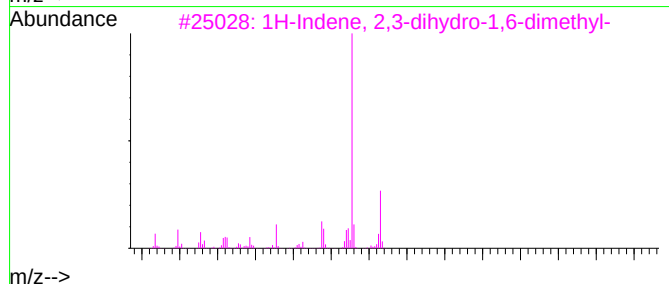
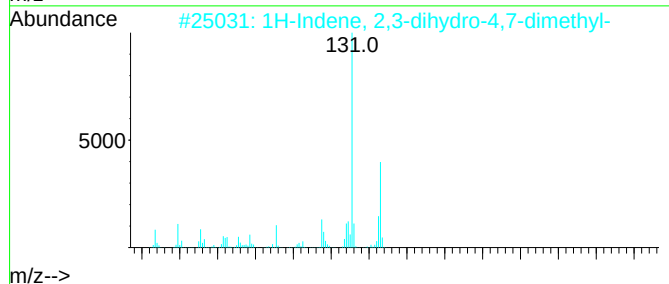
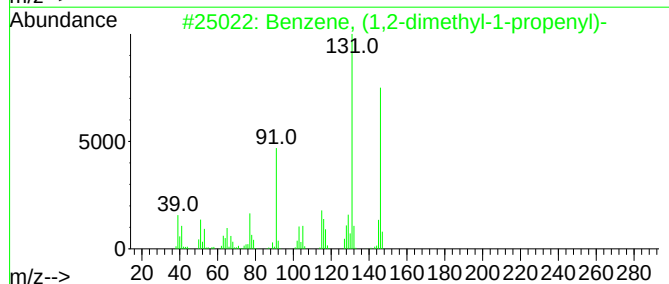
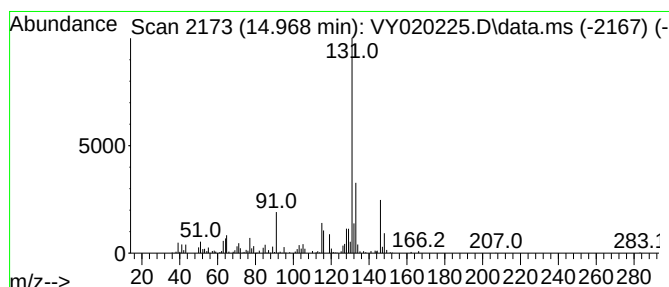
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.968	9.54 ug/l	239777	1,4-Dichlorobenzene-d4	13.347	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	94
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
3	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	93
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90
5	1,3-Cyclopentadiene, 5-(trans-2-...	146	C11H14	079209-36-2	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110824\
Data File : VY020225.D
Acq On : 08 Nov 2024 12:05
Operator : SY/MD
Sample : P4734-01
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-03-11062024

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103024S.M
Quant Title : SW846 8260

