

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY050124\
 Data File : VY018102.D
 Acq On : 01 May 2024 17:03
 Operator : SY/MD
 Sample : P2309-01
 Misc : 5.09g/5.0mL/MSVOA_Y/SOIL/B
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 TR-05-042924

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\82Y050124S.M
 Title : SW846 8260

Signal : TIC: VY018102.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.125	38	50	51	rBV8	16972	48604	3.20%	0.224%
2	4.698	460	472	485	rBV3	26414	85994	5.66%	0.396%
3	5.734	633	642	651	rVB2	10067	31336	2.06%	0.144%
4	6.972	838	845	857	rBV	12055	35835	2.36%	0.165%
5	7.710	953	966	972	rBV	218324	541638	35.65%	2.496%
6	7.783	972	978	990	rVB	367739	834873	54.96%	3.848%
7	8.136	1025	1036	1048	rBV	179623	401472	26.43%	1.850%
8	8.685	1116	1126	1135	rBV	585542	1155403	76.06%	5.325%
9	10.173	1362	1370	1379	rBV	857112	1519151	100.00%	7.001%
10	10.941	1490	1496	1505	rBV2	21116	41131	2.71%	0.190%
11	11.490	1578	1586	1597	rBV	800534	1344081	88.48%	6.194%
12	11.593	1597	1603	1611	rVV7	7128	22316	1.47%	0.103%
13	11.703	1612	1621	1630	rVV4	36666	98888	6.51%	0.456%
14	12.026	1663	1674	1682	rBV4	34808	98667	6.49%	0.455%
15	12.124	1686	1690	1694	rVB	16017	22274	1.47%	0.103%
16	12.203	1698	1703	1705	rBV6	14079	23830	1.57%	0.110%
17	12.239	1705	1709	1714	rVB4	17791	36539	2.41%	0.168%
18	12.325	1719	1723	1726	rBV5	11513	19240	1.27%	0.089%
19	12.410	1734	1737	1739	rBV2	14575	15354	1.01%	0.071%
20	12.477	1743	1748	1754	rVV	597904	941129	61.95%	4.337%
21	12.520	1754	1755	1760	rVB2	23395	20757	1.37%	0.096%
22	12.648	1771	1776	1777	rBV4	14588	21298	1.40%	0.098%
23	12.733	1783	1790	1793	rBV	118359	196612	12.94%	0.906%
24	12.770	1793	1796	1797	rVV	56624	75110	4.94%	0.346%
25	12.806	1797	1802	1807	rVV2	273219	468864	30.86%	2.161%
26	12.861	1807	1811	1815	rVB2	29940	46383	3.05%	0.214%
27	12.971	1822	1829	1833	rBV	82869	162821	10.72%	0.750%
28	13.038	1836	1840	1847	rVB	85185	129446	8.52%	0.597%
29	13.117	1847	1853	1858	rBV	262787	401013	26.40%	1.848%
30	13.166	1858	1861	1867	rVV6	17741	27000	1.78%	0.124%
31	13.245	1867	1874	1878	rVB4	40302	80467	5.30%	0.371%
32	13.312	1878	1885	1889	rBV3	112885	209535	13.79%	0.966%
33	13.361	1889	1893	1896	rVV3	66818	95509	6.29%	0.440%
34	13.422	1896	1903	1913	rVV2	765269	1424413	93.76%	6.564%
35	13.501	1913	1916	1920	rVB2	46437	53930	3.55%	0.249%
36	13.623	1932	1936	1939	rBV2	186039	239251	15.75%	1.103%

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Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y050124S.M
 Title : SW846 8260

37	13.666	1939	1943	1951	rVB3	180636	365051	24.03%	1.682%
38	13.751	1951	1957	1961	rBV	411398	603836	39.75%	2.783%
39	13.818	1961	1968	1974	rBV2	84098	156656	10.31%	0.722%
40	13.885	1975	1979	1981	rBV	99355	143402	9.44%	0.661%
41	13.916	1981	1984	1988	rVB	138876	186187	12.26%	0.858%
42	13.971	1988	1993	1997	rVB	209811	306301	20.16%	1.412%
43	14.020	1997	2001	2005	rVB3	45477	82182	5.41%	0.379%
44	14.081	2005	2011	2015	rVB	244607	386499	25.44%	1.781%
45	14.154	2015	2023	2027	rBV6	50125	138974	9.15%	0.640%
46	14.215	2027	2033	2035	rBV2	72737	147850	9.73%	0.681%
47	14.257	2035	2040	2043	rVV5	124780	248076	16.33%	1.143%
48	14.294	2043	2046	2049	rVV	113605	186956	12.31%	0.862%
49	14.331	2049	2052	2056	rVB	190839	265326	17.47%	1.223%
50	14.379	2056	2060	2063	rBV	137619	175100	11.53%	0.807%
51	14.416	2063	2066	2071	rVB3	114643	153558	10.11%	0.708%
52	14.483	2073	2077	2079	rBV4	38337	59314	3.90%	0.273%
53	14.519	2079	2083	2086	rVB	95921	129567	8.53%	0.597%
54	14.562	2086	2090	2094	rBV	190712	318396	20.96%	1.467%
55	14.611	2094	2098	2103	rVB	325513	458400	30.17%	2.113%
56	14.684	2103	2110	2114	rBV4	446730	960750	63.24%	4.428%
57	14.733	2114	2118	2124	rVB3	135247	250078	16.46%	1.152%
58	14.830	2129	2134	2141	rVB	314149	480292	31.62%	2.213%
59	14.910	2143	2147	2148	rBV4	42322	64305	4.23%	0.296%
60	14.940	2148	2152	2156	rVV3	175081	328938	21.65%	1.516%
61	14.977	2156	2158	2164	rVV	100391	148135	9.75%	0.683%
62	15.056	2164	2171	2178	rVB	179282	385917	25.40%	1.779%
63	15.208	2191	2196	2199	rBV3	78793	141951	9.34%	0.654%
64	15.288	2205	2209	2214	rBV	153238	257687	16.96%	1.188%
65	15.343	2214	2218	2229	rVV4	123359	426031	28.04%	1.963%
66	15.434	2229	2233	2241	rVB4	149497	259956	17.11%	1.198%
67	15.525	2241	2248	2255	rBV3	138614	304824	20.07%	1.405%
68	15.605	2255	2261	2265	rVB3	34388	62956	4.14%	0.290%
69	15.690	2266	2275	2277	rBV3	92894	220332	14.50%	1.015%
70	15.727	2277	2281	2289	rVB	324365	567051	37.33%	2.613%
71	15.812	2291	2295	2300	rBV4	37782	72563	4.78%	0.334%
72	15.885	2301	2307	2314	rBV3	72619	158414	10.43%	0.730%
73	16.031	2324	2331	2338	rBV2	175501	368654	24.27%	1.699%
74	16.227	2356	2363	2369	rVV	162550	320195	21.08%	1.476%
75	16.294	2369	2374	2381	rBV3	77261	158238	10.42%	0.729%
76	16.489	2399	2406	2412	rBV6	59412	167946	11.06%	0.774%

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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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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Title : SW846 8260

77	16.623	2423	2428	2437	rVB3	46386	111757	7.36%	0.515%
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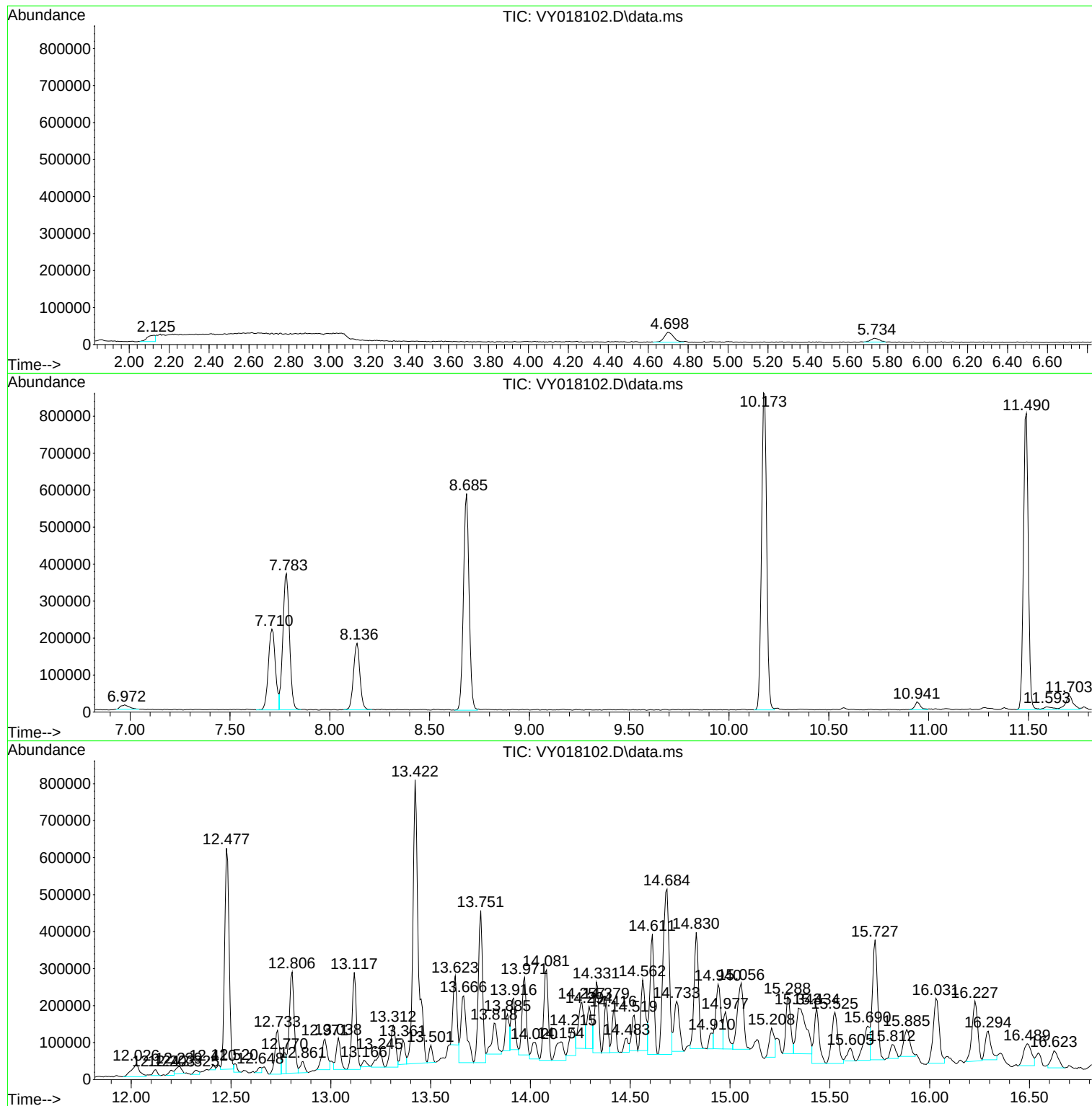
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Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY050124\
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Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y050124S.M
Quant Title : SW846 8260

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



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ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TR-05-042924

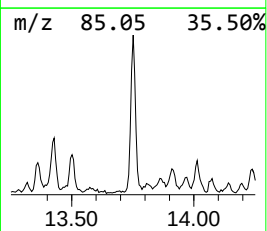
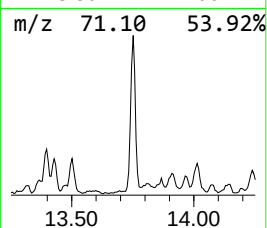
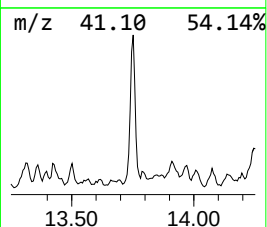
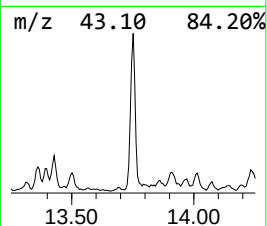
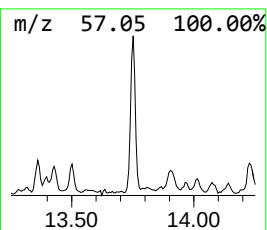
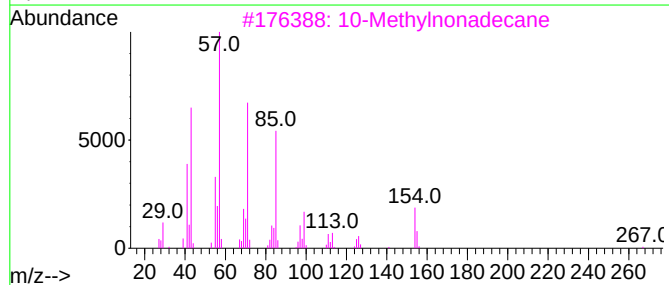
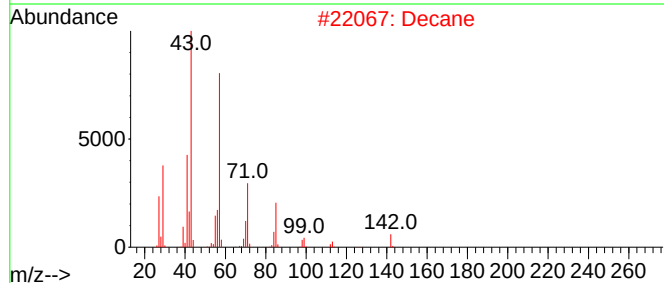
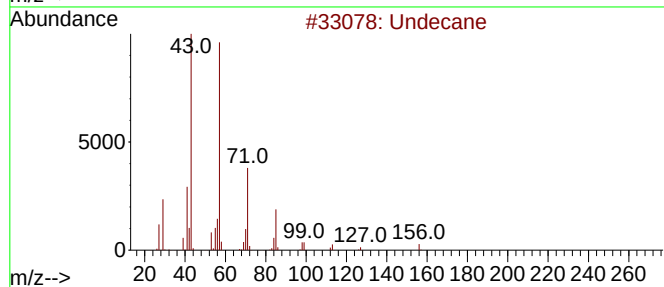
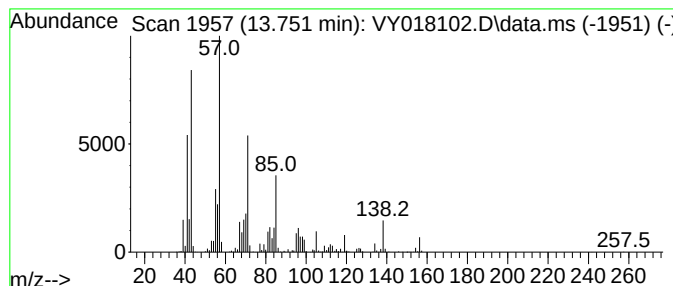
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y050124S.M
Quant Title : SW846 8260