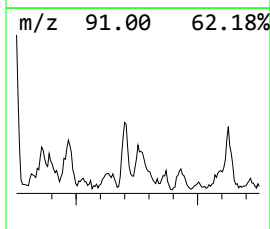
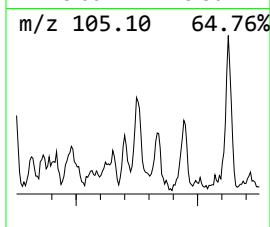
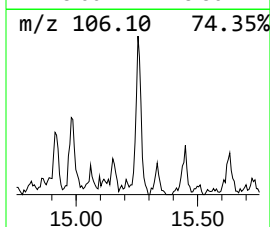
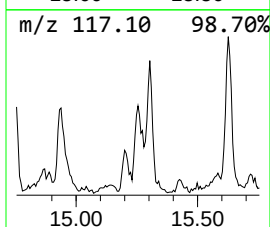
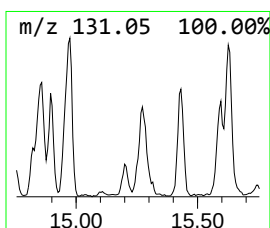
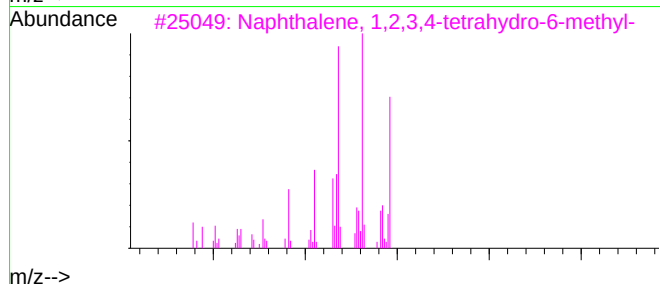
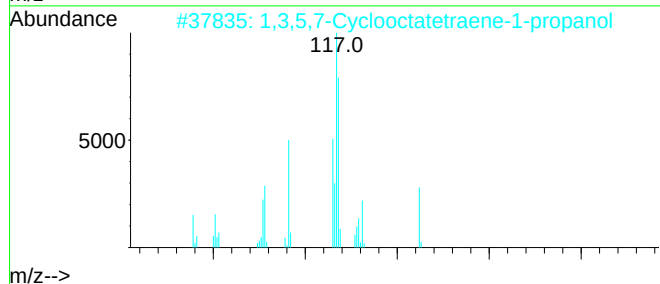
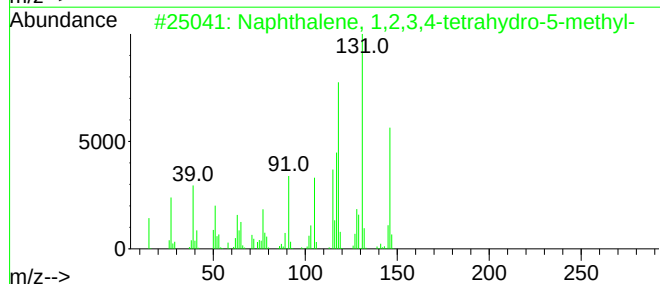
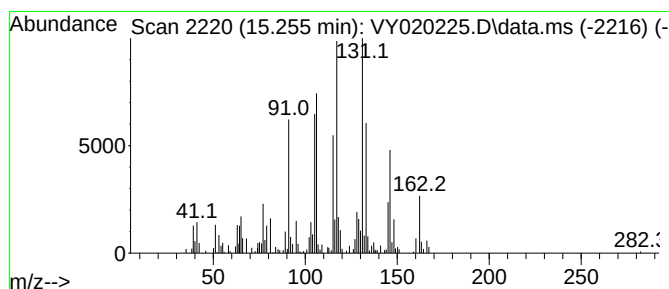
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 8 Naphthalene, 1,2,3,4-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.255	10.92 ug/l	274540	1,4-Dichlorobenzene-d4	13.347

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	50
2	1,3,5,7-Cyclooctatetraene-1-prop...	162	C11H14O	054365-73-0	42
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	41
4	1H-Indene-4-carboxaldehyde, 2,3-...	146	C10H10O	051932-70-8	38
5	Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110824\
Data File : VY020225.D
Acq On : 08 Nov 2024 12:05
Operator : SY/MD
Sample : P4734-01
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-03-11062024

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103024S.M
Quant Title : SW846 8260