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

Peak Number 1 Undecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.751	21.20 ug/l	603836	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	64
2	Decane			142	C10H22	000124-18-5	64
3	10-Methylnonadecane			282	C20H42	056862-62-5	53
4	Carbonic acid, nonyl vinyl ester			214	C12H22O3	1000383-25-6	50
5	Nonane, 5-(1-methylpropyl)-			184	C13H28	062185-54-0	50



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ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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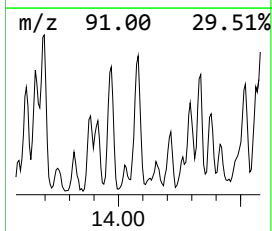
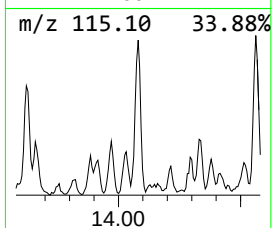
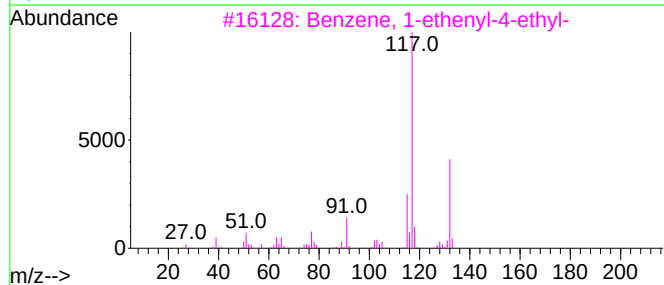
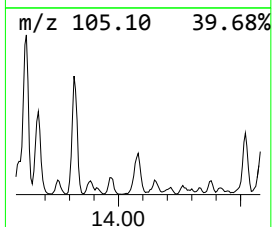
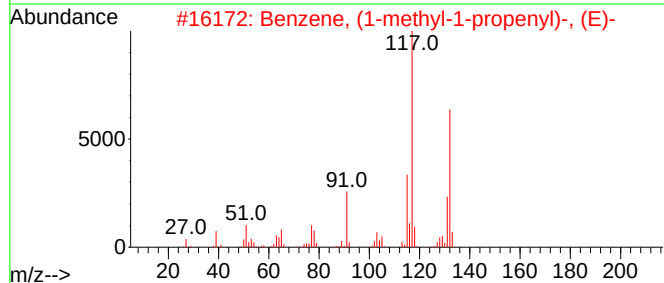
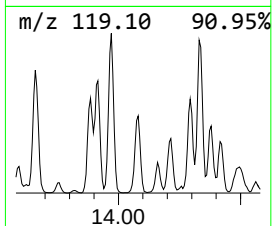
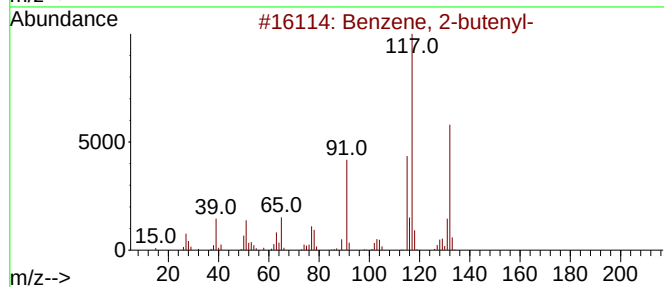
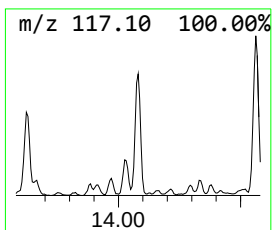
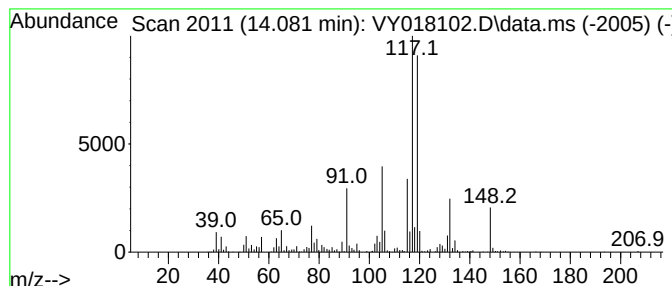
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y050124S.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 2-butenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.081	13.57 ug/l	386499	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
2			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	50
3			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	50
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	50
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	50



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ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TR-05-042924