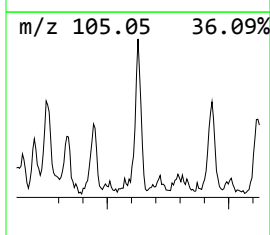
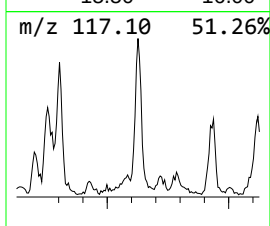
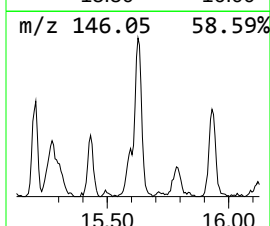
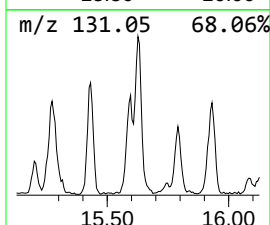
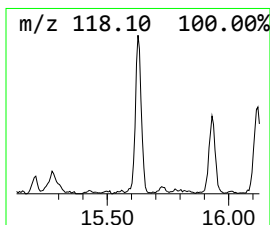
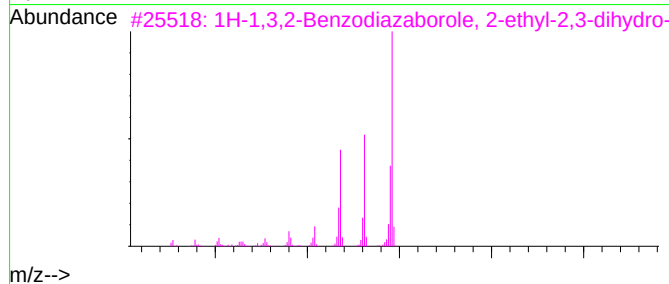
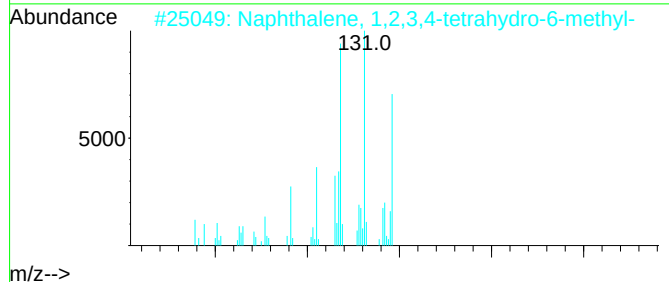
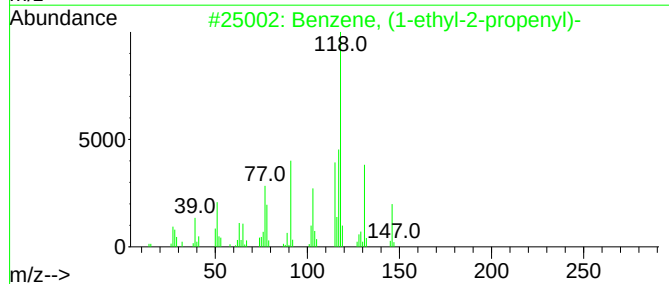
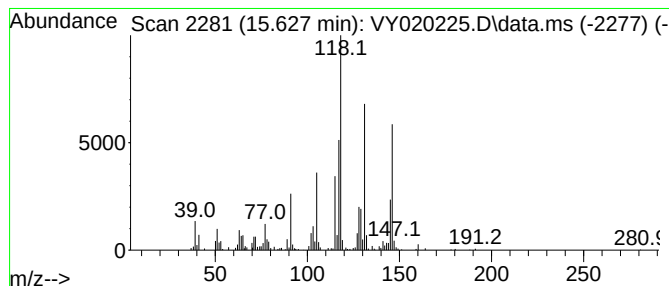
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, (1-ethyl-2-propenyl)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.627	13.71 ug/l	344555	1,4-Dichlorobenzene-d4	13.347

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-ethyl-2-propenyl)-	146	C11H14	019947-22-9	74
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	74
3	1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	64
4	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	50
5	2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1010189-31-0	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110824\
Data File : VY020225.D
Acq On : 08 Nov 2024 12:05
Operator : SY/MD
Sample : P4734-01
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-03-11062024