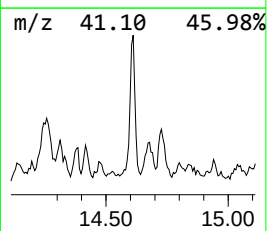
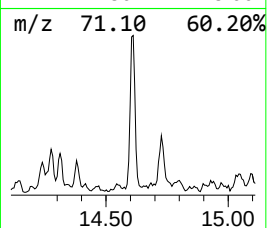
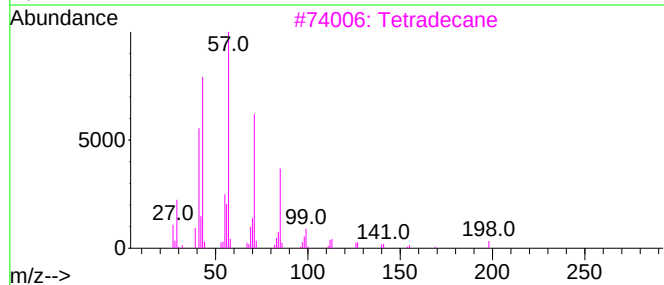
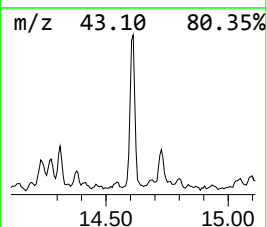
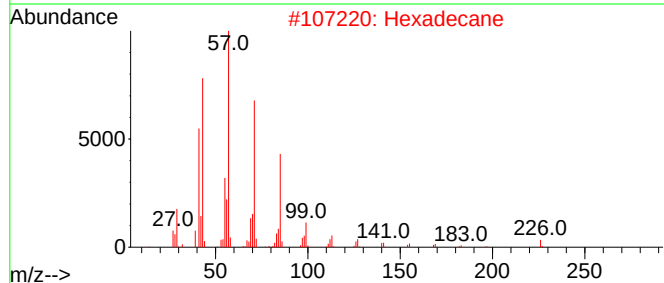
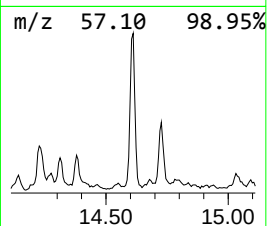
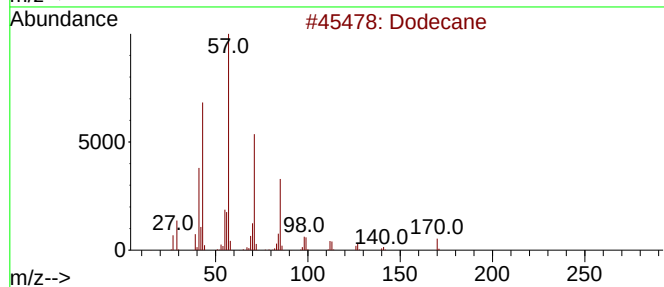
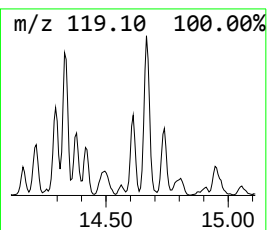
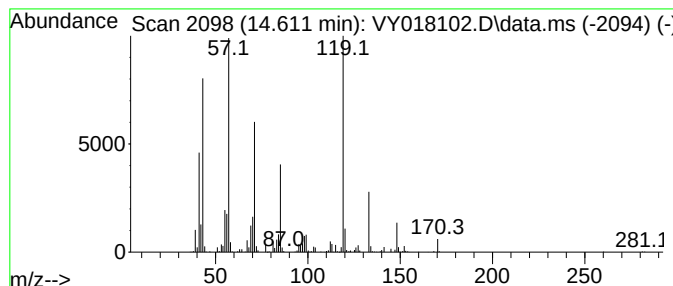
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y050124S.M
Quant Title : SW846 8260

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

Peak Number 3 Dodecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.611	16.09 ug/l	458400	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	94
2			Hexadecane	226	C16H34	000544-76-3	41
3			Tetradecane	198	C14H30	000629-59-4	41
4			Tridecane	184	C13H28	000629-50-5	41
5			Nonadecane	268	C19H40	000629-92-5	38



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ALS Vial : 19 Sample Multiplier: 1

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ClientSampleId :
TR-05-042924

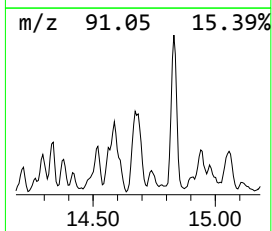
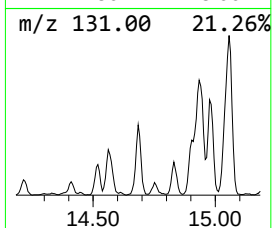
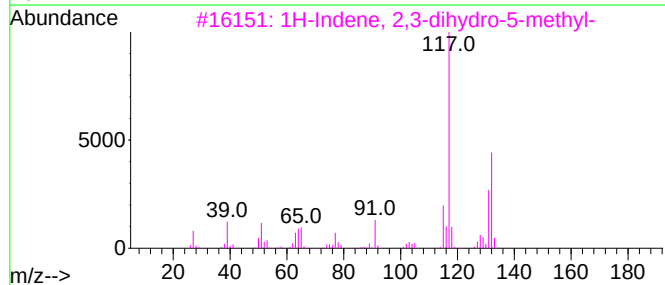
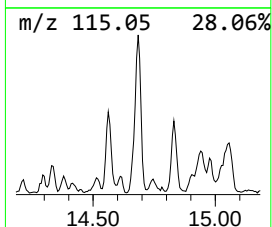
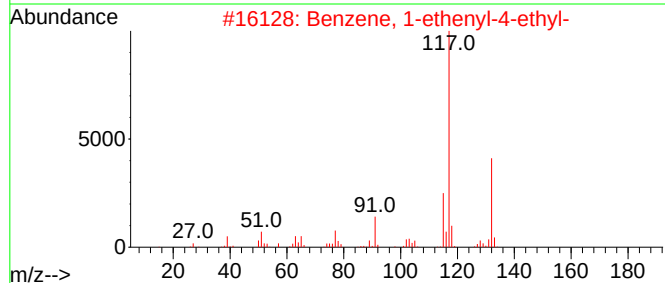
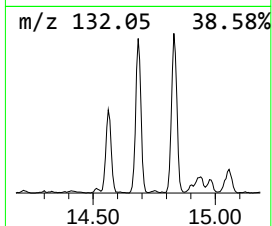
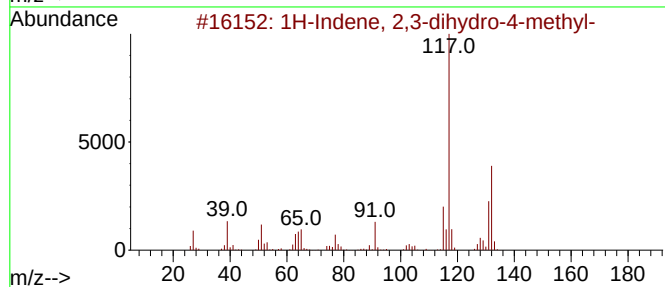
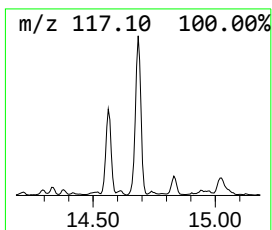
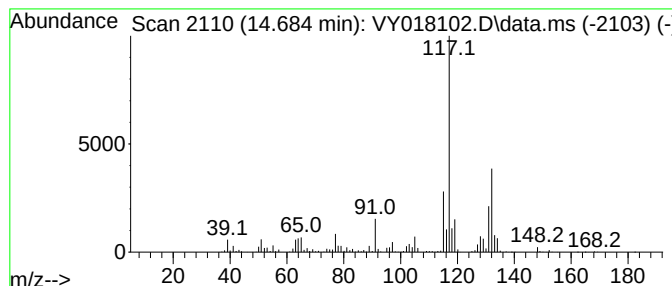
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	33.72 ug/l	960750	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
3			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY050124\
Data File : VY018102.D
Acq On : 01 May 2024 17:03
Operator : SY/MD
Sample : P2309-01
Misc : 5.09g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TR-05-042924

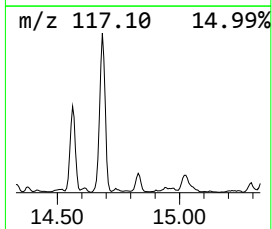
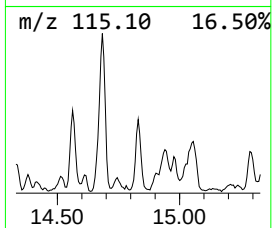
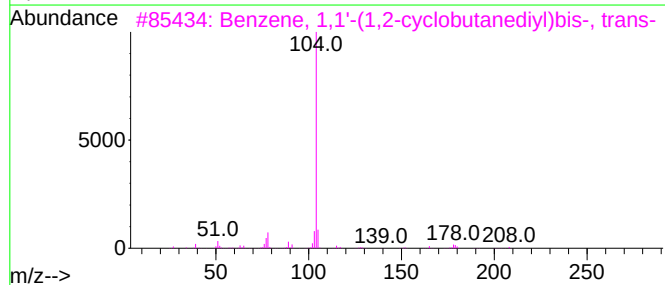
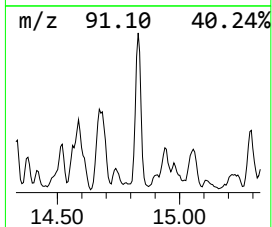
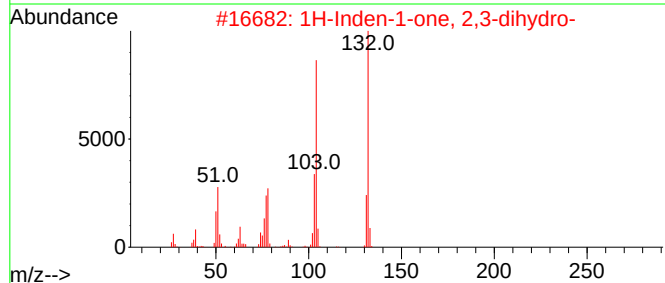
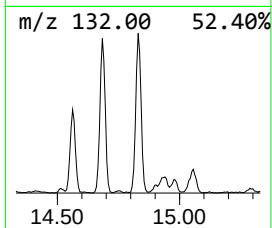
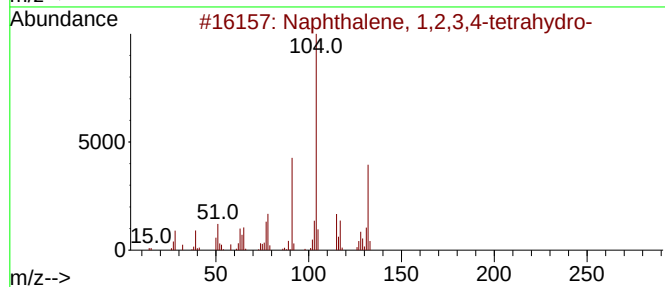
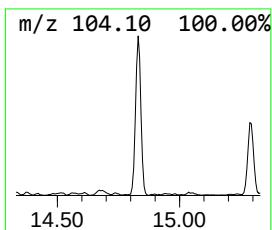
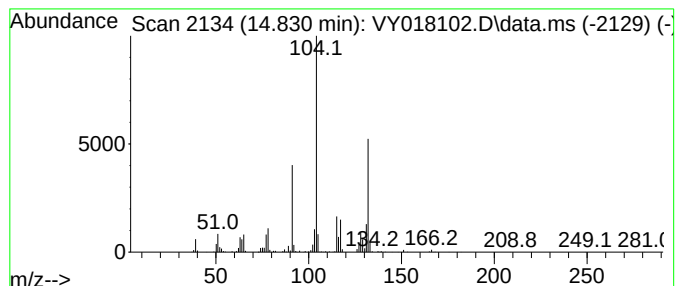
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.830	16.86 ug/l	480292	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2			1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	53
3			Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	38
4			Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	38
5			2-Propenoic acid, 2-phenylethyl ...	176	C11H12O2	003530-36-7	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY050124\
Data File : VY018102.D
Acq On : 01 May 2024 17:03
Operator : SY/MD
Sample : P2309-01
Misc : 5.09g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TR-05-042924