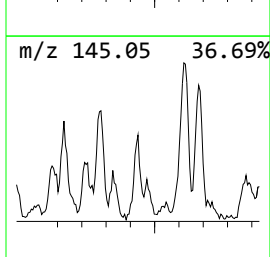
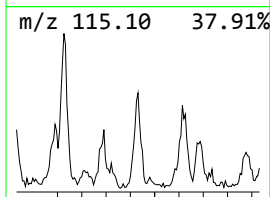
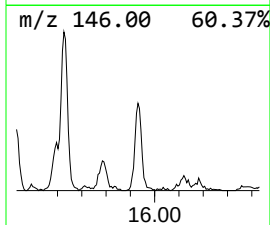
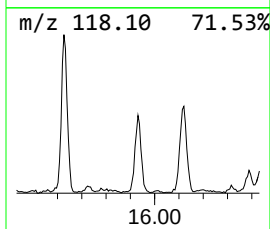
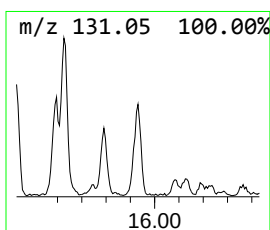
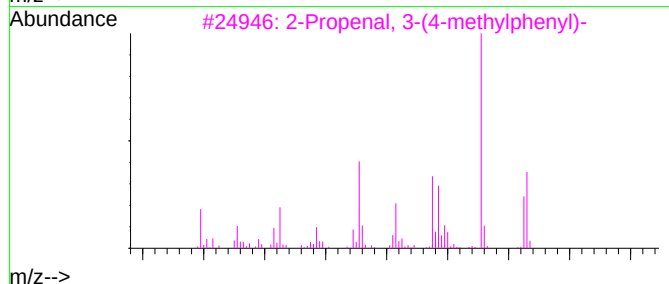
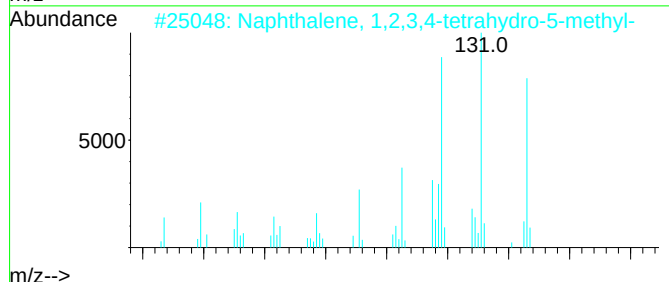
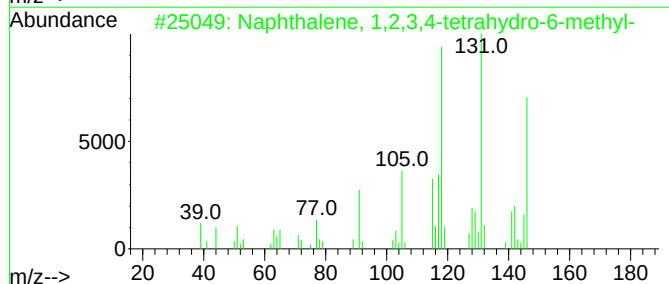
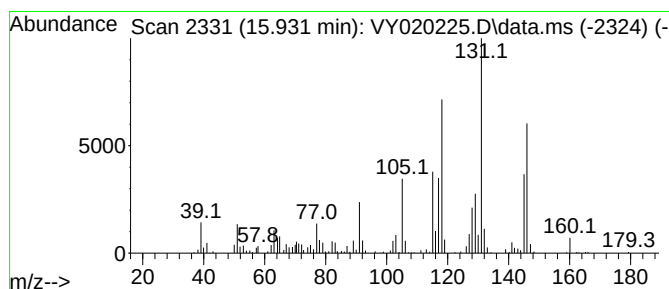
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103024S.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.931	9.46 ug/l	237851	1,4-Dichlorobenzene-d4	13.347

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	93
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	90
3	2-Propenal, 3-(4-methylphenyl)-	146	C10H10O	001504-75-2	62
4	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	58
5	1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110824\
Data File : VY020225.D
Acq On : 08 Nov 2024 12:05
Operator : SY/MD
Sample : P4734-01
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-03-11062024

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103024S.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown13.542	13.542	9.4	ug/l	237434	4	13.347	1256590	50.0
Undecane	13.676	18.9	ug/l	474689	4	13.347	1256590	50.0
Benzene, 2-bute...	13.999	10.4	ug/l	261558	4	13.347	1256590	50.0
unknown14.529	14.529	10.4	ug/l	260881	4	13.347	1256590	50.0
1H-Indene, 2,3-...	14.602	28.1	ug/l	707052	4	13.347	1256590	50.0
Naphthalene, 1,...	14.749	15.7	ug/l	394296	4	13.347	1256590	50.0
Benzene, (1,2-d...	14.968	9.5	ug/l	239777	4	13.347	1256590	50.0
Naphthalene, 1,...	15.255	10.9	ug/l	274540	4	13.347	1256590	50.0
Benzene, (1-eth...	15.627	13.7	ug/l	344555	4	13.347	1256590	50.0
Naphthalene, 1,...	15.931	9.5	ug/l	237851	4	13.347	1256590	50.0