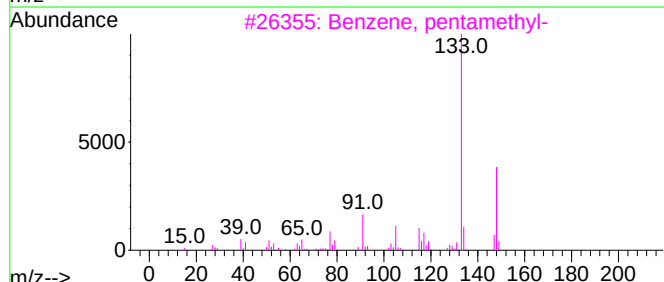
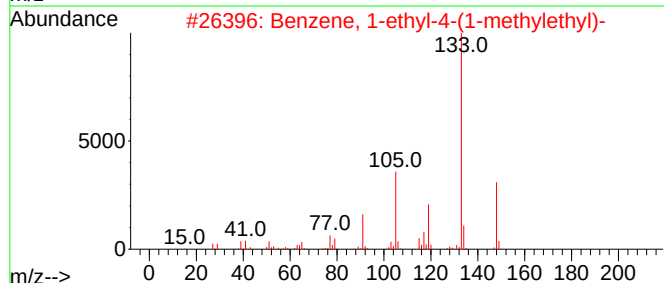
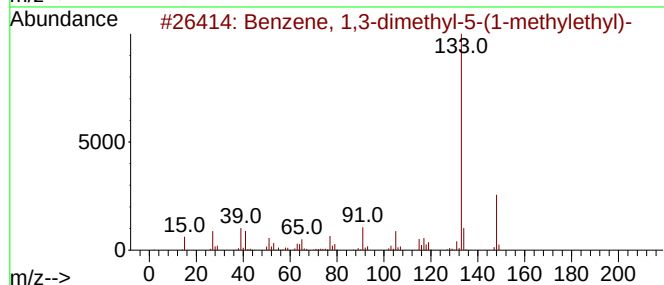
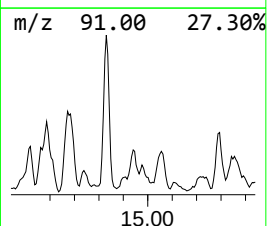
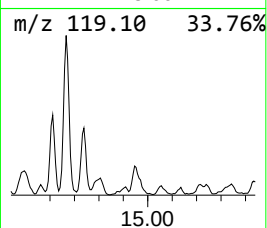
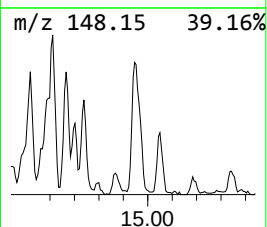
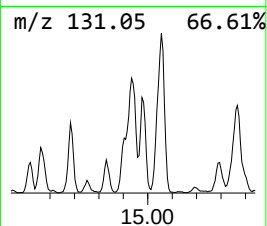
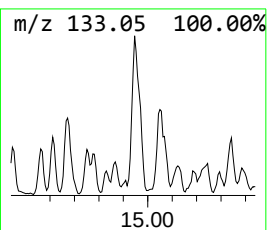
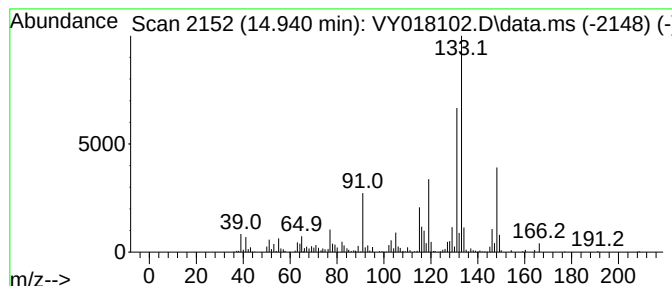
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,3-dimethyl-5-(1-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.940	11.55 ug/l	328938	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	53
2			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	52
3			Benzene, pentamethyl-	148	C11H16	000700-12-9	52
4			Benzene, 1-ethyl-2,4,5-trimethyl-	148	C11H16	017851-27-3	52
5			Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY050124\
Data File : VY018102.D
Acq On : 01 May 2024 17:03
Operator : SY/MD
Sample : P2309-01
Misc : 5.09g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TR-05-042924

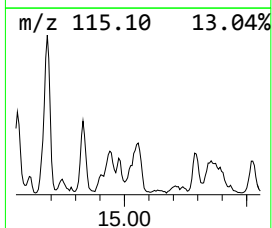
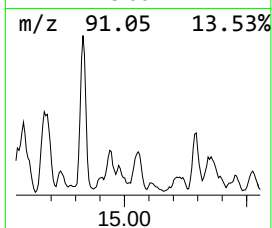
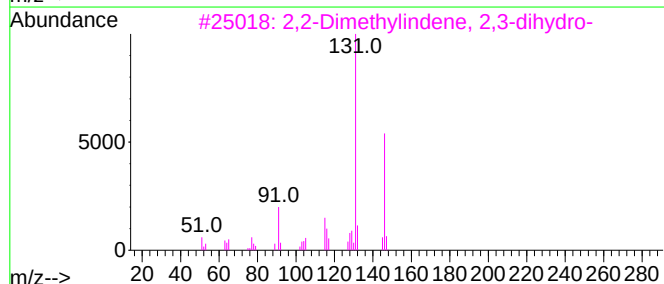
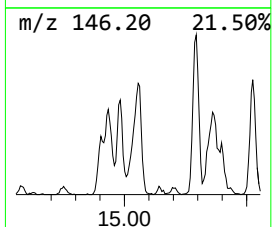
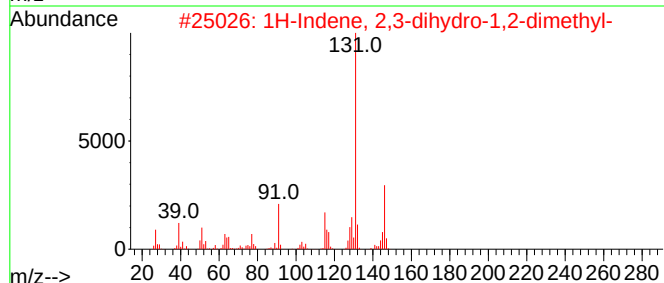
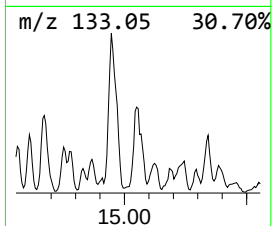
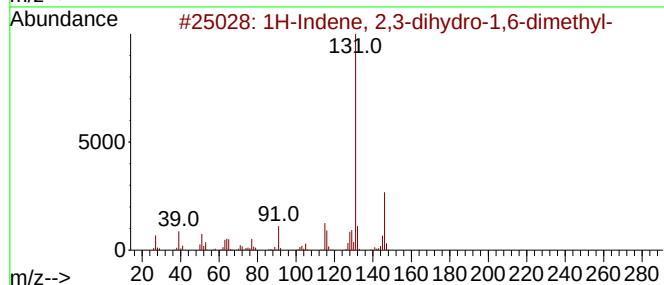
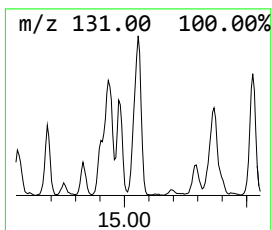
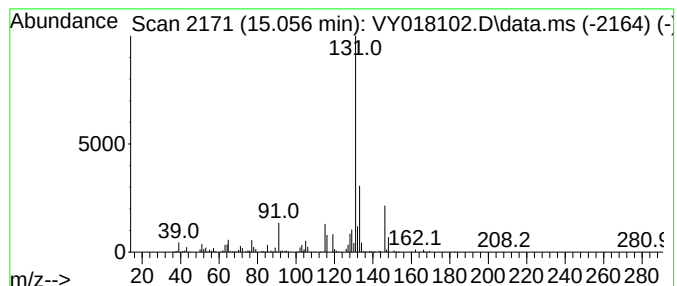
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y050124S.M
Quant Title : SW846 8260

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

Peak Number 7 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.056	13.55 ug/l	385917	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	81
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	81
3			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	81
4			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	81
5			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY050124\
Data File : VY018102.D
Acq On : 01 May 2024 17:03
Operator : SY/MD
Sample : P2309-01
Misc : 5.09g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TR-05-042924

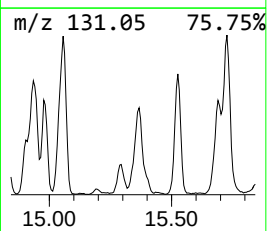
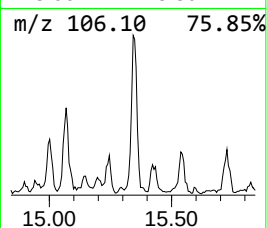
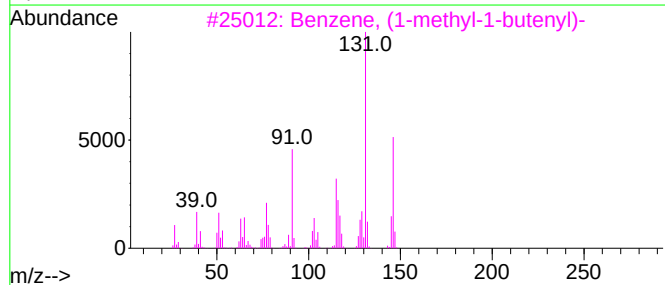
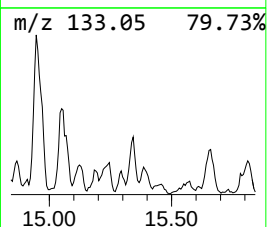
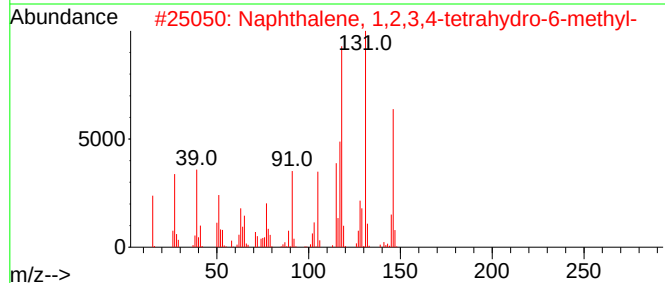
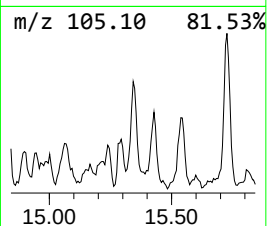
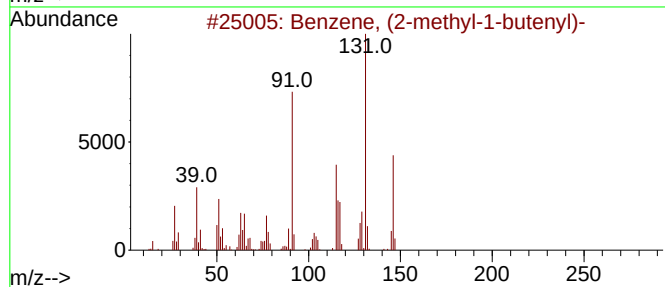
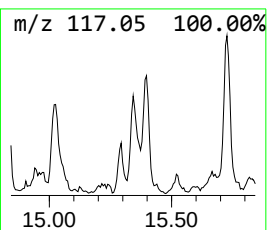
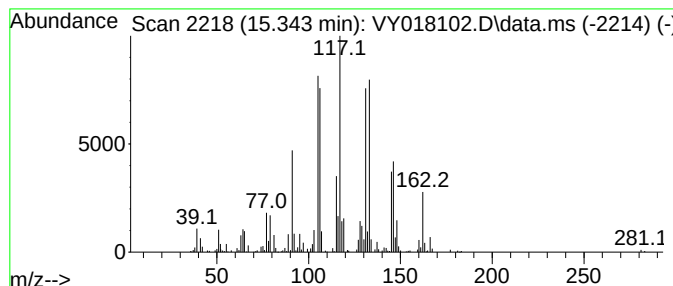
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, (2-methyl-1-butenyl)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.342	14.95 ug/l	426031	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	51
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	42
3			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	38
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	38
5			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY050124\
Data File : VY018102.D
Acq On : 01 May 2024 17:03
Operator : SY/MD
Sample : P2309-01
Misc : 5.09g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TR-05-042924

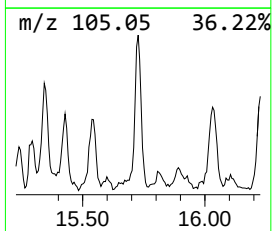
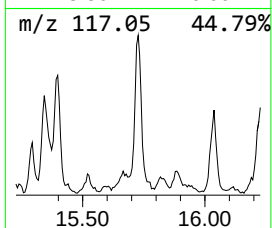
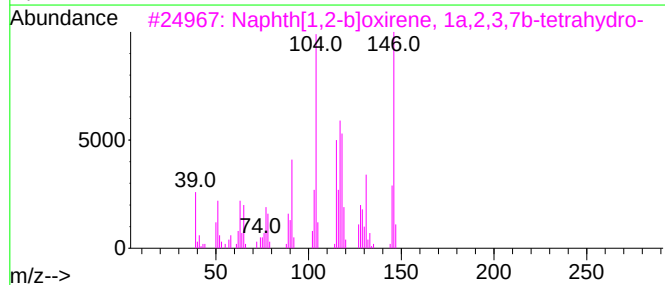
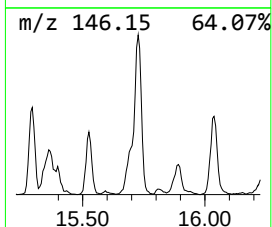
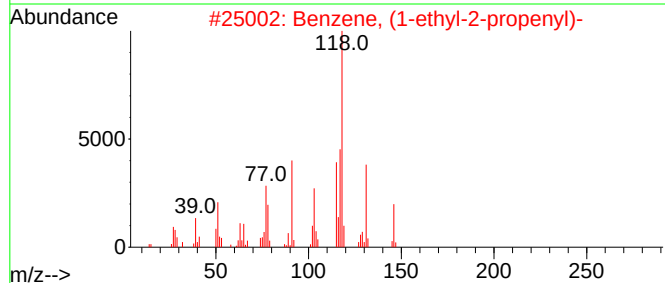
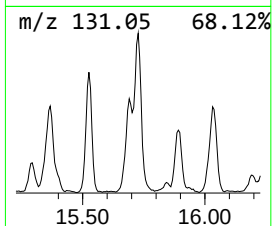
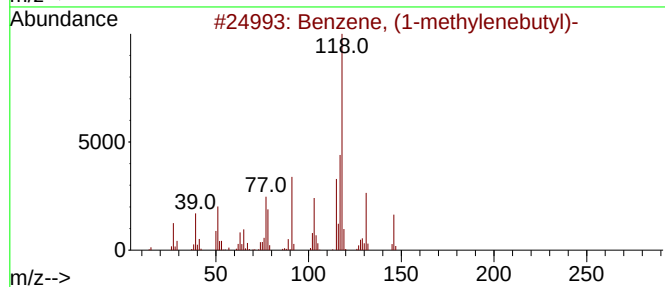
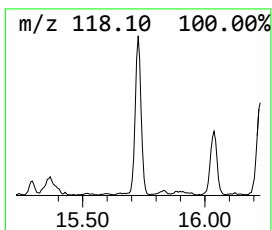
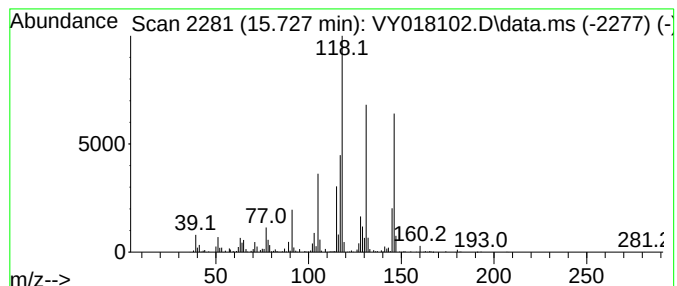
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, (1-methylenebutyl)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.727	19.90 ug/l	567051	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylenebutyl)-	146	C11H14	005676-32-4	87
2			Benzene, (1-ethyl-2-propenyl)-	146	C11H14	019947-22-9	83
3			Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	64
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	64
5			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY050124\
Data File : VY018102.D
Acq On : 01 May 2024 17:03
Operator : SY/MD
Sample : P2309-01
Misc : 5.09g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TR-05-042924

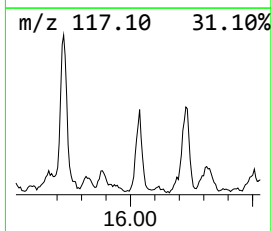
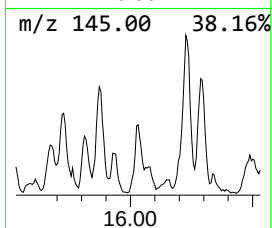
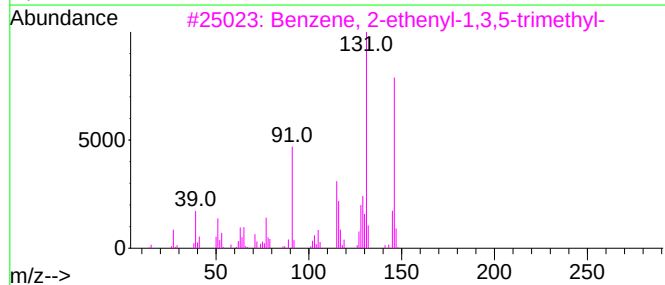
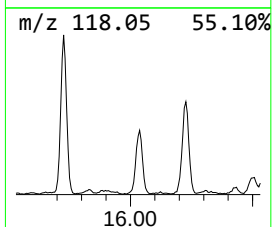
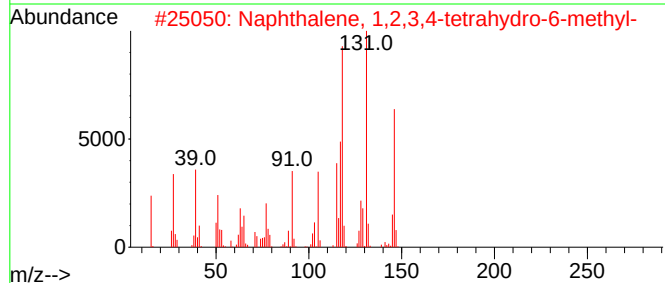
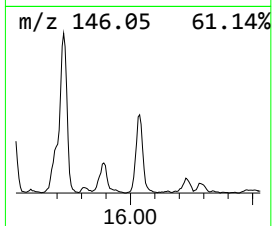
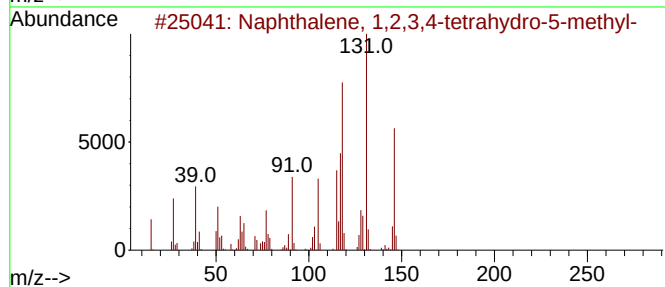
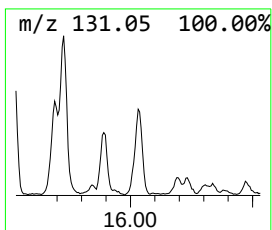
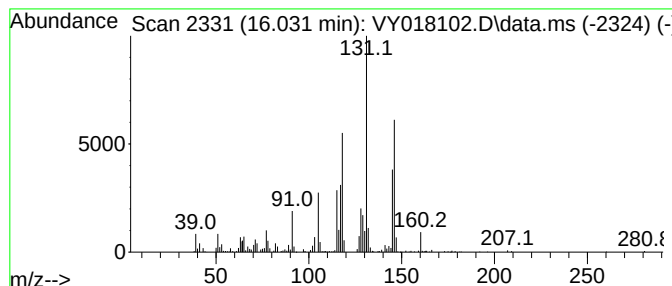
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Quant Title : SW846 8260

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.
16.031	12.94 ug/l	368654	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	93
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	93
3			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	78
4			1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	76
5			1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY050124\
Data File : VY018102.D
Acq On : 01 May 2024 17:03
Operator : SY/MD
Sample : P2309-01
Misc : 5.09g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TR-05-042924

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y050124S.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
Undecane	13.751	21.2	ug/l	603836	4	13.422	1424410	50.0
Benzene, 2-bute...	14.081	13.6	ug/l	386499	4	13.422	1424410	50.0
Dodecane	14.611	16.1	ug/l	458400	4	13.422	1424410	50.0
1H-Indene, 2,3-...	14.684	33.7	ug/l	960750	4	13.422	1424410	50.0
Naphthalene, 1,...	14.830	16.9	ug/l	480292	4	13.422	1424410	50.0
Benzene, 1,3-di...	14.940	11.6	ug/l	328938	4	13.422	1424410	50.0
1H-Indene, 2,3-...	15.056	13.6	ug/l	385917	4	13.422	1424410	50.0
Benzene, (2-met...	15.342	14.9	ug/l	426031	4	13.422	1424410	50.0
Benzene, (1-met...	15.727	19.9	ug/l	567051	4	13.422	1424410	50.0
Naphthalene, 1,...	16.031	12.9	ug/l	368654	4	13.422	1424410	50.0